



Full wwPDB EM Validation Report ⓘ

Jun 17, 2026 – 03:56 pm BST

PDB ID : 8B0X / pdb_00008b0x
EMDB ID : EMD-15793
Title : Translating 70S ribosome in the unrotated state (P and E, tRNAs)
Authors : Fromm, S.A.; O'Connor, K.M.; Purdy, M.; Bhatt, P.R.; Loughran, G.; Atkins, J.F.; Jomaa, A.; Mattei, S.
Deposited on : 2022-09-08
Resolution : 1.55 Å(reported)
Based on initial model : 7K00

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

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

A user guide is available at

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

The types of validation reports are described at

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

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

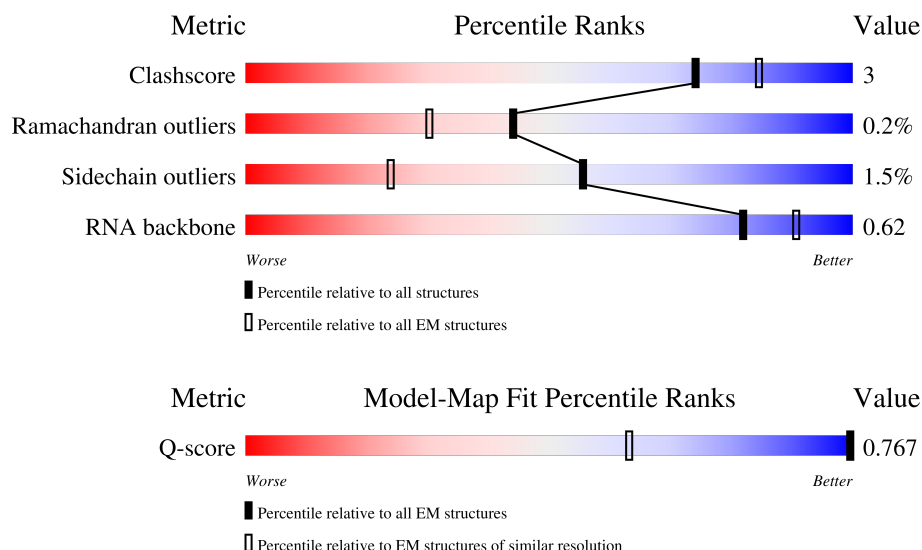
EMDB validation analysis : 0.0.1.dev132
Mogul : 1.8.4, CSD as541be (2020)
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

1 Overall quality at a glance

The following experimental techniques were used to determine the structure:
ELECTRON MICROSCOPY

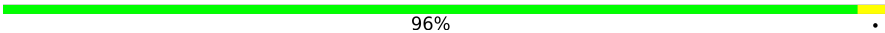
The reported resolution of this entry is 1.55 Å.

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.



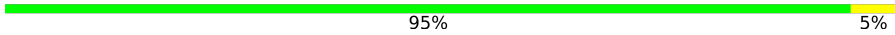
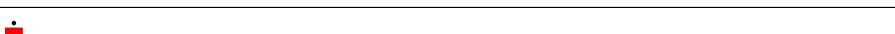
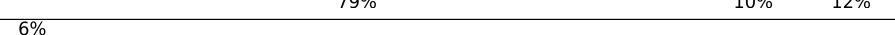
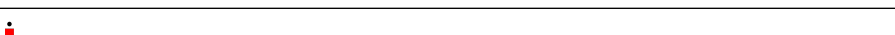
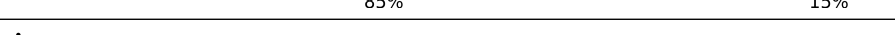
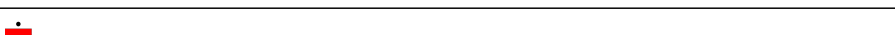
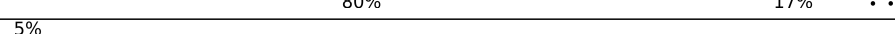
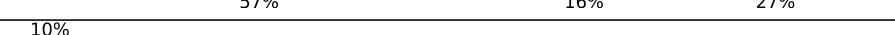
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
RNA backbone	8273	3508	-
Q-score	-	25397	289 (1.05 - 2.05)

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	0	55	 85% 7% 7%
2	1	46	 96% .
3	2	65	 88% 11% .

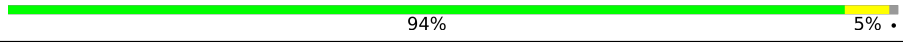
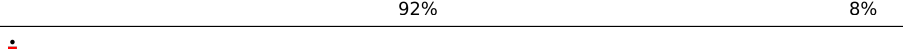
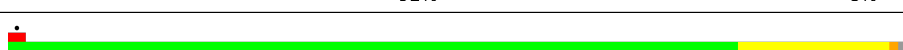
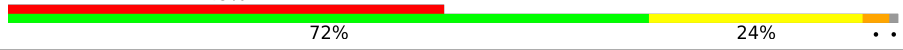
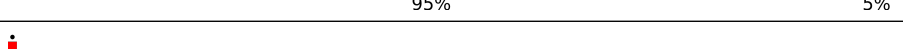
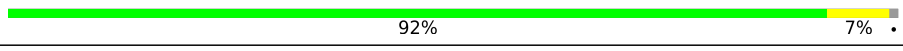
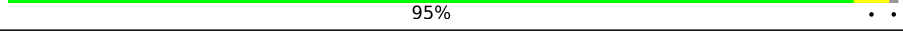
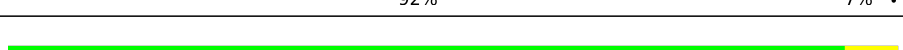
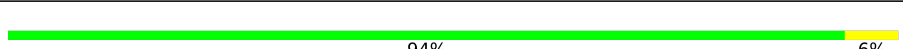
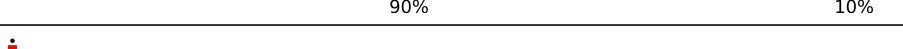
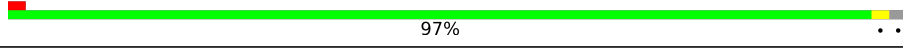
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Mol	Chain	Length	Quality of chain
4	3	38	
5	A	1554	
6	B	241	
7	C	233	
8	D	206	
9	E	167	
10	F	135	
11	G	179	
12	H	130	
13	I	130	
14	J	103	
15	K	129	
16	L	124	
17	M	118	
18	N	101	
19	O	89	
20	P	82	
21	Q	84	
22	R	75	
23	S	92	
24	T	87	
25	U	71	
26	X	6	
27	Z	77	
28	a	2925	

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Mol	Chain	Length	Quality of chain
29	b	119	
30	c	273	
31	d	209	
32	e	201	
33	f	179	
34	g	177	
35	h	149	
36	i	142	
37	j	123	
38	k	144	
39	l	136	
40	m	127	
41	n	117	
42	o	115	
43	p	118	
44	q	103	
45	r	110	
46	s	100	
47	t	104	
48	u	94	
49	v	85	
50	w	78	
51	x	63	
52	y	59	
53	z	57	

2 Entry composition

There are 57 unique types of molecules in this entry. The entry contains 151507 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 protein called 50S ribosomal protein L33.

Mol	Chain	Residues	Atoms				AltConf	Trace
1	0	51	Total	C	N	O	0	0
			417	269	76	72		

- Molecule 2 is a protein called 50S ribosomal protein L34.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	1	46	Total	C	N	O	S	0	0
			377	228	90	57	2		

- Molecule 3 is a protein called 50S ribosomal protein L35.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	2	64	Total	C	N	O	S	0	0
			504	323	105	74	2		

- Molecule 4 is a protein called 50S ribosomal protein L36.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	3	38	Total	C	N	O	S	0	0
			302	185	65	48	4		

- Molecule 5 is a RNA chain called 16S rRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	A	1500	Total	C	N	O	P	0	0
			32211	14373	5920	10418	1500		

- Molecule 6 is a protein called 30S ribosomal protein S2.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	B	224	Total	C	N	O	S	0	0
			1753	1109	315	321	8		

- Molecule 7 is a protein called 30S ribosomal protein S3.

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

- Molecule 8 is a protein called 30S ribosomal protein S4.

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

- Molecule 9 is a protein called 30S ribosomal protein S5.

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

- Molecule 10 is a protein called 30S ribosomal protein S6.

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

- Molecule 11 is a protein called 30S ribosomal protein S7.

Mol	Chain	Residues	Atoms					AltConf	Trace
11	G	153	Total	C	N	O	S	0	0
			1203	750	231	218	4		

- Molecule 12 is a protein called 30S ribosomal protein S8.

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

- Molecule 13 is a protein called 30S ribosomal protein S9.

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

- Molecule 14 is a protein called 30S ribosomal protein S10.

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

- Molecule 15 is a protein called 30S ribosomal protein S11.

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

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
K	119	IAS	ASN	modified residue	UNP P0A7R9

- Molecule 16 is a protein called 30S ribosomal protein S12.

Mol	Chain	Residues	Atoms					AltConf	Trace
16	L	123	Total	C	N	O	S	0	0
			957	591	196	165	5		

- Molecule 17 is a protein called 30S ribosomal protein S13.

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

- Molecule 18 is a protein called 30S ribosomal protein S14.

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

- Molecule 19 is a protein called 30S ribosomal protein S15.

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

- Molecule 20 is a protein called 30S ribosomal protein S16.

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

- Molecule 21 is a protein called 30S ribosomal protein S17.

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

- Molecule 22 is a protein called 30S ribosomal protein S18.

Mol	Chain	Residues	Atoms				AltConf	Trace
22	R	55	Total	C	N	O	0	0
			455	288	86	81		

- Molecule 23 is a protein called 30S ribosomal protein S19.

Mol	Chain	Residues	Atoms					AltConf	Trace
23	S	84	Total	C	N	O	S	0	0
			668	427	127	112	2		

- Molecule 24 is a protein called 30S ribosomal protein S20.

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

- Molecule 25 is a protein called 30S ribosomal protein S21.

Mol	Chain	Residues	Atoms					AltConf	Trace
25	U	70	Total	C	N	O	S	0	0
			589	366	125	97	1		

- Molecule 26 is a RNA chain called mRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
26	X	6	Total	C	N	O	P	0	0
			131	59	27	39	6		

- Molecule 27 is a RNA chain called P-site tRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
27	Z	75	Total	C	N	O	P	0	0
			1603	714	292	522	75		

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
Z	34	U	C	conflict	GB 1804121330
Z	36	C	U	conflict	GB 1804121330

- Molecule 28 is a RNA chain called 23S rRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
28	a	2748	Total	C	N	O	P	0	0
			59025	26336	10876	19065	2748		

- Molecule 29 is a RNA chain called 5S rRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
29	b	119	Total	C	N	O	P	0	0
			2549	1135	466	829	119		

- Molecule 30 is a protein called 50S ribosomal protein L2.

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

- Molecule 31 is a protein called 50S ribosomal protein L3.

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

- Molecule 32 is a protein called 50S ribosomal protein L4.

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

- Molecule 33 is a protein called 50S ribosomal protein L5.

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

- Molecule 34 is a protein called 50S ribosomal protein L6.

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

- Molecule 35 is a protein called 50S ribosomal protein L9.

Mol	Chain	Residues	Atoms					AltConf	Trace
35	h	148	Total	C	N	O	S	0	0
			1101	694	196	210	1		

- Molecule 36 is a protein called 50S ribosomal protein L13.

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

- Molecule 37 is a protein called 50S ribosomal protein L14.

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

- Molecule 38 is a protein called 50S ribosomal protein L15.

Mol	Chain	Residues	Atoms					AltConf	Trace
38	k	144	Total	C	N	O	S	0	0
			1053	654	207	190	2		

- Molecule 39 is a protein called 50S ribosomal protein L16.

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

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
1	82	MS6	MET	modified residue	UNP P0ADY7

- Molecule 40 is a protein called 50S ribosomal protein L17.

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

- Molecule 41 is a protein called 50S ribosomal protein L18.

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

- Molecule 42 is a protein called 50S ribosomal protein L19.

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

- Molecule 43 is a protein called 50S ribosomal protein L20.

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

- Molecule 44 is a protein called 50S ribosomal protein L21.

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

- Molecule 45 is a protein called 50S ribosomal protein L22.

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

- Molecule 46 is a protein called 50S ribosomal protein L23.

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

- Molecule 47 is a protein called 50S ribosomal protein L24.

Mol	Chain	Residues	Atoms					AltConf	Trace
47	t	102	Total	C	N	O	S	0	0
			779	492	146	141			

- Molecule 48 is a protein called 50S ribosomal protein L25.

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

- Molecule 49 is a protein called 50S ribosomal protein L27.

Mol	Chain	Residues	Atoms					AltConf	Trace
49	v	78	Total	C	N	O	S	0	0
			586	362	116	107	1		

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

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

- Molecule 51 is a protein called 50S ribosomal protein L29.

Mol	Chain	Residues	Atoms					AltConf	Trace
51	x	62	Total	C	N	O	S	0	0
			501	308	98	94	1		

- Molecule 52 is a protein called 50S ribosomal protein L30.

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

- Molecule 53 is a protein called 50S ribosomal protein L32.

Mol	Chain	Residues	Atoms					AltConf	Trace
53	z	56	Total	C	N	O	S	0	0
			444	269	94	80	1		

- Molecule 54 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		AltConf
54	3	1	Total	Zn	0
			1	1	

- Molecule 55 is POTASSIUM ION (CCD ID: K) (formula: K).

Mol	Chain	Residues	Atoms		AltConf
55	A	48	Total	K	0
			48	48	
55	a	112	Total	K	0
			112	112	
55	b	1	Total	K	0
			1	1	
55	c	4	Total	K	0
			4	4	
55	d	1	Total	K	0
			1	1	
55	e	1	Total	K	0
			1	1	
55	f	1	Total	K	0
			1	1	

- Molecule 56 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		AltConf
56	A	97	Total	Mg	0
			97	97	
56	X	1	Total	Mg	0
			1	1	
56	a	254	Total	Mg	0
			254	254	
56	b	5	Total	Mg	0
			5	5	
56	c	2	Total	Mg	0
			2	2	
56	d	1	Total	Mg	0
			1	1	

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Mol	Chain	Residues	Atoms		AltConf
56	z	1	Total 1	Mg 1	0

- Molecule 57 is water.

Mol	Chain	Residues	Atoms		AltConf
57	0	5	Total 5	O 5	0
57	1	51	Total 51	O 51	0
57	2	45	Total 45	O 45	0
57	3	13	Total 13	O 13	0
57	A	2886	Total 2886	O 2886	0
57	B	7	Total 7	O 7	0
57	C	16	Total 16	O 16	0
57	D	14	Total 14	O 14	0
57	E	38	Total 38	O 38	0
57	F	5	Total 5	O 5	0
57	G	4	Total 4	O 4	0
57	H	45	Total 45	O 45	0
57	I	11	Total 11	O 11	0
57	J	15	Total 15	O 15	0
57	K	22	Total 22	O 22	0
57	L	35	Total 35	O 35	0
57	M	2	Total 2	O 2	0
57	N	8	Total 8	O 8	0

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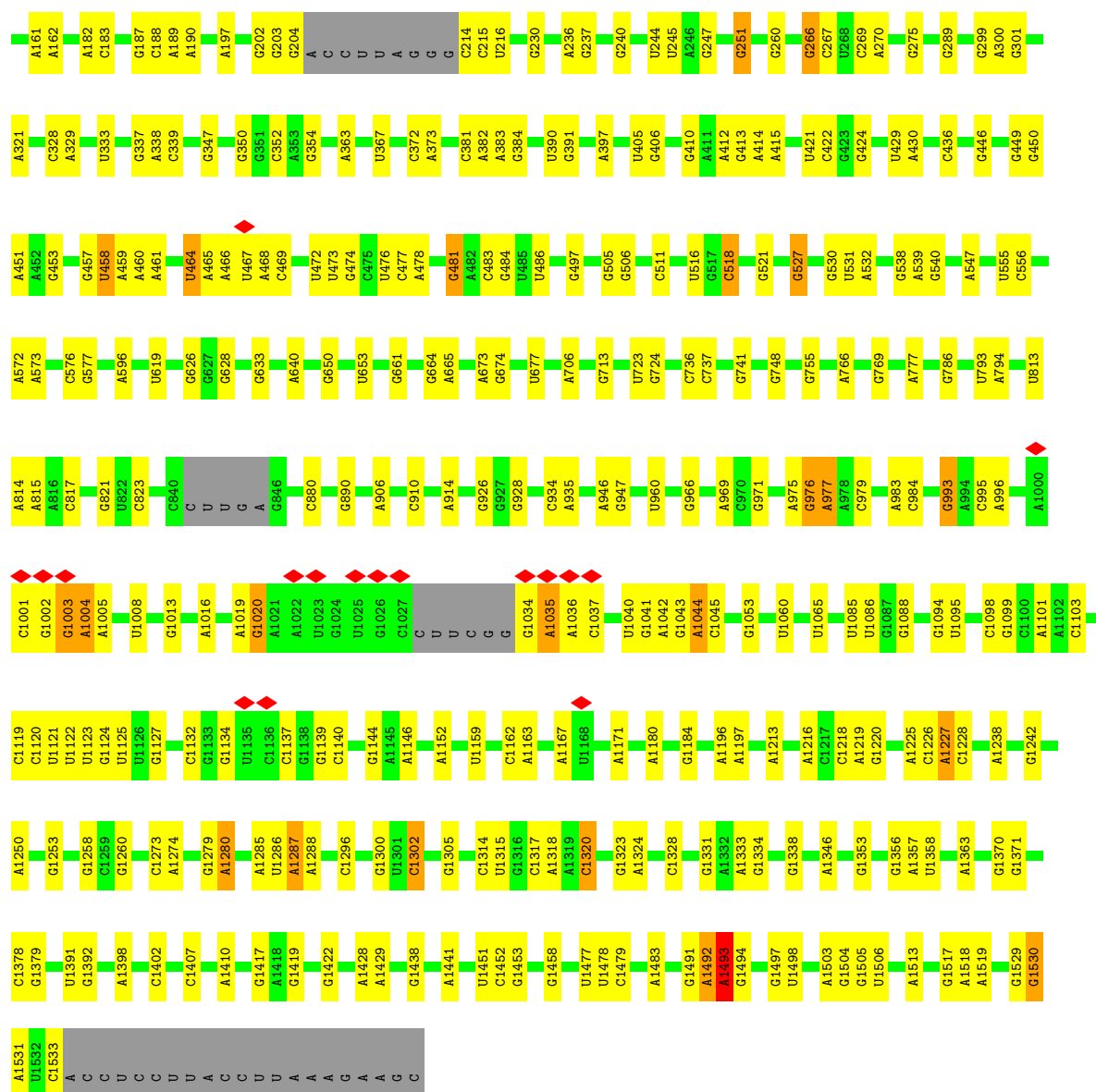
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Mol	Chain	Residues	Atoms		AltConf
57	O	29	Total 29	O 29	0
57	P	22	Total 22	O 22	0
57	Q	9	Total 9	O 9	0
57	R	13	Total 13	O 13	0
57	T	18	Total 18	O 18	0
57	U	17	Total 17	O 17	0
57	X	19	Total 19	O 19	0
57	Z	20	Total 20	O 20	0
57	a	6905	Total 6905	O 6905	0
57	b	81	Total 81	O 81	0
57	c	195	Total 195	O 195	0
57	d	107	Total 107	O 107	0
57	e	80	Total 80	O 80	0
57	h	1	Total 1	O 1	0
57	i	70	Total 70	O 70	0
57	j	58	Total 58	O 58	0
57	k	72	Total 72	O 72	0
57	l	57	Total 57	O 57	0
57	m	61	Total 61	O 61	0
57	n	3	Total 3	O 3	0
57	o	71	Total 71	O 71	0

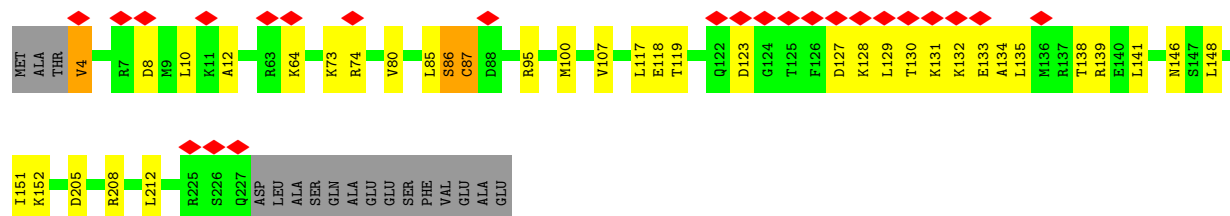
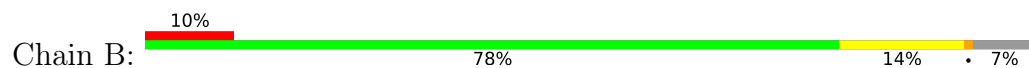
Continued on next page...

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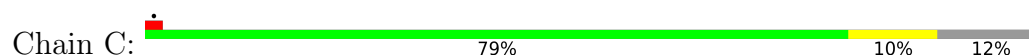
Mol	Chain	Residues	Atoms		AltConf
57	p	66	Total 66	O 66	0
57	q	37	Total 37	O 37	0
57	r	58	Total 58	O 58	0
57	s	24	Total 24	O 24	0
57	t	8	Total 8	O 8	0
57	u	5	Total 5	O 5	0
57	v	33	Total 33	O 33	0
57	w	28	Total 28	O 28	0
57	x	2	Total 2	O 2	0
57	y	16	Total 16	O 16	0
57	z	54	Total 54	O 54	0

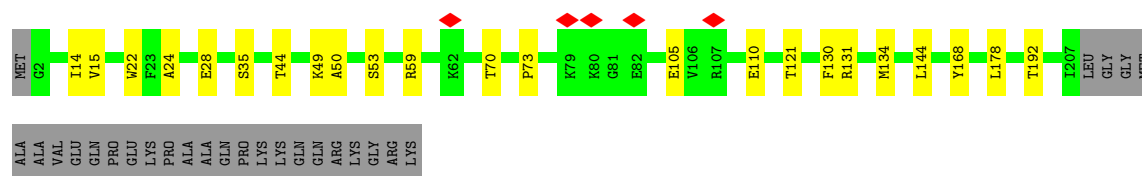


• Molecule 6: 30S ribosomal protein S2

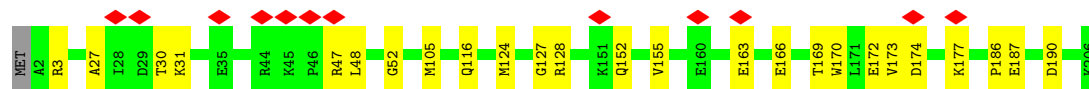
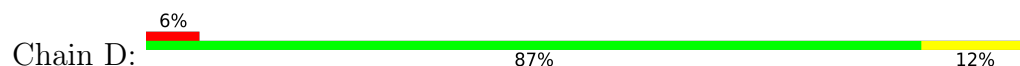


• Molecule 7: 30S ribosomal protein S3

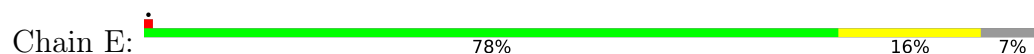




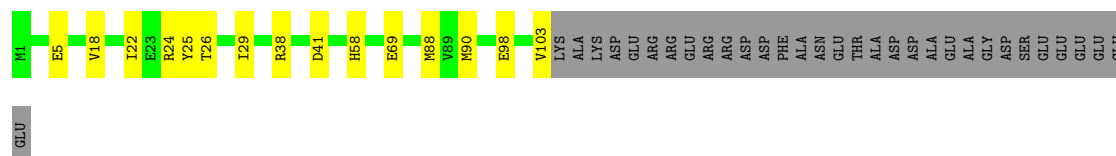
- Molecule 8: 30S ribosomal protein S4



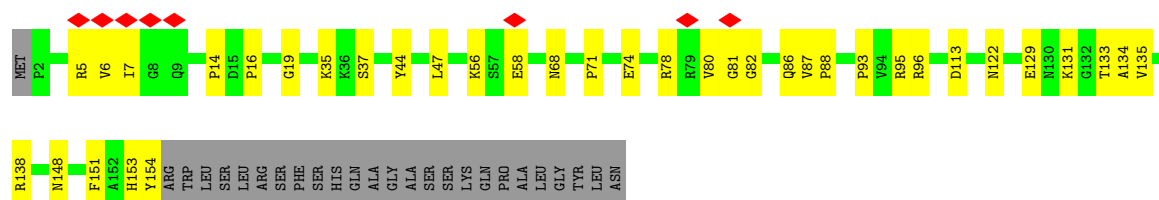
- Molecule 9: 30S ribosomal protein S5



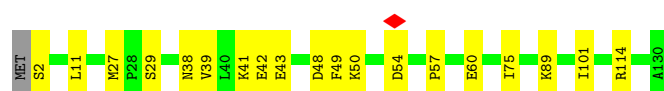
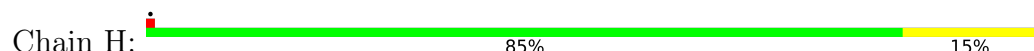
- Molecule 10: 30S ribosomal protein S6



