Full wwPDB EM Validation Report

Nov 26, 2022 – 07:15 PM EST

PDB ID	:	7SYH
EMDB ID	:	EMD-25528
Title	:	Structure of the HCV IRES binding to the 40S ribosomal subunit, closed conformation. Structure 2(delta dII)
Authors	:	Brown, Z.P.; Abaeva, I.S.; De, S.; Hellen, C.U.T.; Pestova, T.V.; Frank, J.
Deposited on	:	2021-11-25
Resolution	:	4.60 Å(reported)
Based on initial models	:	5FLX, 5K0Y, 4UJD, 6D9J

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 (i) 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 (i)) were used in the production of this report:

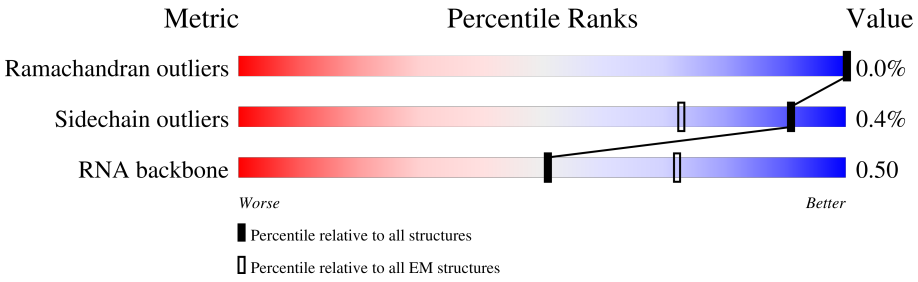
EMDB validation analysis	:	0.0.1.dev43
MolProbity	:	4.02b-467
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ	:	1.9.9
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.31.2

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 4.60 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.



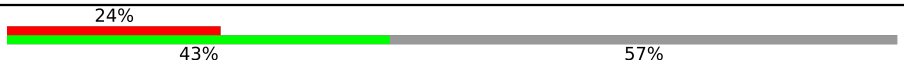
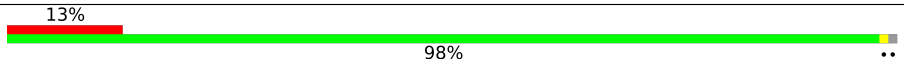
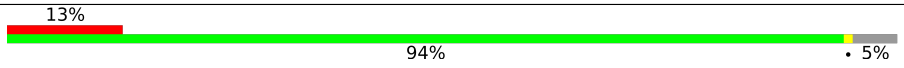
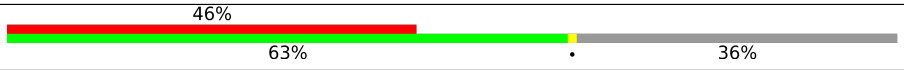
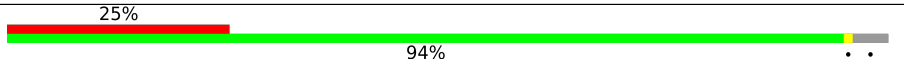
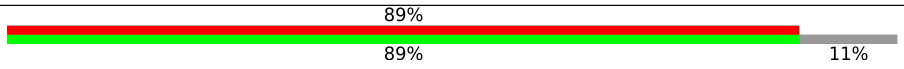
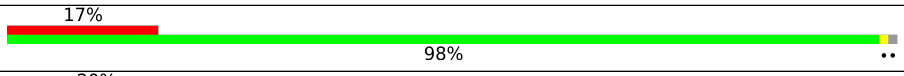
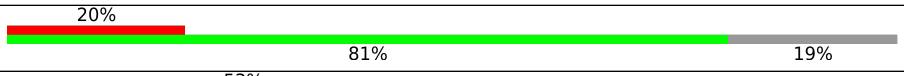
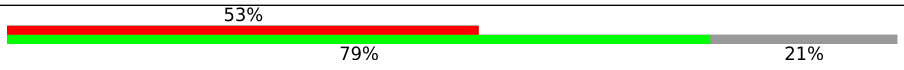
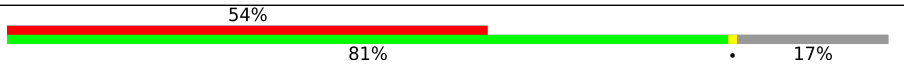
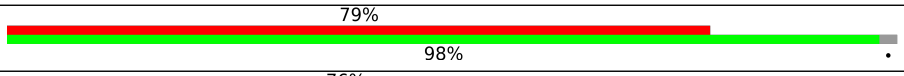
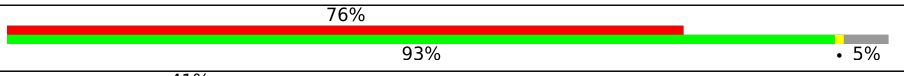
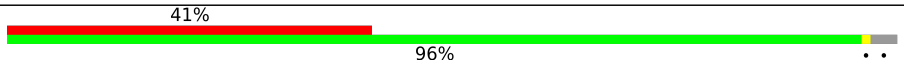