- Molecule 11: 30S ribosomal protein S7

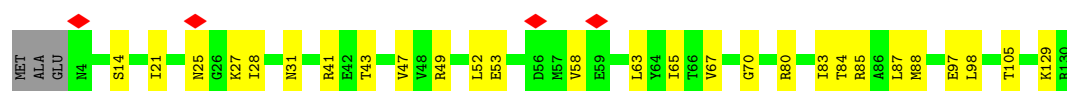


- Molecule 12: 30S ribosomal protein S8



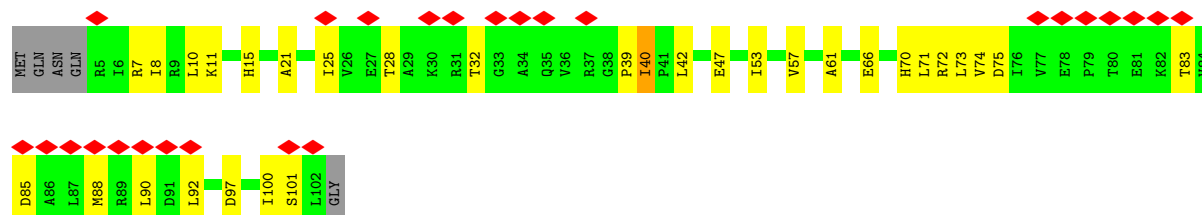
- Molecule 13: 30S ribosomal protein S9

Chain I:  77% 21%



- Molecule 14: 30S ribosomal protein S10

Chain J:  25% 65% 29% 5%

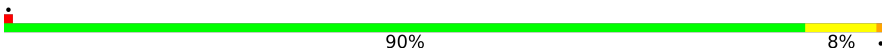


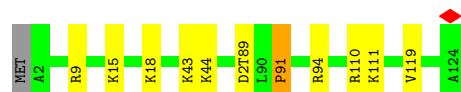
- Molecule 15: 30S ribosomal protein S11

Chain K:  72% 18% 9%



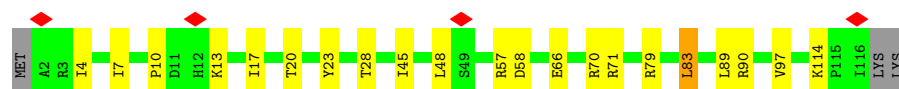
- Molecule 16: 30S ribosomal protein S12

Chain L:  90% 8%



- Molecule 17: 30S ribosomal protein S13

Chain M:  80% 17%



- Molecule 18: 30S ribosomal protein S14

Chain N:  5% 83% 16%



- Molecule 19: 30S ribosomal protein S15

Chain O:  91% 8%



- Molecule 20: 30S ribosomal protein S16

Chain P:  91% 7%



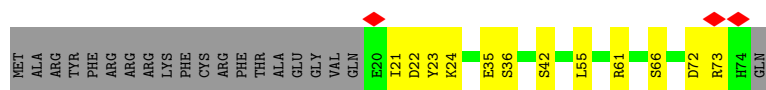
- Molecule 21: 30S ribosomal protein S17

Chain Q:  80% 13% 6%



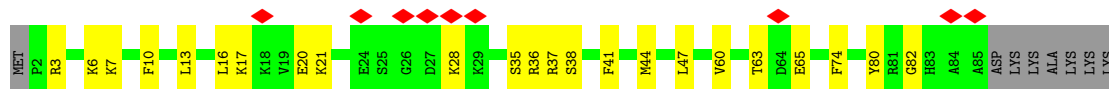
- Molecule 22: 30S ribosomal protein S18

Chain R:  57% 16% 27%



- Molecule 23: 30S ribosomal protein S19

Chain S:  10% 66% 25% 9%



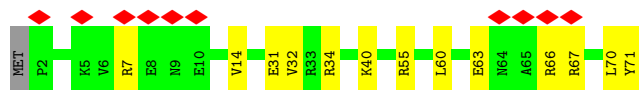
- Molecule 24: 30S ribosomal protein S20

Chain T:  89% 10%



- Molecule 25: 30S ribosomal protein S21

Chain U:  14% 80% 18%



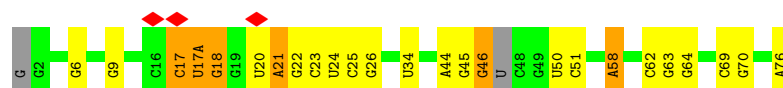
- Molecule 26: mRNA

Chain X:  100%

There are no outlier residues recorded for this chain.

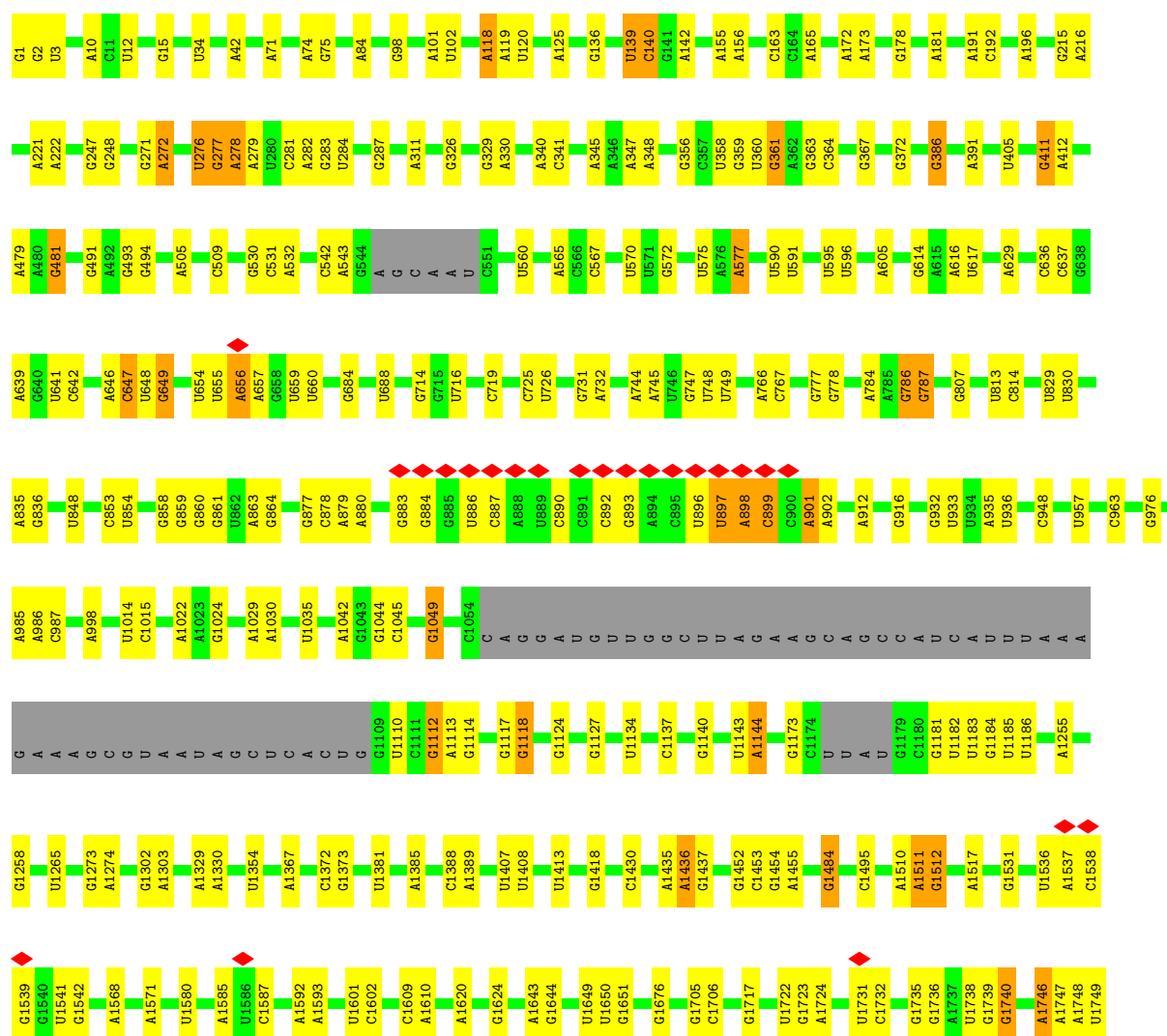
- Molecule 27: P-site tRNA

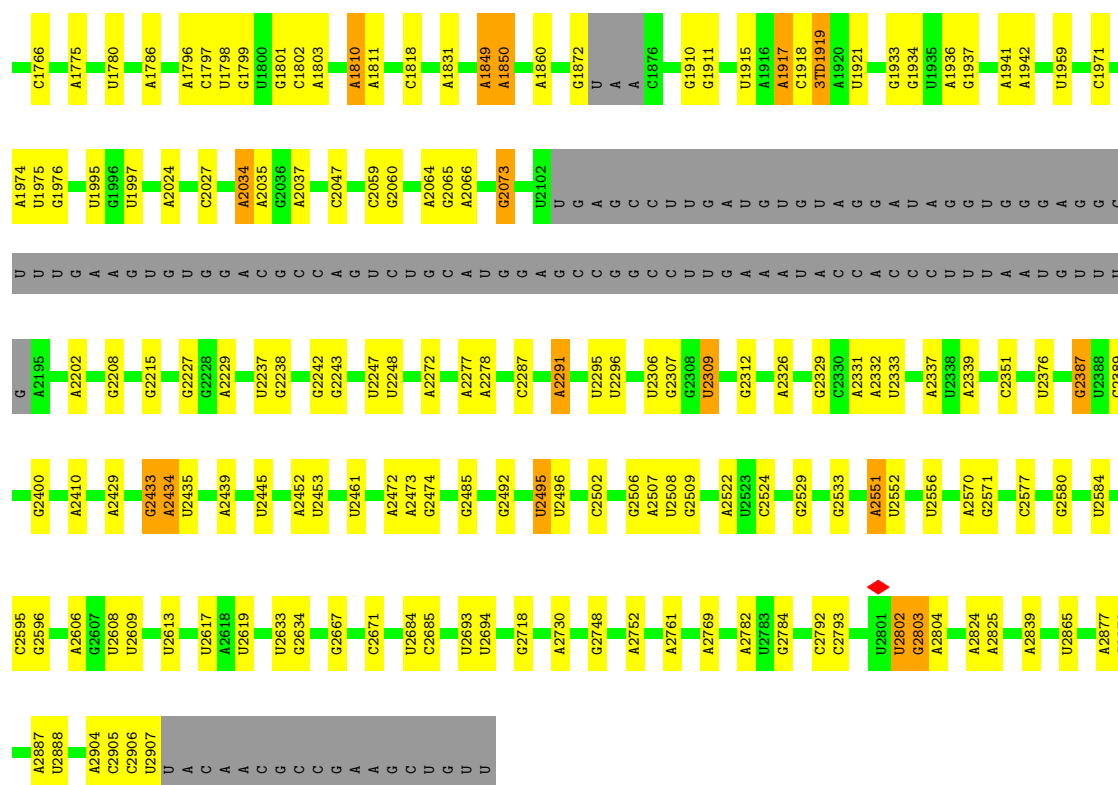
Chain Z:  65% 25% 8%



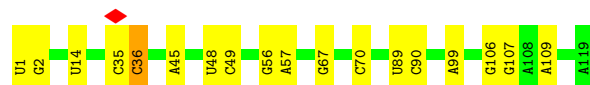
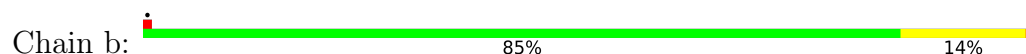
- Molecule 28: 23S rRNA

Chain a:  78% 14% 6%





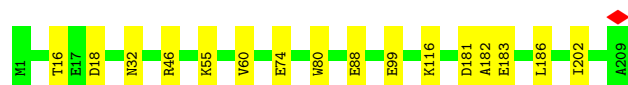
• Molecule 29: 5S rRNA



• Molecule 30: 50S ribosomal protein L2



• Molecule 31: 50S ribosomal protein L3



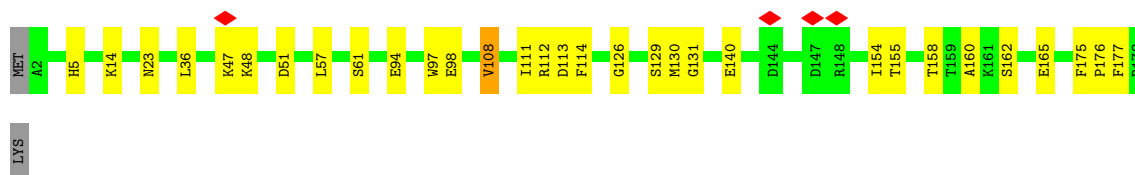
• Molecule 32: 50S ribosomal protein L4





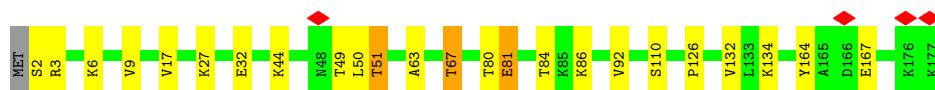
- Molecule 33: 50S ribosomal protein L5

Chain f: 82% 17% ..



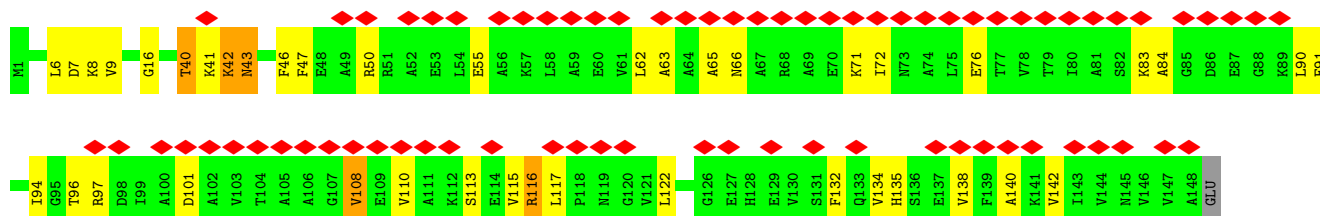
- Molecule 34: 50S ribosomal protein L6

Chain g: 86% 12% ..



- Molecule 35: 50S ribosomal protein L9

Chain h: 49% 72% 24% ..



- Molecule 36: 50S ribosomal protein L13

Chain i: 83% 16% .



- Molecule 37: 50S ribosomal protein L14

Chain j: 95% 5%



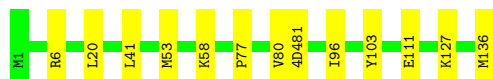
- Molecule 38: 50S ribosomal protein L15

Chain k: 96% .



- Molecule 39: 50S ribosomal protein L16

Chain l: 90% 10%



- Molecule 40: 50S ribosomal protein L17

Chain m: 89% 7%



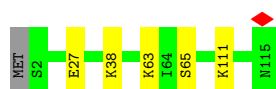
- Molecule 41: 50S ribosomal protein L18

Chain n: 92% 7%



- Molecule 42: 50S ribosomal protein L19

Chain o: 95%



- Molecule 43: 50S ribosomal protein L20

Chain p: 92% 7%



- Molecule 44: 50S ribosomal protein L21

Chain q: 94% 6%

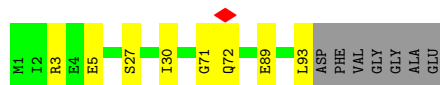
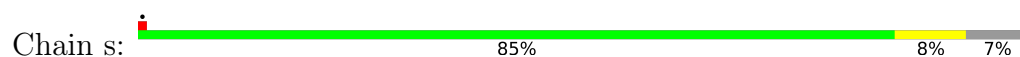


- Molecule 45: 50S ribosomal protein L22

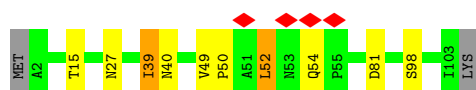
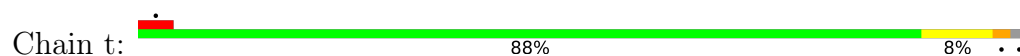
Chain r: 94% 6%



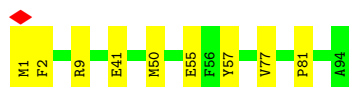
- Molecule 46: 50S ribosomal protein L23



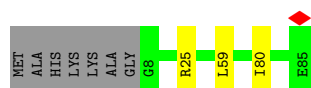
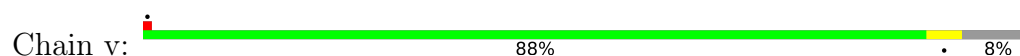
- Molecule 47: 50S ribosomal protein L24



- Molecule 48: 50S ribosomal protein L25



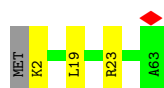
- Molecule 49: 50S ribosomal protein L27



- Molecule 50: 50S ribosomal protein L28



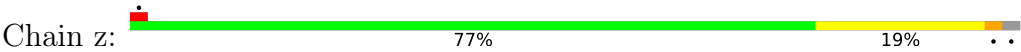
- Molecule 51: 50S ribosomal protein L29



- Molecule 52: 50S ribosomal protein L30



• Molecule 53: 50S ribosomal protein L32



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	506020	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION; per-particle CTF correction	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	40	Depositor
Minimum defocus (nm)	500	Depositor
Maximum defocus (nm)	1500	Depositor
Magnification	165000	Depositor
Image detector	FEI FALCON IV (4k x 4k)	Depositor
Maximum map value	7.947	Depositor
Minimum map value	-2.498	Depositor
Average map value	0.008	Depositor
Map value standard deviation	0.093	Depositor
Recommended contour level	0.2	Depositor
Map size (Å)	409.36002, 409.36002, 409.36002	wwPDB
Map dimensions	560, 560, 560	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.731, 0.731, 0.731	Depositor

5 Model quality

5.1 Standard geometry

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

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	0	0.29	0/424	0.37	0/565
2	1	0.37	0/380	0.42	0/498
3	2	0.34	0/513	0.39	0/676
4	3	0.34	0/303	0.35	0/397
5	A	0.33	0/35787	0.32	2/55816 (0.0%)
6	B	0.19	0/1784	0.33	0/2403
7	C	0.23	0/1651	0.31	0/2225
8	D	0.19	0/1665	0.28	0/2227
9	E	0.29	0/1165	0.34	0/1568
10	F	0.24	0/858	0.32	0/1160
11	G	0.19	0/1219	0.30	0/1635
12	H	0.27	0/989	0.31	0/1326
13	I	0.21	0/1034	0.33	0/1375
14	J	0.26	0/796	0.36	0/1077
15	K	0.30	0/884	0.34	0/1191
16	L	0.25	0/960	0.37	0/1286
17	M	0.18	0/900	0.29	0/1204
18	N	0.22	0/817	0.30	0/1088
19	O	0.25	0/722	0.24	0/964
20	P	0.26	0/653	0.33	0/877
21	Q	0.25	0/650	0.35	0/871
22	R	0.29	0/462	0.36	0/621
23	S	0.19	0/685	0.32	0/922
24	T	0.22	0/676	0.30	0/895
25	U	0.19	0/597	0.26	0/792
26	X	0.28	0/147	0.22	0/227
27	Z	0.23	0/1790	0.31	0/2787
28	a	0.47	0/65531	0.41	1/102222 (0.0%)
29	b	0.37	0/2850	0.31	0/4444
30	c	0.35	0/2121	0.39	0/2852
31	d	0.34	0/1576	0.39	0/2119

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
32	e	0.32	0/1571	0.37	0/2113
33	f	0.22	0/1434	0.30	0/1926
34	g	0.26	0/1343	0.32	0/1816
35	h	0.27	0/1112	0.40	0/1503
36	i	0.34	0/1152	0.36	0/1551
37	j	0.34	0/955	0.38	0/1279
38	k	0.34	0/1062	0.39	0/1413
39	l	0.33	0/1073	0.38	0/1433
40	m	0.35	0/958	0.42	0/1281
41	n	0.27	0/902	0.34	0/1209
42	o	0.34	0/929	0.35	0/1242
43	p	0.35	0/960	0.38	0/1278
44	q	0.34	0/829	0.39	0/1107
45	r	0.34	0/864	0.36	0/1156
46	s	0.31	0/744	0.35	0/994
47	t	0.29	0/787	0.32	0/1051
48	u	0.30	0/766	0.34	0/1025
49	v	0.32	0/593	0.35	0/785
50	w	0.34	0/635	0.37	0/848
51	x	0.28	0/502	0.25	0/667
52	y	0.32	0/453	0.32	0/605
53	z	0.35	0/450	0.42	0/599
All	All	0.38	0/150663	0.37	3/225191 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	1492	A	P-O3'-C3'	-6.86	109.91	120.20
28	a	2433	G	P-O3'-C3'	-6.85	109.93	120.20
5	A	1493	A	P-O3'-C3'	-5.44	112.04	120.20

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	0	417	0	451	5	0
2	1	377	0	418	2	0
3	2	504	0	572	4	0
4	3	302	0	340	1	0
5	A	32211	0	16218	136	0
6	B	1753	0	1780	21	0
7	C	1624	0	1696	15	0
8	D	1643	0	1707	15	0
9	E	1152	0	1196	17	0
10	F	839	0	833	10	0
11	G	1203	0	1254	25	0
12	H	979	0	1031	12	0
13	I	1022	0	1070	20	0
14	J	786	0	828	20	0
15	K	877	0	884	15	0
16	L	957	0	1017	8	0
17	M	891	0	952	19	0
18	N	805	0	844	12	0
19	O	714	0	734	5	0
20	P	643	0	661	4	0
21	Q	641	0	682	9	0
22	R	455	0	478	8	0
23	S	668	0	693	19	0
24	T	670	0	719	7	0
25	U	589	0	629	9	0
26	X	131	0	66	0	0
27	Z	1603	0	816	12	0
28	a	59025	0	29685	137	0
29	b	2549	0	1291	5	0
30	c	2082	0	2154	11	0
31	d	1566	0	1618	10	0
32	e	1552	0	1619	7	0
33	f	1410	0	1444	23	0
34	g	1323	0	1371	12	0
35	h	1101	0	1142	27	0
36	i	1129	0	1162	15	0
37	j	946	0	1023	5	0
38	k	1053	0	1129	4	0
39	l	1075	0	1145	9	0
40	m	945	0	989	2	0
41	n	892	0	923	4	0
42	o	917	0	962	5	0
43	p	947	0	1019	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
44	q	816	0	839	4	0
45	r	857	0	922	4	0
46	s	738	0	807	5	0
47	t	779	0	831	5	0
48	u	753	0	780	7	0
49	v	586	0	596	2	0
50	w	625	0	652	3	0
51	x	501	0	531	2	0
52	y	449	0	488	1	0
53	z	444	0	458	8	0
54	3	1	0	0	0	0
55	A	48	0	0	0	0
55	a	112	0	0	0	0
55	b	1	0	0	0	0
55	c	4	0	0	0	0
55	d	1	0	0	0	0
55	e	1	0	0	0	0
55	f	1	0	0	0	0
56	A	97	0	0	0	0
56	X	1	0	0	0	0
56	a	254	0	0	0	0
56	b	5	0	0	0	0
56	c	2	0	0	0	0
56	d	1	0	0	0	0
56	z	1	0	0	0	0
57	0	5	0	0	1	0
57	1	51	0	0	1	0
57	2	45	0	0	0	0
57	3	13	0	0	0	0
57	A	2886	0	0	27	0
57	B	7	0	0	0	0
57	C	16	0	0	0	0
57	D	14	0	0	0	0
57	E	38	0	0	1	0
57	F	5	0	0	0	0
57	G	4	0	0	2	0
57	H	45	0	0	0	0
57	I	11	0	0	0	0
57	J	15	0	0	0	0
57	K	22	0	0	2	0
57	L	35	0	0	2	0
57	M	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
57	N	8	0	0	0	0
57	O	29	0	0	1	0
57	P	22	0	0	0	0
57	Q	9	0	0	1	0
57	R	13	0	0	1	0
57	T	18	0	0	0	0
57	U	17	0	0	2	0
57	X	19	0	0	0	0
57	Z	20	0	0	0	0
57	a	6905	0	0	10	0
57	b	81	0	0	0	0
57	c	195	0	0	3	0
57	d	107	0	0	1	0
57	e	80	0	0	0	0
57	h	1	0	0	0	0
57	i	70	0	0	2	0
57	j	58	0	0	1	0
57	k	72	0	0	0	0
57	l	57	0	0	0	0
57	m	61	0	0	0	0
57	n	3	0	0	0	0
57	o	71	0	0	2	0
57	p	66	0	0	0	0
57	q	37	0	0	0	0
57	r	58	0	0	0	0
57	s	24	0	0	0	0
57	t	8	0	0	0	0
57	u	5	0	0	0	0
57	v	33	0	0	0	0
57	w	28	0	0	0	0
57	x	2	0	0	0	0
57	y	16	0	0	0	0
57	z	54	0	0	1	0
All	All	151507	0	94149	683	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 3.

All (683) 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)
35:h:46:PHE:CZ	35:h:50:ARG:HD3	2.12	0.84
22:R:36:SER:HA	22:R:72:ASP:HB3	1.60	0.84
5:A:664:G:H22	5:A:741:G:H1	1.30	0.80
33:f:158:THR:HG22	33:f:160:ALA:H	1.46	0.79
11:G:133:THR:HG22	11:G:134:ALA:H	1.50	0.75
5:A:505:G:N7	57:A:1804:HOH:O	2.21	0.74
9:E:101:GLU:OE1	9:E:101:GLU:N	2.17	0.74
5:A:363:A:N7	57:A:1802:HOH:O	2.20	0.74
5:A:880:C:OP1	16:L:9:ARG:NH2	2.21	0.73
28:a:2387:G:N7	57:a:3402:HOH:O	2.22	0.72
22:R:61:ARG:NH1	57:R:101:HOH:O	2.21	0.72
32:e:10:SER:OG	32:e:11:ALA:N	2.22	0.72
35:h:84:ALA:HB2	35:h:90:LEU:HD23	1.71	0.72
28:a:714:G:N7	57:a:3401:HOH:O	2.21	0.72
28:a:284:U:H3	28:a:356:G:H1	1.36	0.72
34:g:86:LYS:HG2	34:g:132:VAL:HG22	1.73	0.71
9:E:65:GLU:OE1	9:E:69:ARG:NH2	2.24	0.70
46:s:3:ARG:NH1	46:s:5:GLU:OE1	2.25	0.70
15:K:72:ASP:OD1	15:K:75:LYS:NZ	2.25	0.70
14:J:88:MET:SD	14:J:88:MET:N	2.65	0.70
14:J:7:ARG:HB2	14:J:101:SER:HB3	1.72	0.70
5:A:1483:A:N6	57:A:1807:HOH:O	2.24	0.69
28:a:1724:A:N6	28:a:1740:G:O2'	2.25	0.69
34:g:2:SER:OG	34:g:3:ARG:N	2.26	0.69
35:h:83:LYS:HE2	35:h:91:PHE:HB2	1.75	0.69
22:R:35:GLU:OE1	22:R:35:GLU:N	2.25	0.68
35:h:71:LYS:HB3	35:h:71:LYS:HZ2	1.59	0.68
28:a:1117:G:O2'	28:a:1118:G:O5'	2.11	0.68
27:Z:21:A:H61	27:Z:46:G:H2'	1.59	0.67
44:q:48:LYS:HE2	44:q:103:ALA:HB1	1.74	0.67
5:A:1458:G:N7	57:A:1809:HOH:O	2.26	0.67
15:K:37:ARG:NH1	15:K:83:GLU:OE2	2.28	0.67
5:A:1242:G:N7	57:A:1817:HOH:O	2.28	0.67
5:A:1410:A:N7	57:A:1814:HOH:O	2.28	0.66
1:O:30:LYS:NZ	28:a:2291:A:OP1	2.28	0.66
34:g:44:LYS:HB2	34:g:51:THR:HG22	1.78	0.66
36:i:9:GLU:HG2	36:i:10:THR:HG23	1.77	0.65
8:D:47:ARG:HA	8:D:47:ARG:NH1	2.11	0.65
28:a:2671:C:N3	34:g:110:SER:OG	2.30	0.65
11:G:113:ASP:OD2	11:G:122:ASN:ND2	2.30	0.65
5:A:506:G:N7	57:A:1825:HOH:O	2.30	0.65
19:O:88:ARG:NH2	28:a:716:U:OP2	2.29	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:S:44:MET:HB2	23:S:47:LEU:HD21	1.79	0.65
5:A:410:G:OP2	8:D:31:LYS:NZ	2.29	0.64
9:E:13:GLU:HG2	9:E:39:VAL:HG12	1.79	0.64
28:a:278:A:OP2	28:a:361:G:N2	2.31	0.64
36:i:1:MET:SD	36:i:2:LYS:N	2.71	0.64
28:a:2472:A:N1	57:a:3410:HOH:O	2.29	0.64
17:M:71:ARG:NH1	33:f:113:ASP:OD2	2.30	0.63
5:A:129:A:N7	57:A:1831:HOH:O	2.31	0.63
28:a:577:A:N6	57:a:3418:HOH:O	2.32	0.63
9:E:66:LYS:O	9:E:70:ASN:ND2	2.32	0.63
15:K:94:GLU:OE1	15:K:94:GLU:N	2.24	0.62
23:S:3:ARG:HH11	23:S:10:PHE:HB2	1.63	0.62
21:Q:76:VAL:HG12	21:Q:77:ARG:HG2	1.81	0.62
35:h:63:ALA:HA	35:h:66:ASN:HD22	1.63	0.62
5:A:677:U:H3	5:A:713:G:H22	1.47	0.61
5:A:1103:C:OP1	6:B:95:ARG:NH2	2.33	0.61
14:J:85:ASP:HA	14:J:88:MET:HE1	1.81	0.61
6:B:130:THR:O	6:B:132:LYS:N	2.33	0.61
20:P:48:GLU:OE1	20:P:51:ARG:NH2	2.27	0.61
14:J:66:GLU:HB3	18:N:99:ALA:HB2	1.82	0.61
5:A:187:G:N2	5:A:190:A:OP2	2.32	0.61
5:A:350:G:N7	57:A:1827:HOH:O	2.30	0.60
11:G:47:LEU:HD13	11:G:58:GLU:HB3	1.82	0.60
28:a:2802:U:H4'	28:a:2803:G:H5'	1.84	0.60
15:K:13:ARG:NH2	15:K:76:GLU:OE1	2.33	0.60
5:A:823:C:HO2'	12:H:2:SER:N	2.00	0.60
36:i:125:TYR:OH	36:i:132:HIS:NE2	2.35	0.60
48:u:9:ARG:HG2	48:u:41:GLU:HG2	1.83	0.60
5:A:1391:U:H2'	5:A:1392:G:C8	2.36	0.59
5:A:1001:C:H2'	5:A:1002:G:H8	1.67	0.59
5:A:1357:A:N7	57:A:1840:HOH:O	2.32	0.59
32:e:144:GLU:HG2	32:e:166:LYS:HE3	1.85	0.59
17:M:90:ARG:HB2	17:M:97:VAL:HG12	1.85	0.59
7:C:24:ALA:HB1	7:C:28:GLU:HG2	1.85	0.59
6:B:129:LEU:HB2	6:B:134:ALA:HB2	1.85	0.59
9:E:88:VAL:HG22	9:E:93:ARG:HG2	1.84	0.59
13:I:47:VAL:HG13	13:I:80:ARG:HD3	1.85	0.59
23:S:35:SER:HG	23:S:38:SER:HG	1.50	0.59
29:b:36:C:N4	29:b:49:C:O2	2.36	0.59
28:a:2473:A:N6	28:a:2485:G:O2'	2.36	0.58
5:A:405:U:OP2	8:D:3:ARG:NH1	2.36	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A:127:G:N7	57:A:1844:HOH:O	2.32	0.58
5:A:1328:C:H5''	17:M:28:THR:HG21	1.85	0.58
2:1:41:ARG:NH2	57:1:101:HOH:O	2.34	0.58
7:C:130:PHE:O	7:C:134:MET:HG3	2.02	0.58
28:a:2309:U:H5''	33:f:131:GLY:HA3	1.84	0.58
6:B:129:LEU:HD13	6:B:133:GLU:HG3	1.86	0.58
33:f:175:PHE:O	33:f:177:PHE:N	2.36	0.58
14:J:28:THR:HG21	14:J:90:LEU:HD22	1.85	0.58
44:q:26:ASP:O	44:q:27:ILE:HD13	2.04	0.58
36:i:39:LYS:NZ	57:i:202:HOH:O	2.37	0.58
14:J:11:LYS:HB2	14:J:97:ASP:HB3	1.86	0.57
33:f:108:VAL:HG11	33:f:176:PRO:HG2	1.86	0.57
17:M:66:GLU:OE1	17:M:70:ARG:NH2	2.36	0.57
35:h:16:GLY:HA2	35:h:47:PHE:CE2	2.39	0.57
39:l:136:MET:HG2	48:u:77:VAL:HG12	1.86	0.57
15:K:17:SER:HA	15:K:79:ILE:HA	1.86	0.57
28:a:2803:G:O2'	28:a:2804:A:H5''	2.04	0.57
13:I:25:ASN:OD1	13:I:27:LYS:NZ	2.36	0.57
5:A:1040:U:H2'	5:A:1041:G:H8	1.70	0.57
33:f:140:GLU:N	33:f:140:GLU:OE1	2.38	0.57
5:A:1305:G:N2	5:A:1331:G:H1'	2.21	0.56
11:G:133:THR:O	11:G:135:VAL:N	2.37	0.56
17:M:17:ILE:O	17:M:20:THR:OG1	2.23	0.56
5:A:458:U:H3	5:A:474:G:H1	1.53	0.56
28:a:883:G:H1	28:a:897:U:H3	1.53	0.56
41:n:76:LYS:O	41:n:80:GLU:HG3	2.04	0.56
5:A:538:G:H5''	16:L:111:LYS:HB2	1.88	0.56
21:Q:11:ARG:HH11	21:Q:11:ARG:HB2	1.69	0.56
45:r:59:GLU:OE1	45:r:66:ILE:HB	2.06	0.56
33:f:162:SER:OG	33:f:165:GLU:OE1	2.24	0.56
42:o:27:GLU:HG2	57:o:208:HOH:O	2.05	0.56
11:G:93:PRO:HA	11:G:96:ARG:HD2	1.88	0.56
30:c:165:VAL:HG21	30:c:181:MET:HE1	1.88	0.56
5:A:1003:G:N2	5:A:1037:C:O2	2.33	0.55
14:J:40:ILE:HB	14:J:73:LEU:HB3	1.87	0.55
23:S:41:PHE:H	23:S:44:MET:HE1	1.71	0.55
1:0:32:GLU:OE1	1:0:32:GLU:N	2.23	0.55
15:K:130:VAL:HA	57:K:318:HOH:O	2.05	0.55
21:Q:11:ARG:HB2	21:Q:11:ARG:NH1	2.21	0.55
5:A:161:A:H2'	5:A:162:A:C8	2.41	0.55
9:E:162:GLU:OE1	9:E:162:GLU:N	2.37	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A:1119:C:OP1	13:I:85:ARG:NH1	2.38	0.55
25:U:55:ARG:NH2	57:U:101:HOH:O	2.38	0.55
28:a:2551:A:H2'	28:a:2552:U:C6	2.42	0.55
5:A:131:A:H2'	5:A:132:C:C6	2.42	0.54
5:A:619:U:H4'	8:D:128:ARG:HH12	1.72	0.54
8:D:105:MET:HG2	8:D:173:VAL:HG22	1.88	0.54
28:a:786:G:H5'	28:a:787:G:OP1	2.07	0.54
5:A:459:A:H2'	5:A:460:A:H8	1.71	0.54
17:M:79:ARG:NE	23:S:65:GLU:OE1	2.26	0.54
5:A:459:A:H2'	5:A:460:A:C8	2.43	0.54
5:A:1422:G:N7	57:A:1847:HOH:O	2.33	0.54
13:I:97:GLU:OE1	13:I:97:GLU:N	2.40	0.54
39:l:77:PRO:HG2	39:l:80:VAL:HG21	1.88	0.54
5:A:1088:G:H4'	25:U:70:LEU:HD13	1.88	0.54
30:c:29:PRO:HG2	30:c:34:LEU:HD11	1.89	0.54
10:F:41:ASP:OD1	10:F:58:HIS:NE2	2.41	0.54
48:u:1:MET:SD	48:u:2:PHE:N	2.80	0.54
15:K:35:THR:HG22	15:K:41:ALA:HA	1.88	0.54
24:T:4:ILE:HG22	24:T:7:ALA:H	1.73	0.54
34:g:84:THR:HG23	34:g:134:LYS:HD3	1.90	0.54
50:w:62:LYS:NZ	50:w:70:GLU:OE2	2.41	0.54
5:A:946:A:H2'	5:A:947:G:C8	2.42	0.54
7:C:35:SER:OG	7:C:59:ARG:NH2	2.41	0.54
28:a:1436:A:H2'	28:a:1437:G:C8	2.43	0.54
10:F:38:ARG:NH1	10:F:98:GLU:O	2.39	0.53
13:I:85:ARG:HA	13:I:88:MET:HE2	1.90	0.53
31:d:55:LYS:HD3	31:d:60:VAL:HG22	1.89	0.53
37:j:43:ILE:HD12	37:j:56:ASP:HB2	1.89	0.53
5:A:673:A:H2'	5:A:674:G:C8	2.43	0.53
10:F:24:ARG:HG3	10:F:24:ARG:HH11	1.73	0.53
14:J:40:ILE:HG13	14:J:73:LEU:HD23	1.90	0.53
23:S:3:ARG:HH21	23:S:7:LYS:HB3	1.74	0.53
5:A:1098:C:O2'	25:U:71:TYR:O	2.25	0.53
4:3:16:ILE:HD13	4:3:25:VAL:HG22	1.90	0.53
57:a:8324:HOH:O	45:r:86:MET:HG2	2.08	0.53
49:v:59:LEU:HD12	49:v:80:ILE:HD12	1.90	0.53
7:C:134:MET:HE2	7:C:168:TYR:HD2	1.73	0.53
18:N:89:MET:HE2	18:N:89:MET:HA	1.90	0.53
19:O:62:GLN:NE2	57:O:101:HOH:O	2.41	0.53
5:A:481:G:O2'	5:A:483:C:N4	2.39	0.53
14:J:10:LEU:HB2	14:J:72:ARG:HB2	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
32:e:123:LYS:HE3	32:e:124:PHE:H	1.74	0.53
10:F:22:ILE:O	10:F:26:THR:HG23	2.09	0.52
13:I:84:THR:HG23	13:I:98:LEU:HD13	1.92	0.52
17:M:7:ILE:HA	33:f:112:ARG:HH12	1.73	0.52
28:a:1484:G:H1'	28:a:1511:A:H61	1.74	0.52
37:j:108:ARG:NH1	57:j:202:HOH:O	2.33	0.52
5:A:769:G:H4'	5:A:1513:A:H4'	1.90	0.52
11:G:5:ARG:NH1	11:G:7:ILE:O	2.43	0.52
27:Z:62:C:H2'	27:Z:63:G:C8	2.45	0.52
47:t:52:LEU:O	47:t:54:GLN:NE2	2.43	0.52
16:L:110:ARG:HB3	16:L:119:VAL:HG21	1.91	0.52
34:g:80:THR:OG1	34:g:81:GLU:N	2.41	0.52
14:J:42:LEU:HB2	14:J:71:LEU:HB3	1.91	0.52
39:l:53:MET:HE1	39:l:103:TYR:CG	2.44	0.52
5:A:1152:A:OP1	14:J:70:HIS:ND1	2.37	0.52
5:A:1323:G:H2'	5:A:1324:A:C8	2.44	0.52
9:E:164:ILE:O	12:H:114:ARG:NH2	2.42	0.52
28:a:12:U:H2'	28:a:12:U:O2	2.09	0.52
5:A:976:G:OP2	5:A:1358:U:O2'	2.28	0.51
5:A:1530:G:H2'	5:A:1531:A:C8	2.45	0.51
30:c:120:VAL:HB	35:h:91:PHE:HB3	1.92	0.51
42:o:63:LYS:NZ	57:o:203:HOH:O	2.42	0.51
45:r:4:ILE:HG12	45:r:106:VAL:HG22	1.92	0.51
53:z:52:ARG:CZ	53:z:54:VAL:HG12	2.40	0.51
5:A:382:A:H2'	5:A:383:A:C8	2.45	0.51
11:G:35:LYS:NZ	57:G:201:HOH:O	2.43	0.51
11:G:86:GLN:HB2	11:G:148:ASN:ND2	2.25	0.51
23:S:17:LYS:O	23:S:21:LYS:HG2	2.10	0.51
27:Z:23:C:H2'	27:Z:24:U:C6	2.45	0.51
45:r:73:LYS:HB2	45:r:106:VAL:HB	1.92	0.51
5:A:203:G:O2'	5:A:465:A:N1	2.42	0.51
5:A:626:G:OP1	20:P:35:ARG:NH2	2.43	0.51
8:D:163:GLU:OE1	8:D:163:GLU:N	2.38	0.51
11:G:80:VAL:O	11:G:82:GLY:N	2.44	0.51
28:a:363:G:H2'	28:a:364:C:C6	2.45	0.51
39:l:136:MET:HE2	48:u:57:TYR:CD1	2.45	0.51
46:s:30:ILE:HD13	46:s:93:LEU:HD23	1.91	0.51
9:E:151:GLU:OE2	9:E:151:GLU:N	2.31	0.51
23:S:3:ARG:NH1	23:S:10:PHE:HB2	2.25	0.51
28:a:139:U:H5''	28:a:140:C:H5	1.75	0.51
28:a:654:U:OP1	28:a:656:A:N6	2.44	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:a:1738:U:H2'	28:a:1739:G:O4'	2.10	0.51
28:a:2332:A:H2'	28:a:2333:U:C6	2.46	0.51
42:o:63:LYS:HE2	42:o:65:SER:HB2	1.91	0.51
6:B:139:ARG:HA	6:B:139:ARG:NE	2.26	0.51
5:A:337:G:H2'	5:A:338:A:C8	2.46	0.51
11:G:129:GLU:HB2	11:G:131:LYS:HD3	1.93	0.51
14:J:21:ALA:HB1	14:J:92:LEU:HD13	1.92	0.51
33:f:158:THR:HG22	33:f:160:ALA:N	2.22	0.50
35:h:7:ASP:OD1	35:h:8:LYS:N	2.41	0.50
6:B:148:LEU:HD22	6:B:151:ILE:HD11	1.94	0.50
17:M:10:PRO:HG2	17:M:13:LYS:HE2	1.92	0.50
27:Z:62:C:H2'	27:Z:63:G:H8	1.76	0.50
28:a:1407:U:H2'	28:a:1408:U:C6	2.46	0.50
28:a:1798:U:H2'	28:a:1799:G:H8	1.77	0.50
5:A:1060:U:H4'	14:J:53:ILE:HG13	1.94	0.50
6:B:85:LEU:O	6:B:87:CYS:N	2.43	0.50
12:H:38:ASN:O	12:H:42:GLU:HG3	2.11	0.50
5:A:1043:G:O2'	5:A:1044:A:O5'	2.30	0.50
21:Q:17:MET:HB2	21:Q:20:SER:HB2	1.92	0.50
28:a:1798:U:H2'	28:a:1799:G:C8	2.47	0.50
28:a:1936:A:H2'	28:a:1937:G:O4'	2.12	0.50
28:a:1029:A:C2	28:a:2492:G:H5'	2.47	0.50
28:a:2024:A:H5'	53:z:9:THR:CG2	2.42	0.50
11:G:133:THR:HG22	11:G:134:ALA:N	2.24	0.50
28:a:1735:G:H2'	28:a:1736:G:H8	1.76	0.49
35:h:72:ILE:HG13	35:h:108:VAL:HG11	1.92	0.49
28:a:2580:G:N9	57:a:3464:HOH:O	2.41	0.49
28:a:1049:G:N2	28:a:1112:G:O2'	2.43	0.49
13:I:49:ARG:HB3	13:I:53:GLU:OE2	2.12	0.49
13:I:52:LEU:HD13	13:I:58:VAL:HA	1.94	0.49
10:F:69:GLU:OE1	10:F:69:GLU:N	2.43	0.49
5:A:1356:G:H2'	5:A:1357:A:C8	2.48	0.49
28:a:1372:C:H2'	28:a:1373:G:O4'	2.12	0.49
29:b:106:G:H2'	29:b:107:G:O4'	2.12	0.49
21:Q:25:ILE:HB	21:Q:42:THR:HG23	1.94	0.49
23:S:3:ARG:NH2	23:S:7:LYS:HE2	2.27	0.49
28:a:282:A:H2'	28:a:283:G:H8	1.77	0.49
33:f:94:GLU:O	33:f:98:GLU:HG3	2.12	0.49
5:A:147:G:H2'	5:A:148:G:C8	2.47	0.49
17:M:58:ASP:OD1	17:M:58:ASP:N	2.43	0.49
28:a:2247:U:H2'	28:a:2248:U:C6	2.48	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A:928:G:O2'	5:A:1533:C:OP1	2.28	0.49
13:I:31:ASN:HD21	13:I:67:VAL:HG22	1.78	0.49
11:G:14:PRO:HB2	11:G:19:GLY:HA2	1.95	0.48
12:H:50:LYS:NZ	12:H:60:GLU:OE2	2.46	0.48
14:J:53:ILE:HD11	14:J:61:ALA:HB1	1.95	0.48
24:T:54:MET:HE2	24:T:79:LEU:HD12	1.94	0.48
19:O:17:ARG:NE	19:O:18:ASP:OD2	2.32	0.48
28:a:1143:U:H4'	28:a:1144:A:O4'	2.13	0.48
5:A:269:C:H2'	5:A:270:A:C8	2.47	0.48
15:K:93:ARG:NH2	15:K:112:ASP:OD2	2.37	0.48
35:h:132:PHE:HB2	35:h:140:ALA:HB3	1.95	0.48
12:H:48:ASP:OD1	12:H:49:PHE:N	2.46	0.48
17:M:71:ARG:HH12	33:f:113:ASP:CG	2.19	0.48
12:H:39:VAL:O	12:H:43:GLU:HG2	2.14	0.48
16:L:94:ARG:NH2	57:L:202:HOH:O	2.46	0.48
27:Z:51:C:H42	27:Z:63:G:H1	1.61	0.48
5:A:1013:G:N2	5:A:1016:A:OP2	2.44	0.48
28:a:2434:A:H2'	28:a:2434:A:N3	2.28	0.48
18:N:46:LEU:HD13	23:S:13:LEU:HD13	1.96	0.48
47:t:49:VAL:O	47:t:54:GLN:HG3	2.13	0.48
5:A:230:G:N3	57:A:1862:HOH:O	2.35	0.48
5:A:1493:A:H1'	28:a:1917:A:N6	2.29	0.48
31:d:16:THR:OG1	31:d:18:ASP:OD1	2.28	0.48
28:a:590:U:H2'	28:a:591:U:C6	2.48	0.47
42:o:38:LYS:HB3	42:o:38:LYS:HE2	1.67	0.47
5:A:1034:G:O2'	5:A:1035:A:H5'	2.13	0.47
5:A:1162:C:H2'	5:A:1163:A:C8	2.49	0.47
5:A:1314:C:H2'	5:A:1315:U:C6	2.49	0.47
28:a:2495:U:H4'	28:a:2496:U:OP1	2.14	0.47
31:d:183:GLU:H	31:d:183:GLU:CD	2.19	0.47
5:A:1124:G:N2	5:A:1125:U:O4	2.33	0.47
7:C:22:TRP:HB3	7:C:59:ARG:HB2	1.95	0.47
28:a:1140:G:N3	36:i:108:MET:HE2	2.29	0.47
28:a:271:G:O2'	28:a:272:A:O5'	2.30	0.47
28:a:932:G:H1'	52:y:25:LEU:HD11	1.96	0.47
5:A:1530:G:O2'	57:A:1803:HOH:O	2.20	0.47
9:E:56:VAL:HB	9:E:57:PRO:HD3	1.96	0.47
28:a:1185:U:H2'	28:a:1186:U:C6	2.49	0.47
31:d:46:ARG:NH2	31:d:88:GLU:O	2.46	0.47
40:m:8:ARG:HD2	40:m:43:GLU:HG2	1.96	0.47
3:2:62:LEU:HB3	3:2:65:ALA:HB2	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:a:84:A:N1	28:a:98:G:O2'	2.42	0.47
28:a:935:A:H5'	28:a:936:U:OP2	2.14	0.47
5:A:244:U:O4	5:A:906:A:H1'	2.15	0.47
5:A:977:A:O2'	5:A:979:C:OP2	2.31	0.47
5:A:1417:G:H1'	57:A:1807:HOH:O	2.14	0.47
17:M:83:LEU:HD22	23:S:74:PHE:HE1	1.78	0.47
19:O:6:GLU:CD	19:O:6:GLU:H	2.23	0.47
28:a:813:U:H2'	38:k:21:ARG:HA	1.97	0.47
28:a:1484:G:H1'	28:a:1511:A:N6	2.28	0.47
28:a:1780:U:H2'	28:a:1786:A:N6	2.30	0.47
28:a:2331:A:H2'	28:a:2332:A:C8	2.50	0.47
34:g:27:LYS:NZ	34:g:32:GLU:OE2	2.44	0.47
34:g:63:ALA:O	34:g:67:THR:HG22	2.15	0.47
34:g:9:VAL:HG22	34:g:50:LEU:HB2	1.95	0.47
35:h:46:PHE:CZ	35:h:50:ARG:CD	2.93	0.47
39:l:6:ARG:HG3	39:l:6:ARG:HH11	1.80	0.47
35:h:62:LEU:O	35:h:66:ASN:ND2	2.48	0.47
35:h:65:ALA:HB1	35:h:135:HIS:HB2	1.97	0.47
51:x:19:LEU:O	51:x:23:ARG:HG3	2.14	0.47
5:A:741:G:N7	57:A:1866:HOH:O	2.36	0.47
5:A:910:C:OP2	16:L:18:LYS:NZ	2.48	0.47
35:h:94:ILE:HB	35:h:122:LEU:HB2	1.96	0.47
35:h:97:ARG:NH1	35:h:101:ASP:OD1	2.48	0.47
5:A:472:U:H2'	5:A:473:U:C6	2.50	0.46
15:K:127:LYS:HE2	15:K:127:LYS:HB2	1.58	0.46
35:h:135:HIS:HB3	35:h:138:VAL:HG12	1.97	0.46
39:l:111:GLU:CD	39:l:111:GLU:H	2.23	0.46
15:K:114:THR:HG21	25:U:32:VAL:HG21	1.97	0.46
28:a:172:A:H2'	28:a:173:A:C8	2.49	0.46
5:A:1493:A:O2'	5:A:1494:G:H8	1.98	0.46
12:H:11:LEU:HD22	12:H:75:ILE:HD11	1.96	0.46
28:a:1592:A:H2'	28:a:1593:A:C8	2.50	0.46
35:h:71:LYS:HE3	35:h:108:VAL:HG13	1.97	0.46
7:C:73:PRO:HG3	7:C:105:GLU:OE1	2.15	0.46
30:c:107:PRO:HD2	30:c:110:LEU:HD22	1.98	0.46
33:f:14:LYS:HE2	33:f:14:LYS:HB2	1.81	0.46
10:F:90:MET:HE1	22:R:23:TYR:CZ	2.50	0.46
13:I:129:LYS:NZ	27:Z:34:U:OP2	2.41	0.46
18:N:64:CYS:HB2	18:N:80:SER:HB3	1.97	0.46
28:a:282:A:H2'	28:a:283:G:C8	2.50	0.46
28:a:1609:C:N4	28:a:1624:G:OP2	2.39	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A:1491:G:H2'	5:A:1492:A:O4'	2.16	0.46
28:a:859:G:H2'	28:a:860:G:O4'	2.15	0.46
5:A:381:C:H2'	5:A:382:A:O4'	2.16	0.46
5:A:1225:A:H2'	5:A:1226:C:C5	2.50	0.46
7:C:134:MET:HE2	7:C:168:TYR:CD2	2.51	0.46
10:F:88:MET:SD	10:F:90:MET:HE2	2.55	0.46
20:P:56:ARG:HD2	20:P:56:ARG:HA	1.62	0.46
6:B:128:LYS:HE3	6:B:129:LEU:HD23	1.97	0.46
30:c:17:VAL:HG12	57:c:498:HOH:O	2.15	0.46
9:E:109:GLY:N	57:E:204:HOH:O	2.49	0.46
21:Q:8:LEU:HB3	21:Q:25:ILE:HD12	1.97	0.46
5:A:275:G:H5'	21:Q:16:LYS:HD3	1.97	0.46
28:a:542:C:H2'	28:a:543:A:C8	2.51	0.46
28:a:898:A:C6	28:a:899:C:H1'	2.50	0.46
36:i:13:ARG:NH2	36:i:49:ASP:O	2.44	0.46
5:A:1287:A:H2'	5:A:1288:A:C8	2.51	0.45
8:D:174:ASP:OD2	8:D:177:LYS:NZ	2.29	0.45
9:E:26:LYS:HB2	9:E:26:LYS:HE3	1.79	0.45
14:J:32:THR:O	14:J:83:THR:OG1	2.35	0.45
28:a:1722:U:H2'	28:a:1723:G:O4'	2.16	0.45
29:b:14:U:OP2	29:b:70:C:O2'	2.32	0.45
3:2:27:ALA:O	3:2:28:ASN:HB2	2.17	0.45
14:J:15:HIS:HB3	14:J:70:HIS:CE1	2.50	0.45
15:K:130:VAL:HG12	57:K:318:HOH:O	2.14	0.45
5:A:188:C:H2'	5:A:189:A:O4'	2.16	0.45
5:A:260:G:N3	57:A:1864:HOH:O	2.35	0.45
9:E:151:GLU:H	9:E:151:GLU:CD	2.22	0.45
27:Z:25:C:H2'	27:Z:26:G:O4'	2.16	0.45
28:a:276:U:H2'	28:a:277:G:C8	2.52	0.45
28:a:2595:C:H2'	28:a:2596:G:C8	2.51	0.45
53:z:36:GLU:OE2	53:z:46:ASP:HB2	2.16	0.45
5:A:460:A:H2'	5:A:461:A:H8	1.82	0.45
5:A:1302:C:C5	17:M:17:ILE:HG13	2.51	0.45
6:B:86:SER:O	6:B:86:SER:OG	2.26	0.45
28:a:2295:U:H2'	28:a:2296:U:C6	2.51	0.45
32:e:7:ASP:CG	32:e:122:GLU:H	2.24	0.45
33:f:47:LYS:HE2	33:f:47:LYS:HB2	1.57	0.45
34:g:3:ARG:HE	34:g:3:ARG:HB2	1.56	0.45
10:F:18:VAL:O	10:F:22:ILE:HG13	2.16	0.45
28:a:1388:C:H2'	28:a:1389:A:C8	2.52	0.45
47:t:81:ASP:OD2	47:t:98:SER:OG	2.32	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:B:12:ALA:HA	6:B:208:ARG:HH11	1.81	0.45
8:D:48:LEU:HD22	8:D:52:GLY:HA3	1.98	0.45
53:z:55:ILE:HG13	53:z:56:ALA:O	2.16	0.45
5:A:202:G:H1	5:A:215:C:H42	1.65	0.45
5:A:1001:C:H2'	5:A:1002:G:C8	2.50	0.45
5:A:1180:A:OP1	13:I:105:THR:OG1	2.33	0.45
16:L:110:ARG:NH1	57:L:201:HOH:O	2.35	0.45
17:M:7:ILE:HG22	33:f:112:ARG:NH1	2.32	0.45
28:a:372:G:O2'	50:w:54:LYS:HE3	2.16	0.45
28:a:595:U:H2'	28:a:596:U:C6	2.52	0.45
28:a:879:A:H2	28:a:902:A:N7	2.15	0.45
35:h:42:LYS:O	35:h:43:ASN:C	2.59	0.45
35:h:113:SER:O	35:h:116:ARG:NH2	2.50	0.45
5:A:301:G:N7	57:A:1867:HOH:O	2.36	0.45
11:G:68:ASN:O	11:G:138:ARG:NH2	2.49	0.45
22:R:21:ILE:HG21	22:R:55:LEU:HD23	1.98	0.45
28:a:2906:C:H2'	28:a:2907:U:O4'	2.16	0.45
5:A:1218:C:H2'	5:A:1219:A:C8	2.51	0.45
7:C:110:GLU:HB2	7:C:144:LEU:HD22	1.99	0.45
8:D:187:GLU:N	8:D:190:ASP:OD2	2.49	0.45
28:a:2784:G:OP2	36:i:120:ARG:HD3	2.17	0.45
21:Q:39:LYS:NZ	57:Q:101:HOH:O	2.50	0.45
28:a:725:C:H2'	28:a:726:U:O4'	2.17	0.45
28:a:858:G:H2'	28:a:859:G:C8	2.52	0.45
35:h:96:THR:HG22	35:h:117:LEU:CD1	2.47	0.45
5:A:1371:G:OP1	13:I:14:SER:OG	2.35	0.44
9:E:81:LEU:HB3	9:E:147:MET:SD	2.57	0.44
21:Q:21:ILE:HG13	21:Q:46:VAL:HB	1.98	0.44
33:f:111:ILE:HB	33:f:114:PHE:HB2	1.99	0.44
5:A:813:U:H2'	57:A:3394:HOH:O	2.17	0.44
24:T:86:LEU:HD23	24:T:86:LEU:HA	1.84	0.44
32:e:147:LEU:HD11	32:e:170:ARG:HG3	2.00	0.44
43:p:49:ASP:HA	43:p:52:GLN:HB2	1.99	0.44
11:G:56:LYS:HB2	11:G:56:LYS:HE2	1.79	0.44
11:G:71:PRO:O	11:G:96:ARG:HG2	2.17	0.44
12:H:29:SER:HB3	12:H:57:PRO:HB2	1.99	0.44
27:Z:17:C:OP2	27:Z:17(A):U:O2'	2.35	0.44
28:a:326:G:N7	57:a:3432:HOH:O	2.36	0.44
32:e:137:LYS:O	32:e:141:MET:HG3	2.16	0.44
5:A:1162:C:H2'	5:A:1163:A:H8	1.81	0.44
6:B:4:VAL:HG21	6:B:212:LEU:HD21	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:a:493:G:H2'	28:a:494:G:O4'	2.18	0.44
28:a:1183:U:H2'	28:a:1184:G:C8	2.53	0.44
38:k:10:GLU:H	38:k:10:GLU:CD	2.25	0.44
1:0:38:LYS:HB3	1:0:38:LYS:HE3	1.77	0.44
5:A:476:U:H2'	5:A:477:C:C6	2.52	0.44
7:C:50:ALA:O	7:C:70:THR:OG1	2.35	0.44
13:I:28:ILE:HG12	13:I:63:LEU:HD22	1.99	0.44
30:c:5:LYS:HB3	30:c:5:LYS:HE3	1.68	0.44
35:h:110:VAL:HG11	35:h:132:PHE:CD2	2.53	0.44
11:G:95:ARG:NH1	57:G:202:HOH:O	2.45	0.44
28:a:191:A:H2'	28:a:192:C:C6	2.53	0.44
5:A:216:U:H4'	5:A:464:U:H4'	1.99	0.44
5:A:736:C:H2'	5:A:737:C:C6	2.53	0.44
5:A:1004:A:H2'	5:A:1005:A:O4'	2.18	0.44
5:A:1086:U:H3	5:A:1099:G:H22	1.66	0.44
11:G:16:PRO:HD2	11:G:44:TYR:OH	2.18	0.44
28:a:358:U:H2'	28:a:359:G:C8	2.53	0.44
31:d:99:GLU:HG2	31:d:182:ALA:HB2	2.00	0.44
33:f:36:LEU:HD22	33:f:154:ILE:HG12	1.99	0.44
34:g:2:SER:O	34:g:6:LYS:HG2	2.17	0.44
47:t:50:PRO:HA	47:t:54:GLN:HE21	1.83	0.44
37:j:38:ILE:HD11	37:j:112:PHE:CZ	2.53	0.44
43:p:58:ARG:HA	43:p:61:TRP:CE3	2.53	0.44
5:A:518:C:O2'	5:A:530:G:N2	2.50	0.44
5:A:1019:A:H2'	5:A:1020:G:O4'	2.18	0.44
6:B:74:ARG:HD3	6:B:74:ARG:HA	1.80	0.44
18:N:83:LYS:HA	18:N:83:LYS:HD2	1.65	0.44
28:a:1030:A:N6	28:a:1127:G:H2'	2.33	0.44
28:a:2237:U:H2'	28:a:2238:G:C8	2.53	0.44
37:j:38:ILE:HD11	37:j:112:PHE:HZ	1.82	0.44
3:2:14:PHE:O	3:2:15:LYS:HD3	2.18	0.43
5:A:1273:C:H2'	5:A:1274:A:O4'	2.18	0.43
28:a:572:G:H2'	28:a:2034:6MZ:N7	2.32	0.43
28:a:853:C:H2'	28:a:854:U:C6	2.53	0.43
48:u:55:GLU:CD	48:u:55:GLU:H	2.26	0.43
5:A:555:U:H2'	5:A:556:C:C6	2.54	0.43
5:A:814:A:H8	57:A:3394:HOH:O	2.01	0.43
6:B:100:MET:HE3	6:B:100:MET:HB2	1.74	0.43
14:J:8:ILE:HB	14:J:74:VAL:HB	2.00	0.43
28:a:744:A:H2'	28:a:745:A:C8	2.54	0.43
28:a:2904:A:H2'	28:a:2905:C:O4'	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:i:11:VAL:HG21	36:i:50:THR:HA	1.99	0.43
41:n:60:GLU:HG3	41:n:61:GLN:OE1	2.18	0.43
5:A:460:A:H2'	5:A:461:A:C8	2.54	0.43
5:A:1144:G:N2	5:A:1146:A:H62	2.16	0.43
35:h:65:ALA:CB	35:h:135:HIS:HB2	2.48	0.43
35:h:115:VAL:O	35:h:116:ARG:NH1	2.51	0.43
36:i:114:LEU:HG	36:i:118:MET:HE3	2.01	0.43
11:G:37:SER:HB3	13:I:41:ARG:HD3	2.00	0.43
5:A:299:G:H2'	5:A:300:A:C8	2.53	0.43
6:B:119:THR:O	6:B:123:ASP:N	2.48	0.43
7:C:14:ILE:HG22	7:C:15:VAL:HG13	2.00	0.43
13:I:21:ILE:HD13	13:I:21:ILE:HA	1.81	0.43
28:a:1511:A:HO2'	28:a:1512:G:H8	1.64	0.43
39:l:20:LEU:HD13	48:u:81:PRO:HG2	2.01	0.43
57:A:3721:HOH:O	9:E:54:ARG:HG2	2.18	0.43
6:B:117:LEU:HB3	6:B:141:LEU:HD23	1.99	0.43
11:G:78:ARG:HD2	11:G:78:ARG:HA	1.94	0.43
15:K:80:LYS:HB3	15:K:80:LYS:HE3	1.74	0.43
17:M:4:ILE:HG23	17:M:57:ARG:HG2	2.00	0.43
28:a:646:A:H2'	28:a:647:C:O4'	2.18	0.43
41:n:83:LEU:HD11	41:n:114:GLY:HA3	2.00	0.43
3:2:54:ASP:HB3	38:k:57:LEU:HD22	2.00	0.43
15:K:24:HIS:HB3	15:K:31:ILE:HB	2.00	0.43
17:M:23:TYR:CE2	17:M:70:ARG:HG3	2.54	0.43
18:N:98:LYS:HB3	18:N:98:LYS:HE3	1.83	0.43
28:a:1849:A:O3'	28:a:1850:A:H8	2.01	0.43
28:a:2277:A:H2'	28:a:2278:A:C8	2.53	0.43
5:A:1040:U:H2'	5:A:1041:G:C8	2.51	0.43
5:A:1127:G:H5'	5:A:1280:A:O2'	2.18	0.43
5:A:1314:C:OP1	23:S:6:LYS:NZ	2.46	0.43
8:D:170:TRP:CD2	8:D:186:PRO:HB3	2.54	0.43
11:G:80:VAL:HG21	11:G:154:TYR:HE2	1.84	0.43
22:R:21:ILE:N	22:R:21:ILE:HD13	2.34	0.43
29:b:1:U:H2'	29:b:2:G:C8	2.54	0.43
30:c:161:TYR:HB3	30:c:194:GLU:HG2	2.00	0.43
5:A:1216:A:H5''	18:N:5:SER:HB3	2.00	0.43
9:E:34:THR:HG22	9:E:52:LYS:HG2	2.00	0.43
18:N:40:ASP:OD1	18:N:41:ARG:N	2.50	0.43
28:a:731:G:C6	30:c:207:LYS:HB2	2.53	0.43
28:a:1705:G:H2'	28:a:1706:C:C6	2.54	0.43
28:a:1746:A:H3'	28:a:1747:A:H8	1.84	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:a:1810:A:H3'	28:a:1811:A:C8	2.54	0.43
35:h:46:PHE:CE1	35:h:50:ARG:HD3	2.53	0.43
49:v:25:ARG:HD2	49:v:25:ARG:HA	1.80	0.43
11:G:74:GLU:O	11:G:88:PRO:HA	2.19	0.43
23:S:16:LEU:O	23:S:20:GLU:HG2	2.19	0.43
28:a:641:U:H2'	28:a:642:C:C6	2.54	0.43
25:U:31:GLU:OE2	25:U:34:ARG:NH2	2.50	0.42
28:a:835:A:H2'	28:a:836:G:C8	2.54	0.42
28:a:1748:A:H2'	28:a:1749:U:C6	2.54	0.42
31:d:116:LYS:NZ	57:d:402:HOH:O	2.41	0.42
36:i:95:ARG:HG2	36:i:96:ARG:HG2	2.01	0.42
28:a:659:U:H2'	28:a:660:U:C6	2.53	0.42
32:e:41:GLN:HG2	32:e:43:THR:HG23	2.01	0.42
33:f:130:MET:HE3	33:f:130:MET:HB2	1.78	0.42
36:i:110:PRO:O	36:i:115:GLY:HA3	2.20	0.42
38:k:132:ARG:HG3	38:k:142:ILE:HD13	2.00	0.42
5:A:1227:A:O2'	17:M:114:LYS:HG3	2.19	0.42
5:A:1320:C:O2	23:S:36:ARG:NH2	2.45	0.42
5:A:1492:A:H2'	5:A:1493:A:C1'	2.49	0.42
12:H:43:GLU:HG3	12:H:101:ILE:HD13	2.02	0.42
28:a:2792:C:H2'	28:a:2793:C:C6	2.54	0.42
30:c:39:LYS:NZ	57:c:415:HOH:O	2.52	0.42
28:a:1541:U:H2'	28:a:1542:G:H8	1.84	0.42
28:a:2306:U:H2'	28:a:2307:G:C8	2.55	0.42
34:g:164:TYR:HB2	34:g:167:GLU:HB2	2.01	0.42
5:A:48:C:OP1	57:A:1805:HOH:O	2.21	0.42
5:A:1120:C:H2'	5:A:1121:U:C6	2.55	0.42
6:B:139:ARG:HA	6:B:139:ARG:CZ	2.50	0.42
14:J:47:GLU:OE1	18:N:76:LYS:NZ	2.40	0.42
20:P:12:LYS:HG2	20:P:13:LYS:HG2	2.02	0.42
27:Z:44:A:H2'	27:Z:45:G:C8	2.55	0.42
46:s:89:GLU:OE1	46:s:89:GLU:HA	2.19	0.42
1:O:38:LYS:NZ	57:O:102:HOH:O	2.52	0.42
57:A:1819:HOH:O	25:U:40:LYS:HE3	2.20	0.42
12:H:89:LYS:HE3	12:H:89:LYS:HB3	1.76	0.42
23:S:44:MET:HB2	23:S:47:LEU:CD2	2.48	0.42
31:d:74:GLU:H	31:d:74:GLU:CD	2.26	0.42
35:h:110:VAL:HG11	35:h:132:PHE:CE2	2.54	0.42
10:F:25:TYR:O	10:F:29:ILE:HD12	2.20	0.42
12:H:41:LYS:HD2	12:H:48:ASP:HA	2.02	0.42
5:A:214:C:H2'	5:A:215:C:H6	1.85	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A:1492:A:OP1	16:L:44:LYS:HE2	2.20	0.42
6:B:73:LYS:NZ	6:B:205:ASP:OD1	2.52	0.42
14:J:25:ILE:HD13	14:J:25:ILE:HA	1.82	0.42
23:S:80:TYR:CZ	23:S:82:GLY:HA2	2.55	0.42
25:U:34:ARG:NH2	57:U:103:HOH:O	2.52	0.42
27:Z:69:C:H2'	27:Z:70:G:C8	2.55	0.42
33:f:57:LEU:O	33:f:61:SER:OG	2.34	0.42
47:t:39:ILE:HG22	47:t:40:ASN:N	2.35	0.42
5:A:640:A:N7	57:A:1877:HOH:O	2.37	0.42
6:B:118:GLU:OE1	6:B:152:LYS:NZ	2.48	0.42
6:B:135:LEU:HA	6:B:138:THR:HG22	2.02	0.42
28:a:2:G:H2'	28:a:3:U:C6	2.54	0.42
28:a:155:A:H2'	28:a:156:A:C8	2.54	0.42
30:c:37:ASN:HB3	57:c:416:HOH:O	2.18	0.42
36:i:58:ASN:HA	36:i:126:ALA:O	2.19	0.42
5:A:1228:C:P	17:M:114:LYS:HZ3	2.43	0.42
5:A:1305:G:H22	5:A:1331:G:H1'	1.85	0.42
13:I:83:ILE:O	13:I:87:LEU:HD12	2.20	0.42
25:U:63:GLU:HA	25:U:66:ARG:HD2	2.02	0.42
28:a:2684:U:O2'	28:a:2685:C:H5'	2.19	0.41
33:f:48:LYS:HA	33:f:51:ASP:OD2	2.20	0.41
6:B:64:LYS:HA	6:B:64:LYS:HD3	1.74	0.41
11:G:47:LEU:HD23	11:G:47:LEU:HA	1.83	0.41
27:Z:50:U:H2'	27:Z:51:C:O4'	2.20	0.41
28:a:340:A:H2'	28:a:341:C:O4'	2.20	0.41
28:a:570:U:H1'	28:a:2034:6MZ:H9C1	2.03	0.41
39:l:41:LEU:HG	39:l:96:ILE:HG13	2.01	0.41
5:A:983:A:H5'	5:A:984:C:OP2	2.20	0.41
5:A:993:G:O2'	5:A:995:C:N4	2.52	0.41
28:a:880:A:N7	28:a:901:A:H2	2.18	0.41
28:a:1643:A:H2'	28:a:1644:G:O4'	2.20	0.41
31:d:80:TRP:CD1	31:d:202:ILE:HD11	2.55	0.41
33:f:126:GLY:O	33:f:158:THR:HB	2.20	0.41
5:A:449:G:H2'	5:A:450:G:C8	2.55	0.41
5:A:1220:G:P	18:N:53:ARG:HH12	2.44	0.41
12:H:27:MET:HE2	12:H:27:MET:HB3	1.94	0.41
28:a:614:G:O2'	28:a:616:A:N7	2.53	0.41
43:p:76:TYR:CZ	43:p:80:ILE:HG13	2.56	0.41
5:A:539:A:H2'	5:A:540:G:C8	2.55	0.41
5:A:1333:A:H2'	5:A:1334:G:O4'	2.20	0.41
7:C:131:ARG:HE	7:C:131:ARG:HB2	1.73	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:C:178:LEU:HD23	7:C:178:LEU:HA	1.93	0.41
28:a:567:C:N4	57:a:3418:HOH:O	2.51	0.41
28:a:1735:G:H2'	28:a:1736:G:C8	2.55	0.41
46:s:93:LEU:HD12	46:s:93:LEU:HA	1.83	0.41
5:A:1123:U:O2'	14:J:39:PRO:O	2.34	0.41
5:A:1250:A:H4'	13:I:70:GLY:N	2.35	0.41
5:A:1438:G:N7	57:A:1874:HOH:O	2.37	0.41
5:A:1478:U:H2'	5:A:1479:C:C6	2.56	0.41
8:D:152:GLN:HB2	8:D:155:VAL:HG23	2.03	0.41
11:G:87:VAL:HG13	11:G:151:PHE:O	2.21	0.41
33:f:112:ARG:HE	33:f:112:ARG:HB2	1.63	0.41
8:D:47:ARG:HA	8:D:47:ARG:HH11	1.85	0.41
13:I:97:GLU:H	13:I:97:GLU:CD	2.28	0.41
27:Z:18:G:H21	27:Z:58:A:H5'	1.86	0.41
28:a:636:C:H2'	28:a:637:C:C6	2.56	0.41
28:a:1601:U:H2'	28:a:1602:C:C6	2.56	0.41
28:a:1801:G:O2'	30:c:180:GLU:OE2	2.33	0.41
31:d:181:ASP:HB3	31:d:186:LEU:HB2	2.02	0.41
5:A:1318:A:H1'	23:S:37:ARG:NH1	2.35	0.41
11:G:47:LEU:HB3	11:G:58:GLU:HG2	2.02	0.41
24:T:53:GLU:H	24:T:53:GLU:CD	2.29	0.41
25:U:67:ARG:HD3	25:U:67:ARG:HA	1.86	0.41
28:a:647:C:H2'	28:a:649:G:C8	2.56	0.41
28:a:1329:A:H2'	28:a:1330:A:O4'	2.20	0.41
57:a:5912:HOH:O	44:q:86:GLN:HG3	2.20	0.41
36:i:101:ILE:HG13	57:i:201:HOH:O	2.19	0.41
39:l:127:LYS:HE3	39:l:127:LYS:HB2	1.68	0.41
41:n:39:VAL:HG12	41:n:48:LEU:HD12	2.01	0.41
53:z:43:ILE:HG22	53:z:49:TYR:HB2	2.02	0.41
5:A:333:U:OP1	24:T:2:ALA:N	2.54	0.41
5:A:706:A:N3	57:A:1882:HOH:O	2.37	0.41
8:D:177:LYS:HB2	8:D:177:LYS:HE2	1.75	0.41
28:a:479:A:N3	28:a:481:G:H5''	2.36	0.41
28:a:877:G:C6	28:a:878:C:C4	3.09	0.41
28:a:1452:G:C6	28:a:1453:C:N4	2.88	0.41
28:a:1796:A:H2'	28:a:1797:C:C6	2.55	0.41
28:a:1919:3TD:O5'	28:a:1919:3TD:H6	2.20	0.41
28:a:2619:U:C2	53:z:4:GLN:HA	2.55	0.41
57:a:8941:HOH:O	43:p:15:LYS:HE2	2.20	0.41
29:b:48:U:H2'	29:b:49:C:C6	2.56	0.41
36:i:104:ALA:O	36:i:108:MET:HG3	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
42:o:111:LYS:HB3	42:o:111:LYS:HE3	1.78	0.41
51:x:2:LYS:O	51:x:2:LYS:HG3	2.20	0.41
5:A:451:A:H61	5:A:481:G:H5'	1.86	0.41
5:A:1296:C:H4'	5:A:1302:C:N4	2.35	0.41
18:N:24:ARG:HD2	18:N:55:SER:OG	2.21	0.41
19:O:73:LYS:HA	19:O:73:LYS:HD2	1.84	0.41
23:S:47:LEU:H	23:S:47:LEU:HD23	1.85	0.41
28:a:1:G:H2'	28:a:2:G:C8	2.56	0.41
28:a:118:A:N3	28:a:178:G:H1'	2.37	0.41
28:a:560:U:OP1	36:i:113:PRO:HD2	2.21	0.41
28:a:654:U:P	28:a:656:A:H61	2.44	0.41
28:a:1022:A:C2	28:a:1143:U:C2	3.09	0.41
5:A:236:A:H2'	5:A:237:G:C8	2.56	0.40
5:A:251:G:C6	5:A:266:G:C6	3.10	0.40
9:E:72:ILE:HD13	9:E:145:GLU:HB2	2.03	0.40
24:T:53:GLU:CD	24:T:53:GLU:N	2.80	0.40
28:a:347:A:H2'	28:a:348:A:C8	2.56	0.40
28:a:391:A:H1'	28:a:411:G:O4'	2.20	0.40
28:a:863:A:H2'	28:a:864:G:O4'	2.22	0.40
28:a:1044:G:C6	28:a:1045:C:C4	3.09	0.40
35:h:40:THR:O	35:h:41:LYS:C	2.63	0.40
37:j:7:MET:HE3	37:j:7:MET:HB3	1.90	0.40
43:p:89:GLU:HG3	44:q:52:PRO:HB3	2.02	0.40
1:0:33:LYS:HA	1:0:33:LYS:HD2	1.82	0.40
57:A:3469:HOH:O	15:K:118:HIS:HD2	2.03	0.40
6:B:100:MET:HA	6:B:107:VAL:HG21	2.03	0.40
7:C:22:TRP:CD1	7:C:59:ARG:HD2	2.57	0.40
8:D:124:MET:HE3	8:D:127:GLY:HA2	2.02	0.40
17:M:48:LEU:HD23	17:M:48:LEU:HA	1.85	0.40
24:T:57:ILE:HD13	24:T:57:ILE:HA	1.82	0.40
28:a:247:G:H4'	28:a:386:G:C5	2.56	0.40
28:a:1184:G:H2'	28:a:1185:U:O4'	2.21	0.40
28:a:1265:U:O2'	53:z:8:PRO:HD2	2.21	0.40
46:s:71:GLY:O	46:s:72:GLN:HB3	2.21	0.40
48:u:2:PHE:CD1	48:u:50:MET:HE3	2.57	0.40
50:w:3:ARG:O	50:w:12:PRO:HD3	2.20	0.40
5:A:390:U:H2'	5:A:391:G:C8	2.57	0.40
7:C:49:LYS:HE3	7:C:49:LYS:HB3	1.81	0.40
7:C:131:ARG:HD3	7:C:168:TYR:OH	2.22	0.40
9:E:160:SER:OG	9:E:163:GLU:HG3	2.21	0.40
10:F:5:GLU:OE2	22:R:24:LYS:NZ	2.53	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
16:L:43:LYS:HE3	16:L:91:PRO:HD3	2.03	0.40
28:a:1181:G:H2'	28:a:1182:U:C6	2.57	0.40
28:a:1435:A:H2'	28:a:1436:A:O4'	2.22	0.40
35:h:8:LYS:HD2	35:h:9:VAL:N	2.36	0.40
40:m:86:ARG:NH2	40:m:117:ASP:OD1	2.43	0.40
2:l:26:ASN:ND2	28:a:684:G:H5'	2.37	0.40
5:A:1121:U:H2'	5:A:1122:U:C6	2.56	0.40
8:D:27:ALA:HB3	8:D:30:THR:HG23	2.03	0.40
11:G:16:PRO:HG3	13:I:43:THR:HG22	2.03	0.40
13:I:63:LEU:HD23	13:I:65:ILE:HD11	2.02	0.40
17:M:45:ILE:HA	17:M:48:LEU:HG	2.03	0.40
18:N:16:LEU:HD23	18:N:16:LEU:HA	1.87	0.40
22:R:22:ASP:OD2	22:R:24:LYS:HE2	2.21	0.40
23:S:28:LYS:HA	23:S:28:LYS:HD3	1.90	0.40
33:f:129:SER:OG	33:f:155:THR:HG23	2.22	0.40
53:z:5:GLN:HG3	57:z:201:HOH:O	2.21	0.40
5:A:465:A:H2'	5:A:466:A:C8	2.57	0.40
5:A:1428:A:H2'	5:A:1429:A:O4'	2.21	0.40
5:A:1477:U:H2'	5:A:1478:U:C6	2.57	0.40
5:A:1504:G:H4'	5:A:1505:G:C4	2.57	0.40
15:K:82:LEU:HD23	15:K:82:LEU:HA	1.93	0.40
28:a:360:U:C4	28:a:361:G:C6	3.10	0.40
28:a:878:C:H2'	28:a:879:A:O4'	2.21	0.40
28:a:896:U:H2'	28:a:897:U:O4'	2.22	0.40
31:d:32:ASN:HD22	31:d:32:ASN:N	2.19	0.40
33:f:5:HIS:HB2	33:f:97:TRP:CD1	2.57	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	0	49/55 (89%)	49 (100%)	0	0	100	100
2	1	44/46 (96%)	44 (100%)	0	0	100	100
3	2	62/65 (95%)	60 (97%)	2 (3%)	0	100	100
4	3	36/38 (95%)	36 (100%)	0	0	100	100
6	B	222/241 (92%)	210 (95%)	9 (4%)	3 (1%)	9	1
7	C	204/233 (88%)	193 (95%)	11 (5%)	0	100	100
8	D	203/206 (98%)	197 (97%)	6 (3%)	0	100	100
9	E	154/167 (92%)	150 (97%)	4 (3%)	0	100	100
10	F	101/135 (75%)	99 (98%)	2 (2%)	0	100	100
11	G	151/179 (84%)	136 (90%)	13 (9%)	2 (1%)	9	1
12	H	127/130 (98%)	122 (96%)	5 (4%)	0	100	100
13	I	125/130 (96%)	120 (96%)	5 (4%)	0	100	100
14	J	96/103 (93%)	88 (92%)	7 (7%)	1 (1%)	12	3
15	K	113/129 (88%)	108 (96%)	5 (4%)	0	100	100
16	L	120/124 (97%)	116 (97%)	3 (2%)	1 (1%)	16	4
17	M	113/118 (96%)	109 (96%)	4 (4%)	0	100	100
18	N	98/101 (97%)	95 (97%)	3 (3%)	0	100	100
19	O	86/89 (97%)	84 (98%)	2 (2%)	0	100	100
20	P	79/82 (96%)	77 (98%)	2 (2%)	0	100	100
21	Q	77/84 (92%)	70 (91%)	7 (9%)	0	100	100
22	R	53/75 (71%)	51 (96%)	1 (2%)	1 (2%)	6	1
23	S	82/92 (89%)	74 (90%)	8 (10%)	0	100	100
24	T	84/87 (97%)	84 (100%)	0	0	100	100
25	U	68/71 (96%)	67 (98%)	1 (2%)	0	100	100
30	c	269/273 (98%)	259 (96%)	10 (4%)	0	100	100
31	d	206/209 (99%)	200 (97%)	6 (3%)	0	100	100
32	e	199/201 (99%)	194 (98%)	5 (2%)	0	100	100
33	f	175/179 (98%)	165 (94%)	10 (6%)	0	100	100
34	g	174/177 (98%)	160 (92%)	13 (8%)	1 (1%)	21	6
35	h	146/149 (98%)	134 (92%)	10 (7%)	2 (1%)	9	1
36	i	140/142 (99%)	140 (100%)	0	0	100	100
37	j	121/123 (98%)	118 (98%)	3 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
38	k	142/144 (99%)	138 (97%)	4 (3%)	0	100	100
39	l	132/136 (97%)	127 (96%)	5 (4%)	0	100	100
40	m	116/127 (91%)	112 (97%)	4 (3%)	0	100	100
41	n	114/117 (97%)	113 (99%)	1 (1%)	0	100	100
42	o	112/115 (97%)	110 (98%)	2 (2%)	0	100	100
43	p	115/118 (98%)	114 (99%)	1 (1%)	0	100	100
44	q	101/103 (98%)	100 (99%)	1 (1%)	0	100	100
45	r	108/110 (98%)	106 (98%)	2 (2%)	0	100	100
46	s	91/100 (91%)	90 (99%)	1 (1%)	0	100	100
47	t	100/104 (96%)	96 (96%)	3 (3%)	1 (1%)	12	3
48	u	92/94 (98%)	90 (98%)	2 (2%)	0	100	100
49	v	76/85 (89%)	75 (99%)	1 (1%)	0	100	100
50	w	75/78 (96%)	75 (100%)	0	0	100	100
51	x	60/63 (95%)	60 (100%)	0	0	100	100
52	y	56/59 (95%)	55 (98%)	1 (2%)	0	100	100
53	z	54/57 (95%)	53 (98%)	1 (2%)	0	100	100
All	All	5521/5843 (94%)	5323 (96%)	186 (3%)	12 (0%)	44	24

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
6	B	87	CYS
6	B	131	LYS
11	G	81	GLY
14	J	57	VAL
34	g	126	PRO
6	B	86	SER
35	h	42	LYS
35	h	76	GLU
11	G	153	HIS
22	R	73	ARG
47	t	39	ILE
16	L	91	PRO

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	0	46/49 (94%)	46 (100%)	0	100	100
2	1	38/38 (100%)	38 (100%)	0	100	100
3	2	51/52 (98%)	51 (100%)	0	100	100
4	3	34/34 (100%)	34 (100%)	0	100	100
6	B	186/199 (94%)	180 (97%)	6 (3%)	34	8
7	C	170/190 (90%)	166 (98%)	4 (2%)	43	15
8	D	172/173 (99%)	168 (98%)	4 (2%)	44	16
9	E	119/126 (94%)	118 (99%)	1 (1%)	73	56
10	F	90/116 (78%)	89 (99%)	1 (1%)	65	43
11	G	126/147 (86%)	125 (99%)	1 (1%)	73	56
12	H	104/105 (99%)	103 (99%)	1 (1%)	68	47
13	I	105/107 (98%)	105 (100%)	0	100	100
14	J	86/90 (96%)	83 (96%)	3 (4%)	32	7
15	K	89/98 (91%)	86 (97%)	3 (3%)	32	7
16	L	102/103 (99%)	101 (99%)	1 (1%)	68	47
17	M	93/96 (97%)	91 (98%)	2 (2%)	45	18
18	N	83/84 (99%)	82 (99%)	1 (1%)	63	40
19	O	76/77 (99%)	75 (99%)	1 (1%)	61	36
20	P	65/65 (100%)	65 (100%)	0	100	100
21	Q	73/78 (94%)	72 (99%)	1 (1%)	59	33
22	R	48/65 (74%)	46 (96%)	2 (4%)	26	4
23	S	72/79 (91%)	70 (97%)	2 (3%)	38	11
24	T	65/66 (98%)	64 (98%)	1 (2%)	57	31
25	U	60/61 (98%)	57 (95%)	3 (5%)	22	2
30	c	216/218 (99%)	216 (100%)	0	100	100
31	d	163/163 (100%)	163 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
32	e	165/165 (100%)	162 (98%)	3 (2%)	51	24
33	f	148/150 (99%)	146 (99%)	2 (1%)	59	33
34	g	137/138 (99%)	131 (96%)	6 (4%)	25	4
35	h	113/114 (99%)	105 (93%)	8 (7%)	13	1
36	i	116/116 (100%)	115 (99%)	1 (1%)	70	51
37	j	104/104 (100%)	104 (100%)	0	100	100
38	k	103/103 (100%)	102 (99%)	1 (1%)	68	47
39	l	107/107 (100%)	106 (99%)	1 (1%)	70	51
40	m	98/103 (95%)	97 (99%)	1 (1%)	68	47
41	n	86/87 (99%)	86 (100%)	0	100	100
42	o	99/100 (99%)	99 (100%)	0	100	100
43	p	89/90 (99%)	89 (100%)	0	100	100
44	q	84/84 (100%)	84 (100%)	0	100	100
45	r	93/93 (100%)	92 (99%)	1 (1%)	65	43
46	s	80/84 (95%)	79 (99%)	1 (1%)	61	36
47	t	83/85 (98%)	80 (96%)	3 (4%)	31	6
48	u	78/78 (100%)	78 (100%)	0	100	100
49	v	58/63 (92%)	58 (100%)	0	100	100
50	w	67/68 (98%)	66 (98%)	1 (2%)	57	31
51	x	54/55 (98%)	54 (100%)	0	100	100
52	y	48/49 (98%)	48 (100%)	0	100	100
53	z	47/48 (98%)	46 (98%)	1 (2%)	47	19
All	All	4589/4763 (96%)	4521 (98%)	68 (2%)	55	31

All (68) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
6	B	4	VAL
6	B	8	ASP
6	B	10	LEU
6	B	80	VAL
6	B	127	ASP
6	B	146	ASN
7	C	44	THR

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Mol	Chain	Res	Type
7	C	53	SER
7	C	121	THR
7	C	192	THR
8	D	116	GLN
8	D	166	GLU
8	D	169	THR
8	D	172	GLU
9	E	55	GLU
10	F	103	VAL
11	G	6	VAL
12	H	54	ASP
14	J	40	ILE
14	J	75	ASP
14	J	100	ILE
15	K	13	ARG
15	K	14	LYS
15	K	108	THR
16	L	15	LYS
17	M	83	LEU
17	M	89	LEU
18	N	49	GLN
19	O	80	GLN
21	Q	42	THR
22	R	42	SER
22	R	66	SER
23	S	60	VAL
23	S	63	THR
24	T	23	SER
25	U	7	ARG
25	U	14	VAL
25	U	60	LEU
32	e	16	GLU
32	e	55	SER
32	e	107	SER
33	f	23	ASN
33	f	108	VAL
34	g	17	VAL
34	g	49	THR
34	g	51	THR
34	g	67	THR
34	g	81	GLU
34	g	92	VAL

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Mol	Chain	Res	Type
35	h	6	LEU
35	h	40	THR
35	h	43	ASN
35	h	55	GLU
35	h	108	VAL
35	h	116	ARG
35	h	134	VAL
35	h	142	VAL
36	i	11	VAL
38	k	109	LYS
39	l	58	LYS
40	m	6	SER
45	r	109	ASP
46	s	27	SER
47	t	15	THR
47	t	27	ASN
47	t	52	LEU
50	w	42	SER
53	z	55	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (30) such sidechains are listed below:

Mol	Chain	Res	Type
6	B	227	GLN
9	E	77	ASN
10	F	17	GLN
11	G	153	HIS
12	H	38	ASN
13	I	110	GLN
14	J	58	ASN
18	N	43	ASN
18	N	60	GLN
21	Q	50	ASN
24	T	3	ASN
24	T	61	GLN
25	U	9	ASN
30	c	86	ASN
30	c	134	ASN
32	e	9	GLN
33	f	5	HIS
34	g	38	ASN
34	g	104	ASN

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Mol	Chain	Res	Type
35	h	66	ASN
36	i	80	HIS
38	k	104	GLN
40	m	73	ASN
41	n	98	GLN
41	n	116	GLN
45	r	9	HIS
47	t	46	GLN
48	u	5	ASN
49	v	76	ASN
51	x	20	ASN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
26	X	5/6 (83%)	0	0
27	Z	73/77 (94%)	11 (15%)	1 (1%)
28	a	2742/2925 (93%)	260 (9%)	0
29	b	118/119 (99%)	10 (8%)	0
5	A	1495/1554 (96%)	172 (11%)	6 (0%)
All	All	4433/4681 (94%)	453 (10%)	7 (0%)

All (453) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
5	A	4	U
5	A	5	U
5	A	9	G
5	A	22	G
5	A	32	A
5	A	39	G
5	A	44	A
5	A	47	C
5	A	48	C
5	A	50	A
5	A	51	A
5	A	72	A
5	A	73	C
5	A	74	A
5	A	77	A
5	A	78	A

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Mol	Chain	Res	Type
5	A	96	U
5	A	120	A
5	A	121	U
5	A	131	A
5	A	143	A
5	A	144	G
5	A	182	A
5	A	183	C
5	A	197	A
5	A	204	G
5	A	240	G
5	A	245	U
5	A	247	G
5	A	251	G
5	A	266	G
5	A	267	C
5	A	289	G
5	A	321	A
5	A	328	C
5	A	329	A
5	A	339	C
5	A	347	G
5	A	352	C
5	A	354	G
5	A	367	U
5	A	372	C
5	A	373	A
5	A	384	G
5	A	397	A
5	A	406	G
5	A	412	A
5	A	413	G
5	A	414	A
5	A	415	A
5	A	421	U
5	A	422	C
5	A	424	G
5	A	429	U
5	A	430	A
5	A	436	C
5	A	446	G
5	A	453	G

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Mol	Chain	Res	Type
5	A	457	G
5	A	458	U
5	A	464	U
5	A	467	U
5	A	468	A
5	A	469	C
5	A	478	A
5	A	481	G
5	A	484	G
5	A	486	U
5	A	497	G
5	A	511	C
5	A	518	C
5	A	521	G
5	A	527	G7M
5	A	531	U
5	A	532	A
5	A	547	A
5	A	572	A
5	A	573	A
5	A	576	C
5	A	577	G
5	A	596	A
5	A	628	G
5	A	633	G
5	A	650	G
5	A	653	U
5	A	661	G
5	A	665	A
5	A	723	U
5	A	724	G
5	A	748	G
5	A	755	G
5	A	777	A
5	A	793	U
5	A	794	A
5	A	815	A
5	A	817	C
5	A	821	G
5	A	890	G
5	A	914	A
5	A	926	G

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Mol	Chain	Res	Type
5	A	934	C
5	A	935	A
5	A	960	U
5	A	966	2MG
5	A	969	A
5	A	971	G
5	A	975	A
5	A	976	G
5	A	977	A
5	A	993	G
5	A	996	A
5	A	1003	G
5	A	1004	A
5	A	1008	U
5	A	1020	G
5	A	1035	A
5	A	1036	A
5	A	1042	A
5	A	1044	A
5	A	1045	C
5	A	1053	G
5	A	1065	U
5	A	1085	U
5	A	1094	G
5	A	1095	U
5	A	1101	A
5	A	1132	C
5	A	1134	G
5	A	1137	C
5	A	1139	G
5	A	1140	C
5	A	1159	U
5	A	1167	A
5	A	1171	A
5	A	1184	G
5	A	1196	A
5	A	1197	A
5	A	1213	A
5	A	1227	A
5	A	1238	A
5	A	1253	G
5	A	1258	G

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Mol	Chain	Res	Type
5	A	1260	G
5	A	1279	G
5	A	1280	A
5	A	1285	A
5	A	1286	U
5	A	1287	A
5	A	1300	G
5	A	1302	C
5	A	1317	C
5	A	1320	C
5	A	1338	G
5	A	1346	A
5	A	1353	G
5	A	1363	A
5	A	1370	G
5	A	1378	C
5	A	1379	G
5	A	1398	A
5	A	1419	G
5	A	1441	A
5	A	1451	U
5	A	1452	C
5	A	1453	G
5	A	1493	A
5	A	1497	G
5	A	1503	A
5	A	1506	U
5	A	1517	G
5	A	1529	G
5	A	1530	G
27	Z	6	G
27	Z	9	G
27	Z	17	C
27	Z	18	G
27	Z	20	U
27	Z	21	A
27	Z	22	G
27	Z	46	G
27	Z	58	A
27	Z	64	G
27	Z	76	A
28	a	10	A

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Mol	Chain	Res	Type
28	a	15	G
28	a	34	U
28	a	42	A
28	a	71	A
28	a	74	A
28	a	75	G
28	a	101	A
28	a	102	U
28	a	118	A
28	a	119	A
28	a	120	U
28	a	125	A
28	a	136	G
28	a	139	U
28	a	140	C
28	a	142	A
28	a	163	C
28	a	165	A
28	a	181	A
28	a	196	A
28	a	215	G
28	a	216	A
28	a	221	A
28	a	222	A
28	a	248	G
28	a	272	A
28	a	276	U
28	a	277	G
28	a	278	A
28	a	279	A
28	a	281	C
28	a	287	G
28	a	311	A
28	a	329	G
28	a	330	A
28	a	345	A
28	a	361	G
28	a	367	G
28	a	386	G
28	a	405	U
28	a	411	G
28	a	412	A

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Mol	Chain	Res	Type
28	a	481	G
28	a	491	G
28	a	505	A
28	a	509	C
28	a	530	G
28	a	531	C
28	a	532	A
28	a	565	A
28	a	575	U
28	a	577	A
28	a	605	A
28	a	617	U
28	a	629	A
28	a	639	A
28	a	647	C
28	a	648	U
28	a	649	G
28	a	655	U
28	a	656	A
28	a	657	A
28	a	688	U
28	a	719	C
28	a	732	A
28	a	749	5MU
28	a	766	A
28	a	767	C
28	a	777	G
28	a	778	G
28	a	784	A
28	a	786	G
28	a	787	G
28	a	807	G
28	a	814	C
28	a	829	U
28	a	830	U
28	a	848	U
28	a	861	G
28	a	884	G
28	a	886	U
28	a	887	C
28	a	890	C
28	a	892	C

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Mol	Chain	Res	Type
28	a	893	G
28	a	897	U
28	a	898	A
28	a	899	C
28	a	901	A
28	a	912	A
28	a	916	G
28	a	933	U
28	a	948	C
28	a	963	C
28	a	976	G
28	a	985	A
28	a	986	A
28	a	987	C
28	a	998	A
28	a	1014	U
28	a	1015	C
28	a	1024	G
28	a	1035	U
28	a	1042	A
28	a	1049	G
28	a	1110	U
28	a	1112	G
28	a	1113	A
28	a	1114	G
28	a	1118	G
28	a	1124	G
28	a	1134	U
28	a	1137	C
28	a	1144	A
28	a	1173	G
28	a	1255	A
28	a	1258	G
28	a	1273	G
28	a	1274	A
28	a	1302	G
28	a	1303	A
28	a	1354	U
28	a	1367	A
28	a	1381	U
28	a	1385	A
28	a	1413	U

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Mol	Chain	Res	Type
28	a	1418	G
28	a	1430	C
28	a	1436	A
28	a	1454	G
28	a	1455	A
28	a	1484	G
28	a	1495	C
28	a	1510	A
28	a	1511	A
28	a	1512	G
28	a	1517	A
28	a	1531	G
28	a	1536	U
28	a	1537	A
28	a	1538	C
28	a	1539	G
28	a	1568	A
28	a	1571	A
28	a	1580	U
28	a	1585	A
28	a	1587	C
28	a	1610	A
28	a	1649	U
28	a	1650	U
28	a	1651	G
28	a	1676	G
28	a	1717	G
28	a	1731	U
28	a	1732	C
28	a	1740	G
28	a	1746	A
28	a	1766	C
28	a	1775	A
28	a	1802	C
28	a	1803	A
28	a	1810	A
28	a	1818	C
28	a	1831	A
28	a	1849	A
28	a	1850	A
28	a	1860	A
28	a	1872	G

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Mol	Chain	Res	Type
28	a	1910	G
28	a	1911	G
28	a	1917	A
28	a	1918	C
28	a	1933	G
28	a	1934	G
28	a	1941	A
28	a	1942	A
28	a	1959	U
28	a	1971	C
28	a	1974	A
28	a	1975	U
28	a	1976	G
28	a	1995	U
28	a	1997	U
28	a	2027	C
28	a	2035	A
28	a	2037	A
28	a	2047	C
28	a	2059	C
28	a	2060	G
28	a	2064	A
28	a	2065	G
28	a	2066	A
28	a	2073	G7M
28	a	2202	A
28	a	2208	G
28	a	2215	G
28	a	2227	G
28	a	2229	A
28	a	2242	G
28	a	2243	G
28	a	2272	A
28	a	2287	C
28	a	2291	A
28	a	2309	U
28	a	2312	G
28	a	2326	A
28	a	2329	G
28	a	2337	A
28	a	2339	A
28	a	2351	C

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Mol	Chain	Res	Type
28	a	2376	U
28	a	2387	G
28	a	2389	C
28	a	2400	G
28	a	2410	A
28	a	2429	A
28	a	2433	G
28	a	2434	A
28	a	2435	U
28	a	2439	A
28	a	2445	U
28	a	2452	A
28	a	2474	G
28	a	2495	U
28	a	2506	G
28	a	2509	G
28	a	2522	A
28	a	2524	C
28	a	2529	G
28	a	2533	G
28	a	2551	A
28	a	2570	A
28	a	2571	G
28	a	2577	C
28	a	2606	A
28	a	2613	U
28	a	2617	U
28	a	2633	U
28	a	2634	G
28	a	2667	G
28	a	2693	U
28	a	2694	U
28	a	2718	G
28	a	2730	A
28	a	2748	G
28	a	2752	A
28	a	2761	A
28	a	2769	A
28	a	2782	A
28	a	2802	U
28	a	2803	G
28	a	2824	A

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Mol	Chain	Res	Type
28	a	2825	A
28	a	2839	A
28	a	2865	U
28	a	2877	A
28	a	2878	C
28	a	2887	A
28	a	2888	U
29	b	35	C
29	b	36	C
29	b	45	A
29	b	56	G
29	b	57	A
29	b	67	G
29	b	89	U
29	b	90	C
29	b	99	A
29	b	109	A

All (7) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
5	A	182	A
5	A	429	U
5	A	467	U
5	A	766	A
5	A	786	G
5	A	1035	A
27	Z	17(A)	U

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

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

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
28	2MG	a	1837	28	23,26,27	0.69	0	32,38,41	0.56	0
5	UR3	A	1498	5	19,22,23	2.56	6 (31%)	26,32,35	1.37	3 (11%)
28	PSU	a	1921	28	18,21,22	1.03	1 (5%)	22,30,33	1.83	4 (18%)
39	4D4	l	81	39	9,11,12	0.54	0	8,13,15	1.03	1 (12%)
28	OMU	a	2556	56,55,28	19,22,23	3.07	7 (36%)	26,31,34	1.67	4 (15%)
15	IAS	K	119	15	6,7,8	0.98	0	6,8,10	1.22	1 (16%)
5	MA6	A	1519	5	23,26,27	1.38	5 (21%)	34,38,41	2.88	13 (38%)
28	PSU	a	748	56,28	18,21,22	1.11	2 (11%)	22,30,33	1.61	3 (13%)
28	PSU	a	2584	28	18,21,22	1.10	2 (11%)	22,30,33	1.76	3 (13%)
28	5MU	a	749	28	19,22,23	0.71	0	28,32,35	0.55	0
39	MS6	l	82	39	5,7,8	0.40	0	2,7,9	0.18	0
5	5MC	A	967	5	18,22,23	0.73	0	26,32,35	0.61	0
5	G7M	A	527	5,55	23,26,27	2.73	8 (34%)	35,39,42	1.70	9 (25%)
28	1MG	a	747	28	22,26,27	2.59	6 (27%)	33,39,42	2.07	7 (21%)
5	2MG	A	966	5	23,26,27	0.56	0	32,38,41	0.56	0
28	H2U	a	2453	28	18,21,22	0.65	0	21,30,33	1.05	3 (14%)
28	PSU	a	2608	28	18,21,22	1.06	1 (5%)	22,30,33	1.86	3 (13%)
28	2MA	a	2507	56,55,28	22,25,26	3.73	11 (50%)	33,37,40	2.78	11 (33%)
28	PSU	a	2508	55,28	18,21,22	1.03	1 (5%)	22,30,33	1.87	4 (18%)
5	2MG	A	1516	5	23,26,27	0.65	0	32,38,41	0.63	0
5	MA6	A	1518	5	23,26,27	1.37	5 (21%)	34,38,41	2.80	14 (41%)
28	2MG	a	2449	28	23,26,27	0.78	0	32,38,41	0.59	0
16	D2T	L	89	16	7,9,10	1.34	1 (14%)	6,11,13	1.79	2 (33%)
28	3TD	a	1919	28	18,22,23	4.03	7 (38%)	22,32,35	1.79	3 (13%)
28	OMC	a	2502	56,28	19,22,23	0.81	1 (5%)	26,31,34	0.66	0
28	PSU	a	1915	28	18,21,22	1.02	1 (5%)	22,30,33	1.77	3 (13%)
5	PSU	A	516	5,56	18,21,22	1.07	1 (5%)	22,30,33	1.91	6 (27%)
28	5MU	a	1943	55,28	19,22,23	0.80	0	28,32,35	0.46	0
5	4OC	A	1402	5	20,23,24	2.99	8 (40%)	26,32,35	0.94	1 (3%)
28	G7M	a	2073	55,28	23,26,27	2.66	9 (39%)	35,39,42	1.69	9 (25%)
28	6MZ	a	1620	28	22,25,26	2.85	7 (31%)	30,36,39	2.72	10 (33%)
28	PSU	a	957	28	18,21,22	1.09	1 (5%)	22,30,33	1.76	2 (9%)
5	2MG	A	1207	5,55	23,26,27	0.55	0	32,38,41	0.54	0
28	OMG	a	2255	27,55,28	23,26,27	0.72	0	33,38,41	0.54	0
28	PSU	a	2609	28	18,21,22	1.05	1 (5%)	22,30,33	1.91	4 (18%)
31	MEQ	d	150	31	8,9,10	0.81	0	5,10,12	0.78	0
28	5MC	a	1966	55,28	18,22,23	0.82	0	26,32,35	0.52	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	5MC	A	1407	5	18,22,23	0.82	1 (5%)	26,32,35	0.57	0
28	6MZ	a	2034	28	22,25,26	2.79	6 (27%)	30,36,39	2.86	11 (36%)
28	PSU	a	2461	28	18,21,22	1.12	1 (5%)	22,30,33	1.82	4 (18%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
28	2MG	a	1837	28	-	0/9/27/28	0/3/3/3
5	UR3	A	1498	5	-	0/7/25/26	0/2/2/2
28	PSU	a	1921	28	-	0/7/25/26	0/2/2/2
39	4D4	l	81	39	-	6/11/12/14	-
28	OMU	a	2556	56,55,28	-	0/9/27/28	0/2/2/2
15	IAS	K	119	15	-	0/7/7/8	-
5	MA6	A	1519	5	-	3/11/29/30	0/3/3/3
28	PSU	a	748	56,28	-	1/7/25/26	0/2/2/2
28	PSU	a	2584	28	-	0/7/25/26	0/2/2/2
28	5MU	a	749	28	-	0/7/25/26	0/2/2/2
39	MS6	l	82	39	-	1/4/6/8	-
5	5MC	A	967	5	-	0/7/25/26	0/2/2/2
5	G7M	A	527	5,55	-	3/7/25/26	0/3/3/3
28	1MG	a	747	28	-	0/7/25/26	0/3/3/3
5	2MG	A	966	5	-	0/9/27/28	0/3/3/3
28	H2U	a	2453	28	-	0/7/38/39	0/2/2/2
28	PSU	a	2608	28	-	0/7/25/26	0/2/2/2
28	2MA	a	2507	56,55,28	-	2/7/25/26	0/3/3/3
28	PSU	a	2508	55,28	-	0/7/25/26	0/2/2/2
5	2MG	A	1516	5	-	0/9/27/28	0/3/3/3
5	MA6	A	1518	5	-	0/11/29/30	0/3/3/3
28	2MG	a	2449	28	-	0/9/27/28	0/3/3/3
16	D2T	L	89	16	-	3/7/12/14	-
28	3TD	a	1919	28	-	2/7/25/26	0/2/2/2
28	OMC	a	2502	56,28	-	0/9/27/28	0/2/2/2
28	PSU	a	1915	28	-	0/7/25/26	0/2/2/2
5	PSU	A	516	5,56	-	0/7/25/26	0/2/2/2
28	5MU	a	1943	55,28	-	0/7/25/26	0/2/2/2
5	4OC	A	1402	5	-	0/9/29/30	0/2/2/2
28	G7M	a	2073	55,28	-	2/7/25/26	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
28	6MZ	a	1620	28	-	0/9/27/28	0/3/3/3
28	PSU	a	957	28	-	0/7/25/26	0/2/2/2
5	2MG	A	1207	5,55	-	0/9/27/28	0/3/3/3
28	OMG	a	2255	27,55,28	-	1/9/27/28	0/3/3/3
28	PSU	a	2609	28	-	0/7/25/26	0/2/2/2
31	MEQ	d	150	31	-	2/8/9/11	-
28	5MC	a	1966	55,28	-	0/7/25/26	0/2/2/2
5	5MC	A	1407	5	-	0/7/25/26	0/2/2/2
28	6MZ	a	2034	28	-	2/9/27/28	0/3/3/3
28	PSU	a	2461	28	-	0/7/25/26	0/2/2/2

All (100) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
28	a	1919	3TD	C6-C5	12.20	1.49	1.35
28	a	1620	6MZ	C6-N6	11.63	1.46	1.34
28	a	2034	6MZ	C6-N6	11.17	1.46	1.34
28	a	2507	2MA	C4-N3	10.38	1.47	1.34
28	a	1919	3TD	C2-N1	8.60	1.48	1.37
28	a	2507	2MA	C6-N1	6.44	1.44	1.35
28	a	747	1MG	C2-N3	6.33	1.45	1.34
5	A	1402	4OC	C4-N3	6.31	1.43	1.32
5	A	527	G7M	C2-N2	6.31	1.49	1.34
28	a	2507	2MA	C2-N3	6.28	1.45	1.34
5	A	1498	UR3	C2-N1	6.27	1.47	1.38
28	a	2556	OMU	C2-N3	6.25	1.49	1.38
5	A	527	G7M	C4-N3	6.13	1.48	1.34
5	A	1402	4OC	C6-C5	6.05	1.49	1.35
28	a	2073	G7M	C2-N2	5.99	1.48	1.34
28	a	747	1MG	C4-N3	5.98	1.48	1.34
28	a	2556	OMU	C2-N1	5.93	1.48	1.38
28	a	2507	2MA	C2-N1	5.83	1.44	1.34
28	a	747	1MG	C2-N2	5.80	1.44	1.34
5	A	1498	UR3	C6-C5	5.75	1.48	1.35
5	A	1402	4OC	C2-N3	5.66	1.47	1.36
28	a	2073	G7M	C4-N3	5.56	1.47	1.34
28	a	1919	3TD	C6-N1	5.36	1.45	1.36
5	A	527	G7M	C2-N3	5.35	1.46	1.33
28	a	2073	G7M	C5-N7	-5.32	1.33	1.39
28	a	2556	OMU	C6-C5	5.26	1.47	1.35
28	a	2556	OMU	O4-C4	-5.14	1.14	1.24
28	a	2073	G7M	C2-N3	4.80	1.44	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	1402	4OC	C4-N4	4.69	1.45	1.35
28	a	2507	2MA	C5-C6	4.67	1.54	1.41
5	A	1498	UR3	C2-N3	4.56	1.47	1.39
5	A	527	G7M	C5-N7	-4.49	1.34	1.39
28	a	2556	OMU	O2-C2	-4.46	1.14	1.23
28	a	1919	3TD	C2-N3	3.98	1.47	1.38
28	a	2507	2MA	C4-N9	-3.98	1.28	1.37
28	a	2507	2MA	C5-N7	-3.88	1.31	1.39
28	a	2556	OMU	C4-N3	3.71	1.45	1.38
5	A	1402	4OC	C5-C4	3.71	1.48	1.40
5	A	527	G7M	C5-C6	3.70	1.53	1.43
28	a	747	1MG	C5-N7	-3.69	1.31	1.39
28	a	2034	6MZ	C5-N7	-3.60	1.32	1.39
28	a	2507	2MA	C8-N7	3.56	1.38	1.31
5	A	1402	4OC	C2-N1	3.54	1.47	1.40
5	A	1402	4OC	O2-C2	-3.43	1.17	1.23
5	A	1518	MA6	C5-C4	-3.42	1.32	1.39
28	a	2073	G7M	C5-C6	3.37	1.52	1.43
28	a	2073	G7M	O6-C6	-3.36	1.17	1.23
28	a	1620	6MZ	C5-N7	-3.35	1.32	1.39
28	a	2556	OMU	C6-N1	3.29	1.45	1.38
5	A	1519	MA6	C5-C4	-3.26	1.32	1.39
5	A	516	PSU	C6-C5	3.11	1.39	1.35
28	a	748	PSU	C6-C5	3.07	1.38	1.35
28	a	2461	PSU	C6-C5	3.00	1.38	1.35
28	a	1921	PSU	C6-C5	2.99	1.38	1.35
28	a	957	PSU	C6-C5	2.92	1.38	1.35
28	a	1915	PSU	C6-C5	2.90	1.38	1.35
28	a	2508	PSU	C6-C5	2.90	1.38	1.35
28	a	747	1MG	C5-C6	2.88	1.52	1.45
5	A	1402	4OC	C6-N1	2.83	1.44	1.38
5	A	527	G7M	O6-C6	-2.81	1.18	1.23
5	A	1498	UR3	O2-C2	-2.76	1.17	1.22
5	A	527	G7M	C2-N1	2.75	1.44	1.37
5	A	1498	UR3	O4-C4	-2.71	1.17	1.23
28	a	2034	6MZ	C5-C4	-2.63	1.34	1.39
5	A	1498	UR3	C6-N1	2.62	1.44	1.38
28	a	2584	PSU	C6-C5	2.62	1.38	1.35
5	A	1518	MA6	C6-N6	2.60	1.44	1.36
28	a	1620	6MZ	C5-C4	-2.60	1.34	1.39
5	A	1519	MA6	C6-N6	2.58	1.44	1.36
28	a	2034	6MZ	C8-N9	-2.58	1.32	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	527	G7M	C6-N1	2.58	1.43	1.38
28	a	747	1MG	O6-C6	-2.57	1.17	1.23
28	a	2608	PSU	C6-C5	2.56	1.38	1.35
28	a	2609	PSU	C6-C5	2.55	1.38	1.35
28	a	2507	2MA	C6-N6	-2.49	1.27	1.34
5	A	1519	MA6	C5-N7	-2.49	1.34	1.39
28	a	1919	3TD	O4-C4	-2.48	1.17	1.23
5	A	1518	MA6	C8-N9	-2.40	1.33	1.37
28	a	2034	6MZ	C6-N1	-2.37	1.31	1.35
28	a	2073	G7M	C2-N1	2.35	1.43	1.37
28	a	1919	3TD	C1'-C5	2.34	1.55	1.50
28	a	1620	6MZ	C8-N9	-2.34	1.33	1.37
28	a	1620	6MZ	C6-N1	-2.32	1.31	1.35
28	a	2584	PSU	O4'-C1'	-2.32	1.40	1.43
5	A	1518	MA6	C5-N7	-2.29	1.34	1.39
28	a	2034	6MZ	C2-N1	2.28	1.38	1.33
5	A	1519	MA6	C8-N9	-2.28	1.33	1.37
5	A	1519	MA6	C4-N9	-2.22	1.32	1.37
28	a	1919	3TD	C4-N3	2.21	1.45	1.40
28	a	2507	2MA	C8-N9	-2.18	1.33	1.37
5	A	1518	MA6	C4-N9	-2.16	1.32	1.37
28	a	2507	2MA	C5-C4	-2.13	1.35	1.39
28	a	2502	OMC	C4-N3	-2.10	1.30	1.34
28	a	2073	G7M	C4-N9	-2.09	1.32	1.38
28	a	2073	G7M	C6-N1	2.09	1.42	1.38
28	a	1620	6MZ	C2-N1	2.03	1.37	1.33
5	A	1407	5MC	C4-N3	-2.02	1.30	1.34
28	a	1620	6MZ	C5-C6	-2.01	1.36	1.41
28	a	748	PSU	O4'-C1'	-2.00	1.41	1.43
16	L	89	D2T	CB1-SB	-2.00	1.75	1.79

All (138) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	1519	MA6	N1-C6-N6	-9.39	106.81	117.08
5	A	1518	MA6	N1-C6-N6	-8.86	107.40	117.08
28	a	2507	2MA	C1'-N9-C8	-8.83	107.21	127.14
28	a	2034	6MZ	C4-C5-C6	7.13	122.34	116.81
28	a	2507	2MA	C4-N9-C1'	6.92	143.07	126.59
28	a	1620	6MZ	C4-C5-C6	6.64	121.95	116.81
28	a	2034	6MZ	C5-C4-N3	-6.28	118.56	126.75
28	a	1620	6MZ	C5-C4-N3	-6.17	118.69	126.75

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
28	a	2507	2MA	N9-C8-N7	-5.95	105.77	113.91
5	A	1519	MA6	N1-C2-N3	-5.93	119.32	128.60
28	a	1919	3TD	N1-C2-N3	5.89	120.79	116.14
5	A	1518	MA6	N1-C2-N3	-5.85	119.45	128.60
28	a	747	1MG	C5-C4-N3	-5.60	119.37	128.46
28	a	747	1MG	C1'-N9-C4	-5.45	110.30	126.50
28	a	1620	6MZ	N1-C2-N3	-5.25	120.39	128.60
28	a	747	1MG	C1'-N9-C8	5.22	141.55	126.70
5	A	1518	MA6	N9-C8-N7	-5.13	106.89	113.91
28	a	2556	OMU	C4-N3-C2	-5.08	119.88	126.58
28	a	2609	PSU	N1-C2-N3	5.07	120.87	115.13
28	a	1620	6MZ	C9-N6-C6	5.07	127.23	122.87
28	a	2034	6MZ	C9-N6-C6	5.04	127.21	122.87
28	a	2461	PSU	N1-C2-N3	5.03	120.83	115.13
28	a	2034	6MZ	N3-C4-N9	4.99	135.30	127.08
28	a	2508	PSU	N1-C2-N3	4.95	120.73	115.13
5	A	1519	MA6	N9-C8-N7	-4.94	107.15	113.91
5	A	1519	MA6	C5-C6-N6	4.92	133.87	125.30
28	a	2034	6MZ	N1-C2-N3	-4.91	120.92	128.60
5	A	516	PSU	C4-N3-C2	-4.90	119.28	126.34
28	a	957	PSU	C4-N3-C2	-4.86	119.34	126.34
28	a	2608	PSU	C4-N3-C2	-4.85	119.35	126.34
28	a	1921	PSU	C4-N3-C2	-4.82	119.39	126.34
5	A	1498	UR3	C4-N3-C2	-4.80	120.05	124.56
28	a	1915	PSU	C4-N3-C2	-4.80	119.43	126.34
28	a	2609	PSU	C4-N3-C2	-4.79	119.43	126.34
5	A	516	PSU	N1-C2-N3	4.79	120.55	115.13
28	a	2608	PSU	N1-C2-N3	4.74	120.50	115.13
28	a	2508	PSU	C4-N3-C2	-4.67	119.61	126.34
28	a	1921	PSU	N1-C2-N3	4.67	120.42	115.13
28	a	2584	PSU	C4-N3-C2	-4.67	119.61	126.34
28	a	1620	6MZ	N3-C4-N9	4.62	134.70	127.08
28	a	957	PSU	N1-C2-N3	4.59	120.33	115.13
5	A	1518	MA6	C5-C6-N6	4.49	133.13	125.30
28	a	748	PSU	C4-N3-C2	-4.47	119.90	126.34
28	a	2461	PSU	C4-N3-C2	-4.46	119.92	126.34
28	a	2073	G7M	C2-N3-C4	4.44	120.21	112.30
28	a	2034	6MZ	N9-C8-N7	-4.44	107.84	113.91
28	a	2584	PSU	N1-C2-N3	4.41	120.13	115.13
28	a	1915	PSU	N1-C2-N3	4.41	120.12	115.13
5	A	527	G7M	C2-N3-C4	4.21	119.80	112.30
28	a	2556	OMU	N3-C2-N1	4.19	120.45	114.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
28	a	748	PSU	N1-C2-N3	4.18	119.86	115.13
5	A	1519	MA6	C5-C4-N3	-4.15	121.34	126.75
5	A	1519	MA6	C2-N1-C6	4.13	121.52	111.75
5	A	1518	MA6	C2-N1-C6	4.05	121.33	111.75
28	a	1919	3TD	C4-N3-C2	-3.97	120.30	124.61
28	a	747	1MG	N9-C4-N3	3.96	133.89	125.94
5	A	527	G7M	C5-C6-N1	3.92	119.97	111.79
5	A	1518	MA6	C5-C4-N3	-3.91	121.65	126.75
28	a	2073	G7M	C5-C6-N1	3.83	119.78	111.79
28	a	2507	2MA	CM2-C2-N1	3.80	123.08	117.15
28	a	2073	G7M	C5-C4-N3	-3.73	121.00	128.15
28	a	747	1MG	C2-N3-C4	3.69	120.27	111.98
28	a	1620	6MZ	N9-C8-N7	-3.67	108.90	113.91
5	A	1518	MA6	C1'-N9-C8	-3.65	118.89	127.14
28	a	2507	2MA	C4-N9-C8	3.64	109.67	105.73
5	A	527	G7M	C5-C4-N3	-3.58	121.28	128.15
28	a	2507	2MA	N3-C2-N1	-3.58	119.13	125.72
28	a	1620	6MZ	C2-N3-C4	3.51	120.04	111.75
28	a	2034	6MZ	C6-C5-N7	-3.48	128.63	132.39
5	A	1518	MA6	C5-N7-C8	3.40	108.34	103.51
28	a	2034	6MZ	C2-N3-C4	3.35	119.66	111.75
5	A	1518	MA6	C4-N9-C8	3.35	109.36	105.73
5	A	1519	MA6	C2-N3-C4	3.32	119.60	111.75
28	a	1620	6MZ	C6-C5-N7	-3.31	128.81	132.39
5	A	1519	MA6	C5-N7-C8	3.30	108.20	103.51
28	a	2034	6MZ	C5-N7-C8	3.28	108.17	103.51
5	A	527	G7M	O6-C6-C5	-3.26	120.69	128.06
5	A	1519	MA6	C1'-N9-C8	-3.23	119.85	127.14
28	a	2508	PSU	O2-C2-N1	-3.19	119.27	122.79
5	A	1518	MA6	C2-N3-C4	3.18	119.27	111.75
28	a	2073	G7M	O6-C6-C5	-3.13	121.01	128.06
28	a	2507	2MA	C5-N7-C8	3.12	107.94	103.51
5	A	1519	MA6	C4-N9-C8	3.02	109.00	105.73
28	a	2556	OMU	C5-C4-N3	3.01	119.35	114.84
28	a	2507	2MA	C5-C4-N9	3.00	109.27	105.78
5	A	527	G7M	C2-N1-C6	-2.88	119.85	125.10
28	a	2461	PSU	C6-N1-C2	-2.85	119.77	122.68
28	a	2034	6MZ	C4-N9-C8	2.85	108.82	105.73
28	a	2073	G7M	C2-N1-C6	-2.84	119.92	125.10
5	A	527	G7M	CN7-N7-C5	2.75	130.18	126.77
5	A	1519	MA6	C4-C5-C6	2.69	118.89	115.88
28	a	2034	6MZ	C4-N9-C1'	-2.69	120.19	126.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
28	a	2507	2MA	C4-C5-N7	-2.69	107.35	110.62
5	A	1498	UR3	C6-N1-C2	-2.67	119.40	121.79
28	a	1620	6MZ	C5-N7-C8	2.63	107.25	103.51
5	A	1518	MA6	C4-C5-C6	2.62	118.82	115.88
28	a	2508	PSU	C6-N1-C2	-2.62	120.01	122.68
28	a	2609	PSU	C6-N1-C2	-2.60	120.02	122.68
5	A	1402	4OC	C6-C5-C4	2.57	120.11	116.96
5	A	527	G7M	N9-C4-N3	2.55	131.06	125.94
28	a	1921	PSU	O2-C2-N1	-2.53	120.01	122.79
5	A	516	PSU	O2-C2-N1	-2.52	120.02	122.79
5	A	516	PSU	C6-C5-C4	2.47	119.92	118.20
28	a	2073	G7M	N9-C4-N3	2.46	130.89	125.94
28	a	747	1MG	N9-C8-N7	-2.45	108.78	113.39
16	L	89	D2T	OD2-CG-CB	2.41	118.36	113.15
28	a	2556	OMU	O4-C4-C5	-2.40	120.94	125.16
28	a	1919	3TD	C6-C5-C4	2.38	119.86	118.22
28	a	1915	PSU	O2-C2-N1	-2.36	120.19	122.79
5	A	1518	MA6	C4-C5-N7	-2.33	107.78	110.62
28	a	2584	PSU	O4'-C1'-C2'	2.32	108.42	105.14
28	a	2507	2MA	N6-C6-N1	2.29	120.27	117.03
5	A	1519	MA6	C4-C5-N7	-2.29	107.83	110.62
5	A	1519	MA6	N3-C4-N9	2.28	130.84	127.08
28	a	747	1MG	C5-C6-N1	2.28	119.31	114.91
16	L	89	D2T	O-C-CA	-2.26	118.85	124.78
39	l	81	4D4	O-C-CA	-2.24	118.92	124.78
28	a	748	PSU	C6-N1-C2	-2.23	120.41	122.68
28	a	2453	H2U	C4-N3-C2	2.21	127.63	125.79
5	A	1518	MA6	N3-C4-N9	2.21	130.72	127.08
5	A	1498	UR3	C1'-N1-C2	2.20	120.70	116.99
28	a	2073	G7M	N2-C2-N1	2.20	121.39	116.71
28	a	2507	2MA	C6-C5-N7	2.20	136.11	132.02
28	a	2609	PSU	O2-C2-N1	-2.20	120.37	122.79
28	a	2461	PSU	O2-C2-N3	-2.18	117.71	121.82
28	a	2453	H2U	C5-C4-N3	-2.18	114.20	116.65
28	a	2608	PSU	C6-C5-C4	2.17	119.71	118.20
5	A	1518	MA6	C4-N9-C1'	2.16	131.73	126.59
15	K	119	IAS	OD1-CG-CB	-2.15	119.15	125.43
5	A	527	G7M	CN7-N7-C8	-2.15	121.52	124.84
28	a	2073	G7M	CN7-N7-C5	2.14	129.43	126.77
5	A	516	PSU	O4'-C1'-C2'	2.14	108.16	105.14
28	a	2453	H2U	O2-C2-N3	-2.11	117.57	121.50
28	a	1620	6MZ	C4-N9-C8	2.11	108.01	105.73

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
28	a	2073	G7M	N9-C8-N7	-2.10	107.01	112.21
5	A	516	PSU	C6-N1-C2	-2.09	120.55	122.68
5	A	527	G7M	N9-C8-N7	-2.06	107.12	112.21
28	a	1921	PSU	C6-N1-C2	-2.01	120.63	122.68

There are no chirality outliers.

All (28) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
39	l	81	4D4	NH1-CZ-NE-CD
28	a	2034	6MZ	O4'-C4'-C5'-O5'
28	a	2255	OMG	C1'-C2'-O2'-CM2
28	a	2034	6MZ	C3'-C4'-C5'-O5'
31	d	150	MEQ	OE1-CD-CG-CB
31	d	150	MEQ	NE2-CD-CG-CB
5	A	527	G7M	C3'-C4'-C5'-O5'
5	A	1519	MA6	O4'-C4'-C5'-O5'
5	A	1519	MA6	C3'-C4'-C5'-O5'
28	a	1919	3TD	O4'-C4'-C5'-O5'
5	A	1519	MA6	C5-C6-N6-C10
5	A	527	G7M	O4'-C4'-C5'-O5'
39	l	81	4D4	NH2-CZ-NE-CD
39	l	82	MS6	CB-CG-SD-CE
28	a	2073	G7M	C4'-C5'-O5'-P
5	A	527	G7M	C4'-C5'-O5'-P
39	l	81	4D4	OB-CB-CG-CD
39	l	81	4D4	CG-CD-NE-CZ
39	l	81	4D4	NE-CD-CG-CB
16	L	89	D2T	CA-CB-SB-CB1
28	a	2507	2MA	C4'-C5'-O5'-P
28	a	1919	3TD	C3'-C4'-C5'-O5'
16	L	89	D2T	SB-CB-CG-OD2
16	L	89	D2T	CG-CB-SB-CB1
28	a	748	PSU	O4'-C1'-C5-C6
28	a	2073	G7M	O4'-C4'-C5'-O5'
28	a	2507	2MA	O4'-C4'-C5'-O5'
39	l	81	4D4	O-C-CA-CB

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
28	a	1919	3TD	1	0
28	a	2034	6MZ	2	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 530 ligands modelled in this entry, 530 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

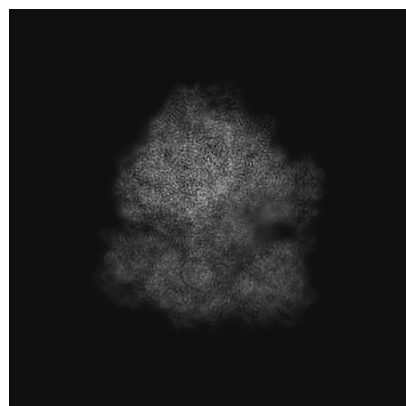
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-15793. These allow visual inspection of the internal detail of the map and identification of artifacts.

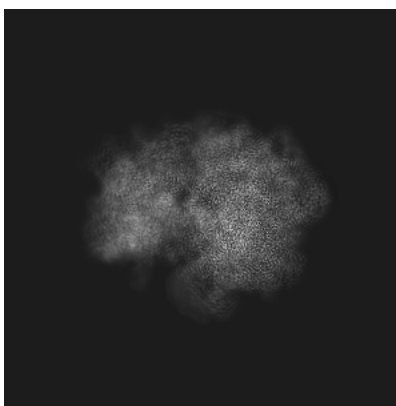
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

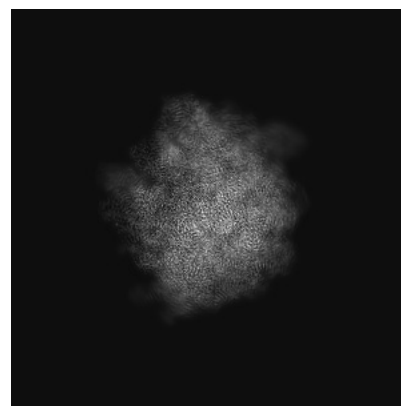
6.1.1 Primary map



X

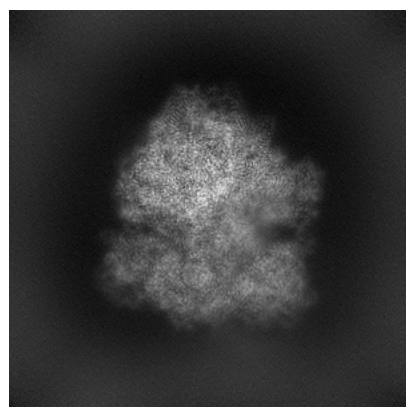


Y

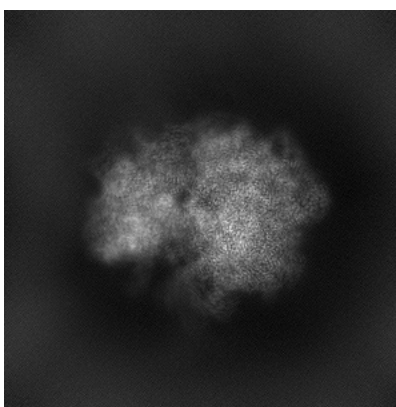


Z

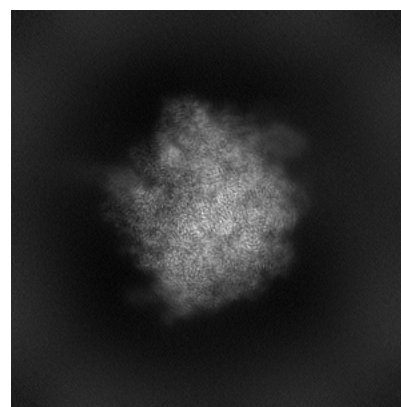
6.1.2 Raw map



X



Y

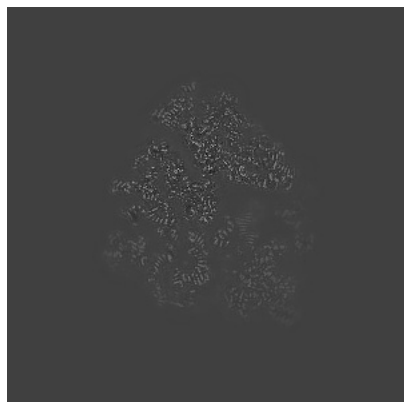


Z

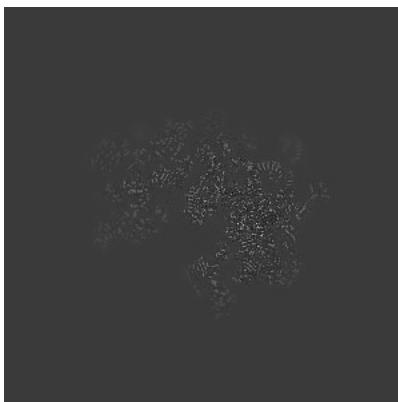
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

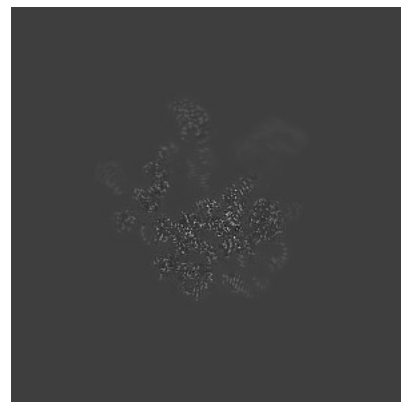
6.2.1 Primary map



X Index: 280

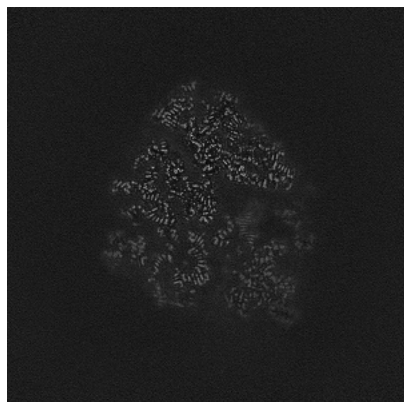


Y Index: 280

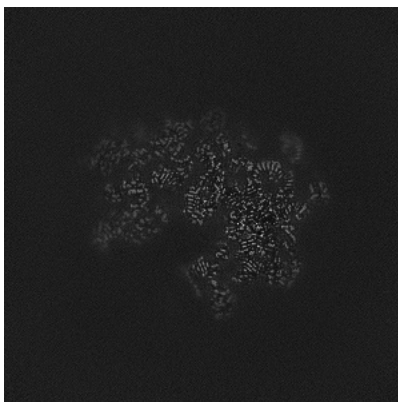


Z Index: 280

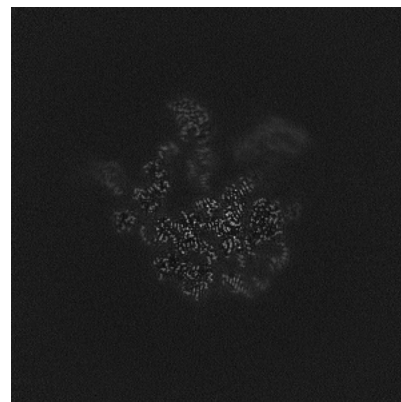
6.2.2 Raw map



X Index: 280



Y Index: 280

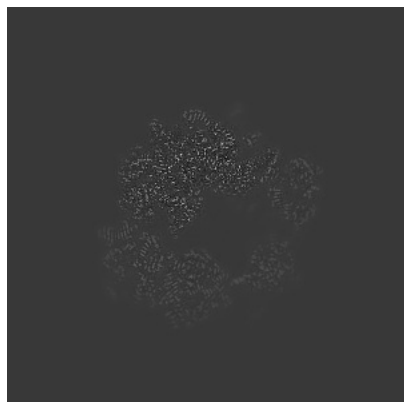


Z Index: 280

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

6.3 Largest variance slices [i](#)

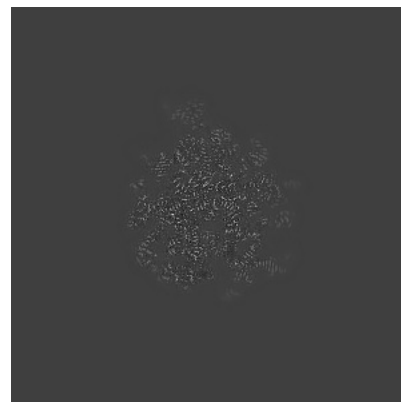
6.3.1 Primary map



X Index: 246

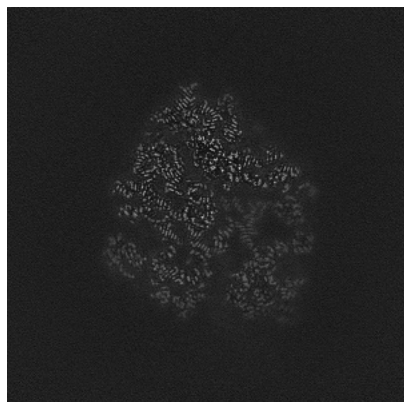


Y Index: 260

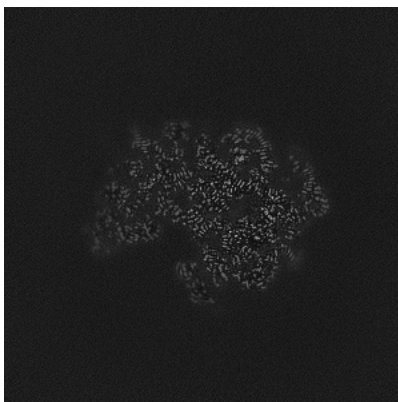


Z Index: 342

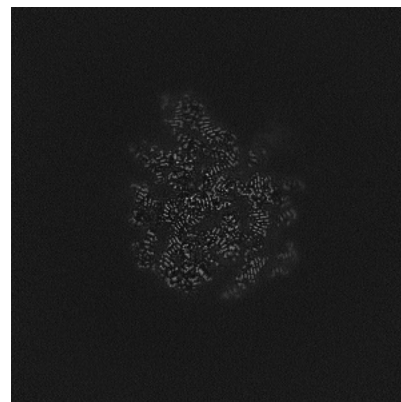
6.3.2 Raw map



X Index: 274



Y Index: 260

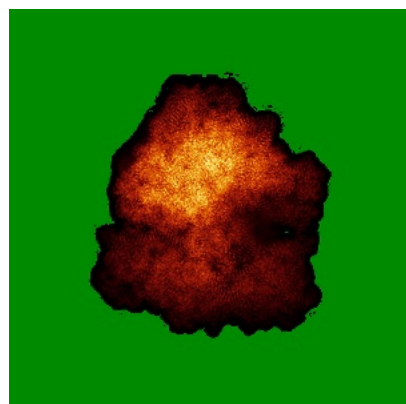


Z Index: 334

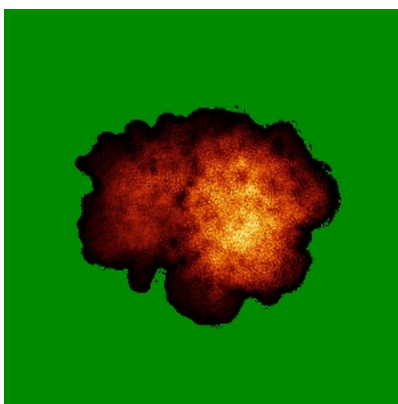
The images above show the largest variance slices of the map in three orthogonal directions.

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

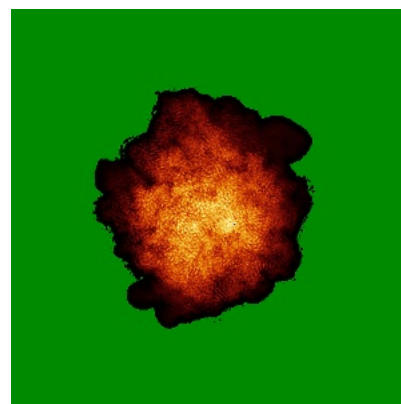
6.4.1 Primary map



X

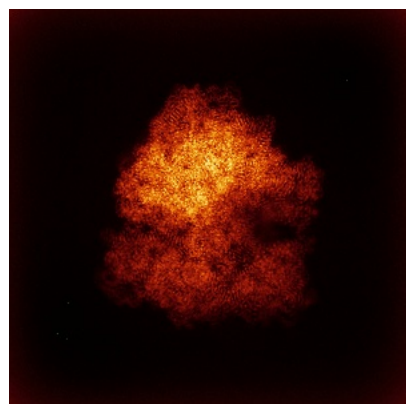


Y

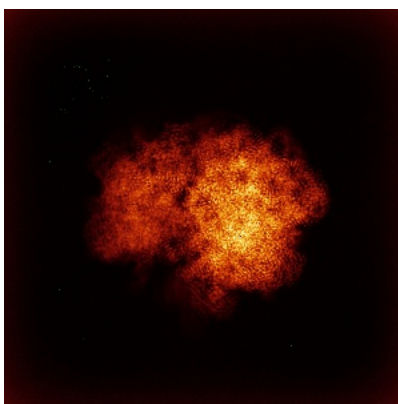


Z

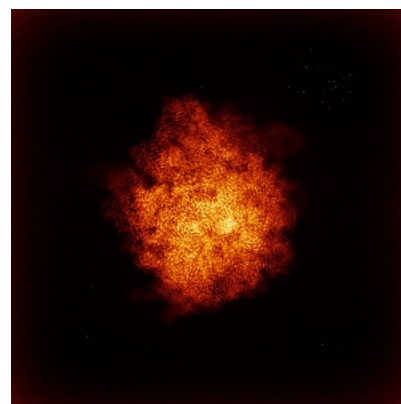
6.4.2 Raw map



X



Y

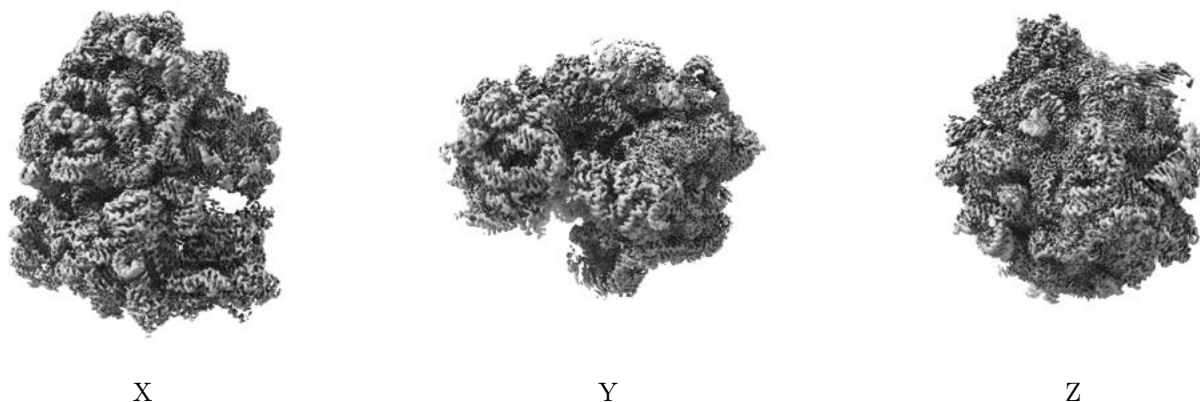


Z

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.

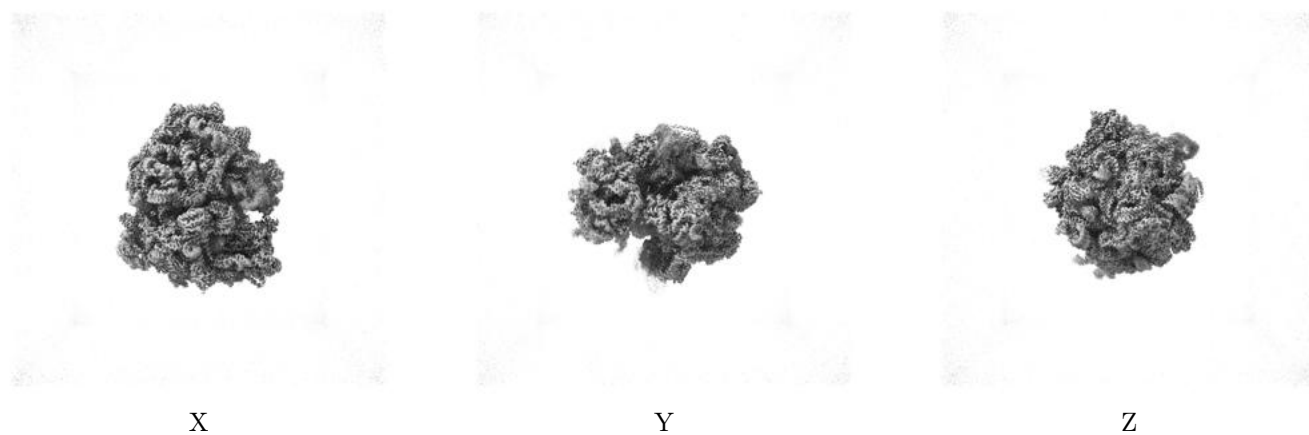
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.2. 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.

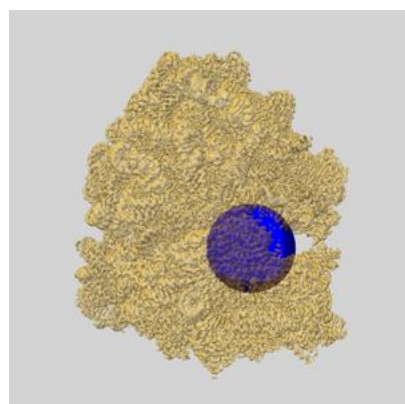
6.6 Mask visualisation [i](#)

This section shows the 3D surface view of the primary map at 50% transparency overlaid with the specified mask at 0% transparency

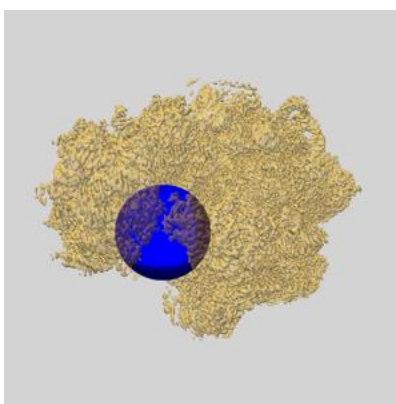
A mask typically either:

- Encompasses the whole structure
- Separates out a domain, a functional unit, a monomer or an area of interest from a larger structure

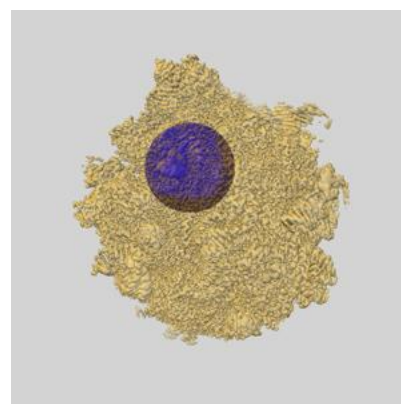
6.6.1 emd_15793_msk_1.map [i](#)



X

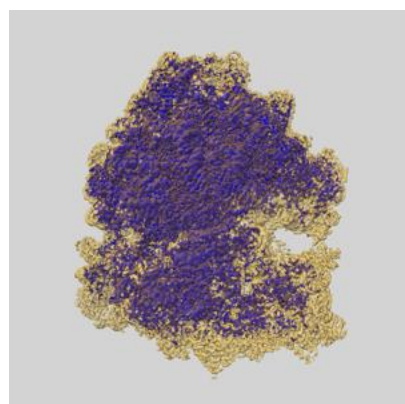


Y

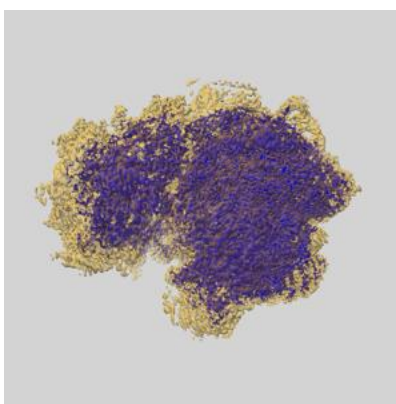


Z

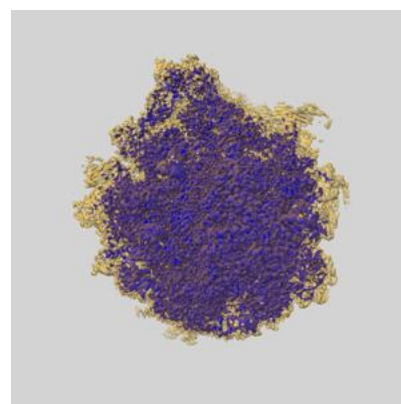
6.6.2 emd_15793_msk_2.map [i](#)



X



Y

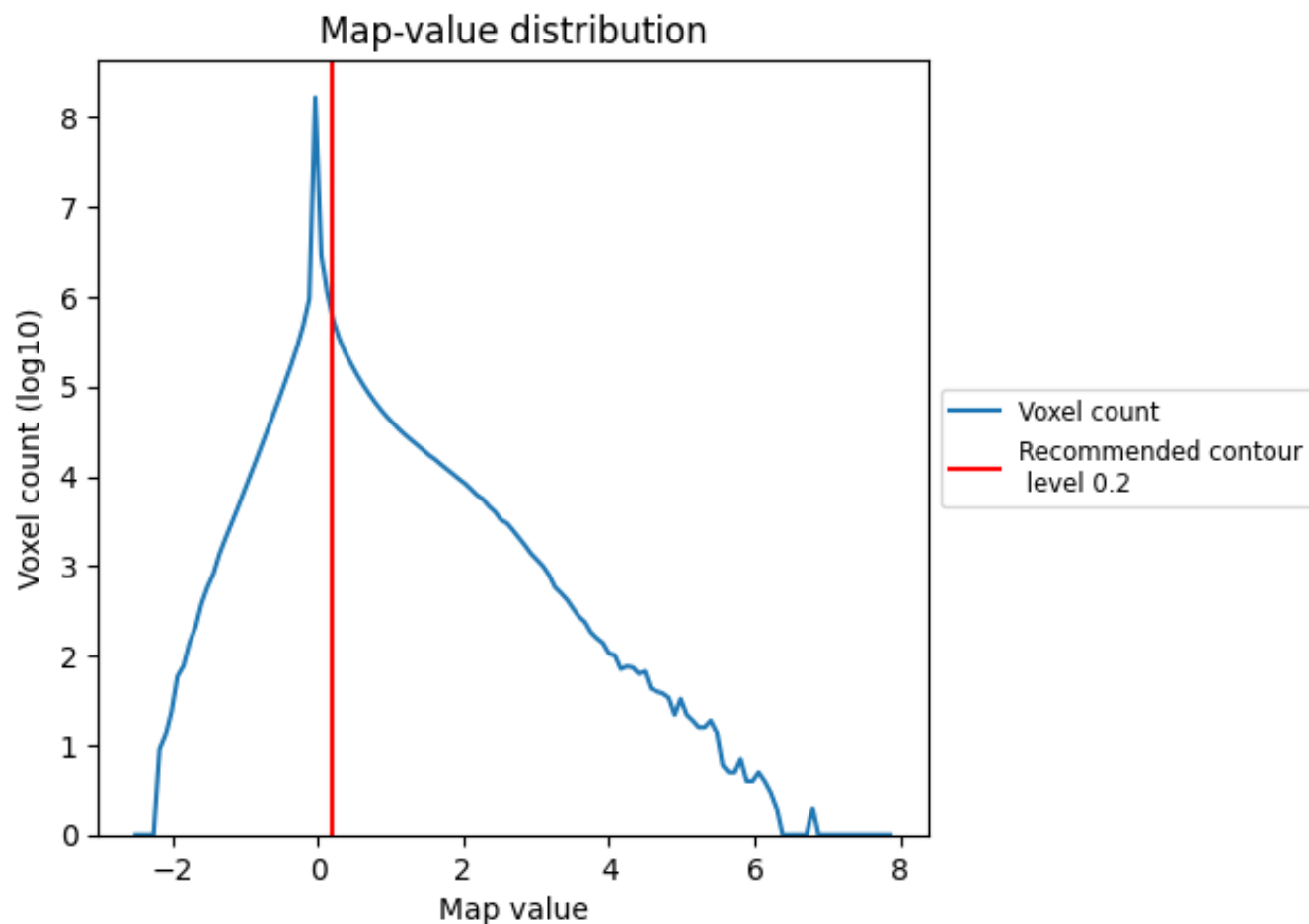


Z

7 Map analysis [i](#)

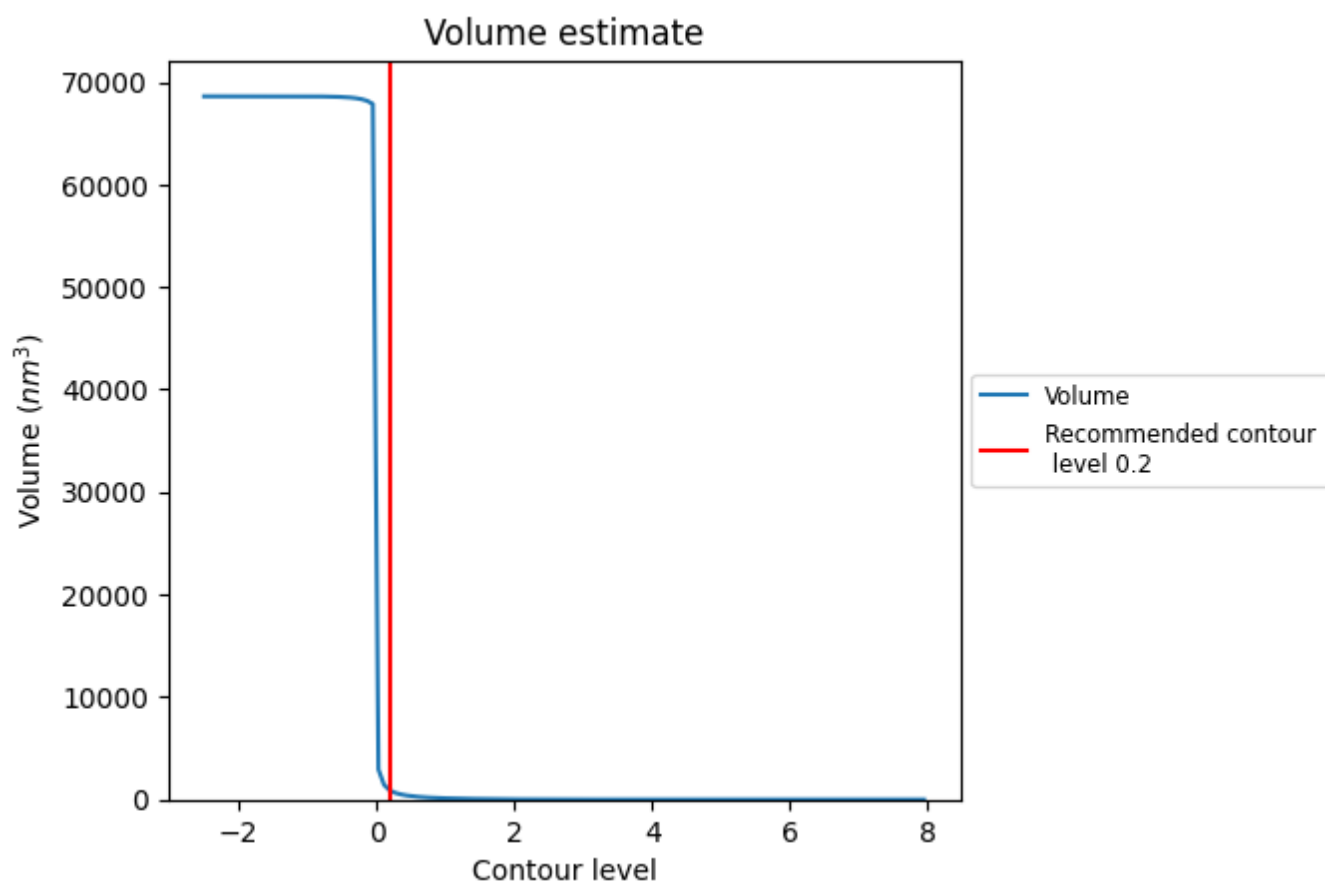
This section contains the results of statistical analysis of the map.

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.

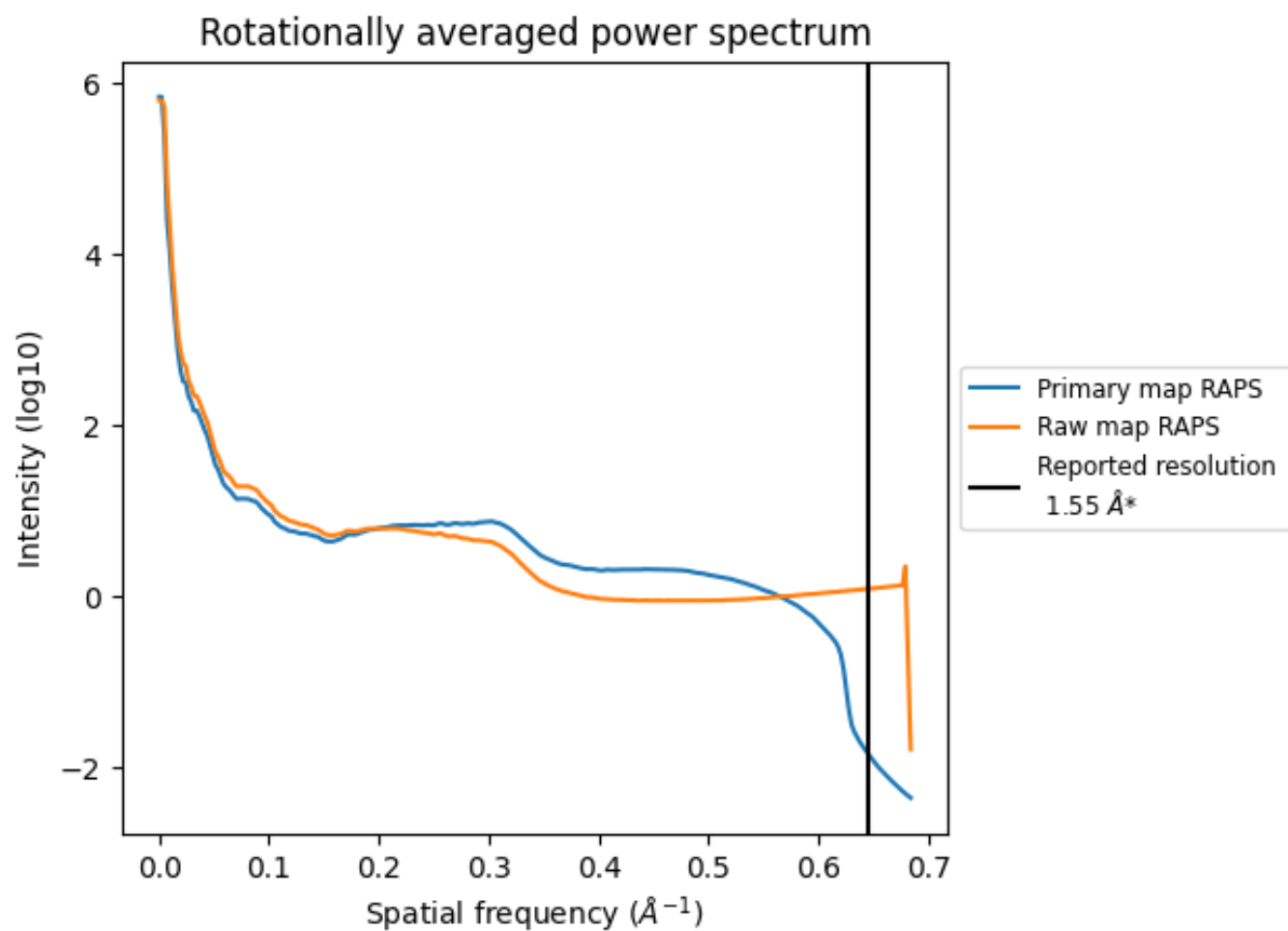
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 882 nm^3 ; this corresponds to an approximate mass of 797 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 ⓘ

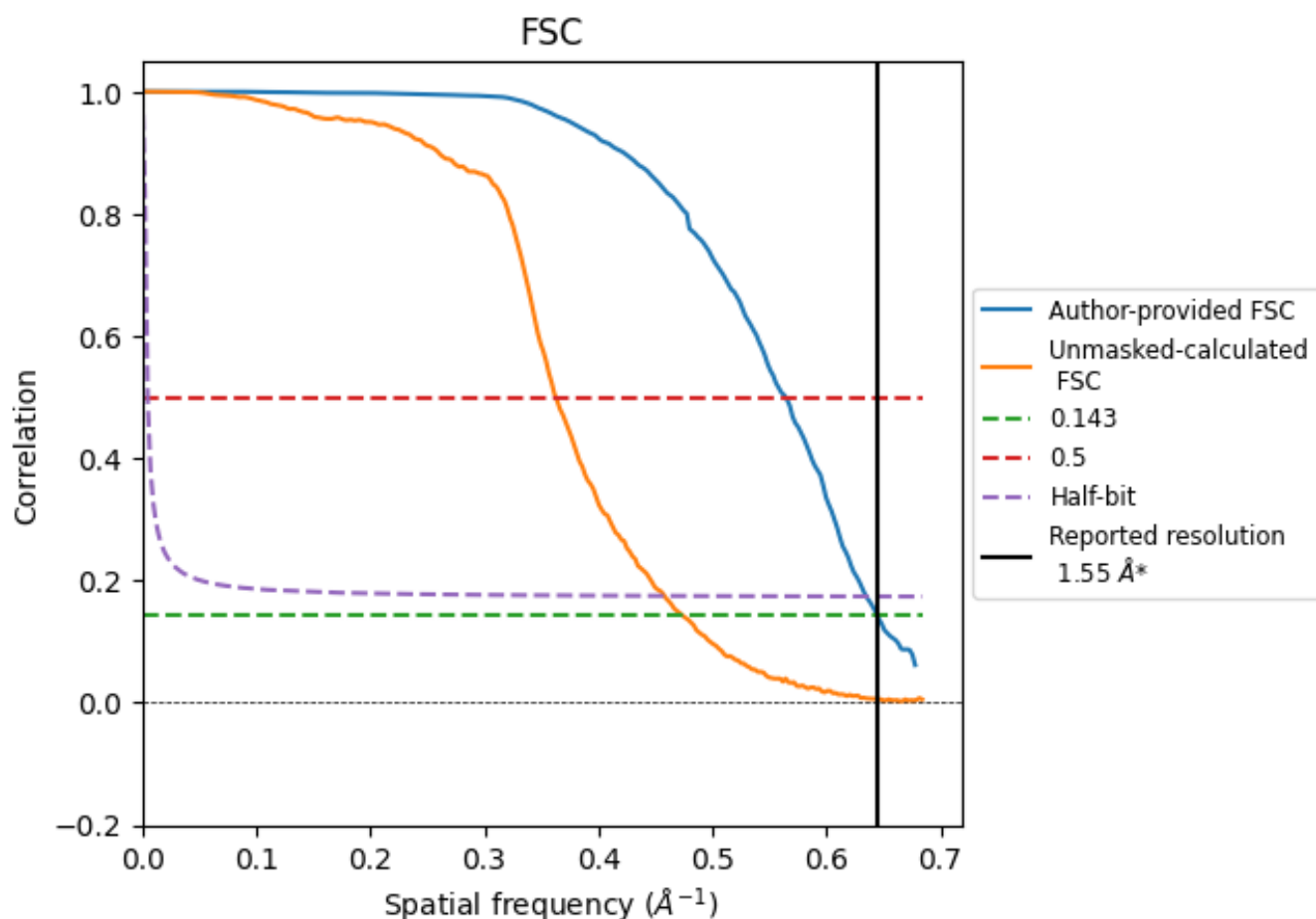


*Reported resolution corresponds to spatial frequency of 0.645 Å⁻¹

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 0.645 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

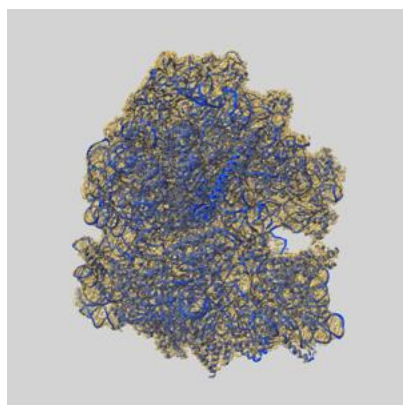
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	1.55	-	-
Author-provided FSC curve	1.55	1.77	1.57
Unmasked-calculated*	2.11	2.76	2.18

*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 2.11 differs from the reported value 1.55 by more than 10 %

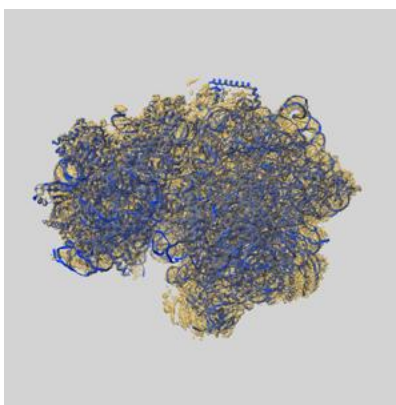
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-15793 and PDB model 8B0X. Per-residue inclusion information can be found in section [3](#) on page [17](#).

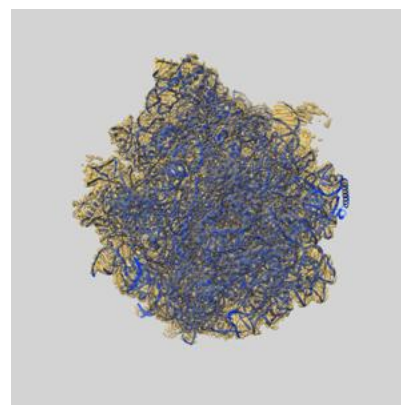
9.1 Map-model overlay [i](#)



X



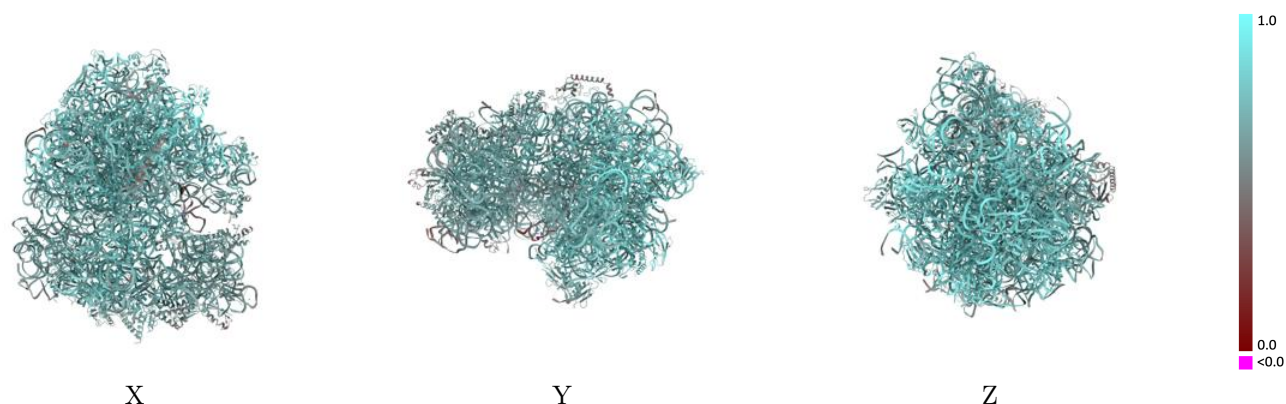
Y



Z

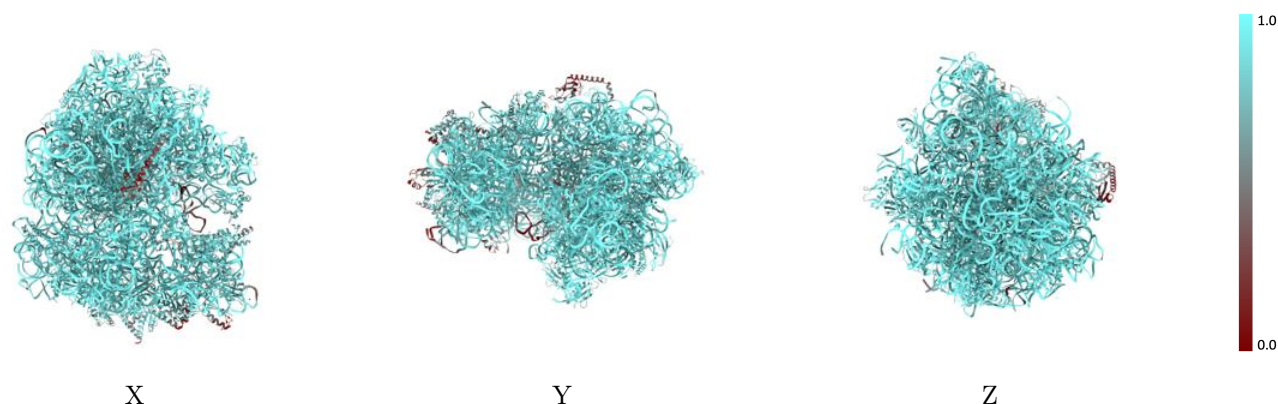
The images above show the 3D surface view of the map at the recommended contour level 0.2 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



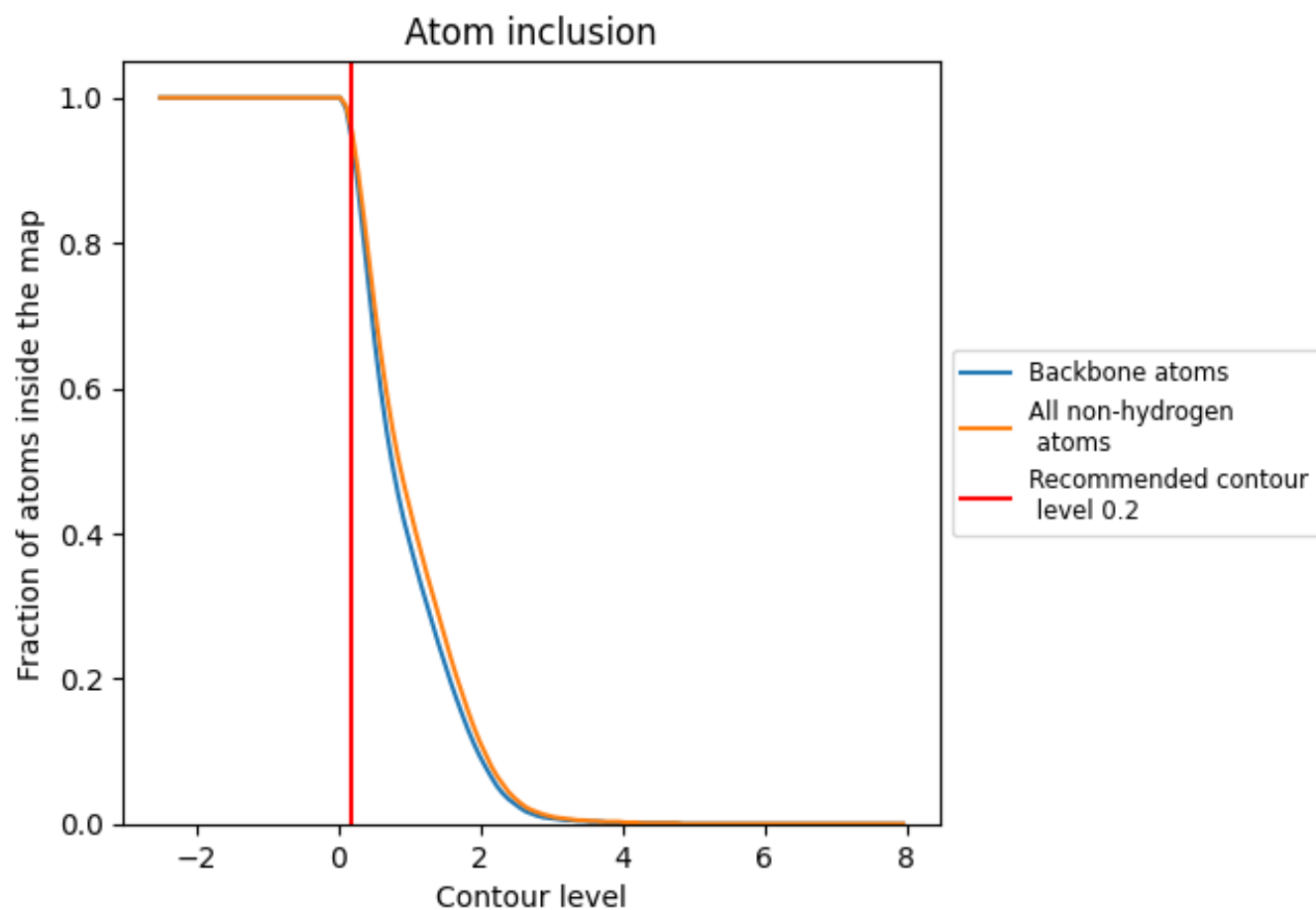
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.2).

9.4 Atom inclusion [i](#)



At the recommended contour level, 94% of all backbone atoms, 95% 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 (0.2) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.9520	 0.7670
0	 0.9660	 0.7960
1	 0.9890	 0.8720
2	 0.9960	 0.8680
3	 0.9760	 0.8180
A	 0.9670	 0.7190
B	 0.7510	 0.6360
C	 0.8730	 0.6750
D	 0.8490	 0.6700
E	 0.9500	 0.7590
F	 0.8970	 0.6690
G	 0.8030	 0.6290
H	 0.9460	 0.7570
I	 0.8410	 0.6420
J	 0.6610	 0.5890
K	 0.9300	 0.7010
L	 0.9450	 0.7400
M	 0.8060	 0.6250
N	 0.8770	 0.6570
O	 0.9380	 0.7380
P	 0.9400	 0.7110
Q	 0.8910	 0.6720
R	 0.9110	 0.7090
S	 0.7590	 0.5930
T	 0.9240	 0.7140
U	 0.6890	 0.6250
X	 0.9550	 0.7130
Z	 0.8790	 0.6240
a	 0.9830	 0.8160
b	 0.9820	 0.7330
c	 0.9910	 0.8530
d	 0.9770	 0.8440
e	 0.9460	 0.7990
f	 0.8600	 0.6460
g	 0.8880	 0.6690



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Chain	Atom inclusion	Q-score
h	 0.4910	 0.5610
i	 0.9860	 0.8440
j	 0.9780	 0.8380
k	 0.9780	 0.8330
l	 0.9850	 0.8240
m	 0.9990	 0.8690
n	 0.9470	 0.7390
o	 0.9640	 0.8300
p	 0.9970	 0.8660
q	 0.9550	 0.8090
r	 0.9810	 0.8460
s	 0.9380	 0.7810
t	 0.9230	 0.7540
u	 0.9310	 0.7500
v	 0.9720	 0.8250
w	 0.9770	 0.8260
x	 0.9240	 0.7440
y	 0.9610	 0.8310
z	 0.9600	 0.8260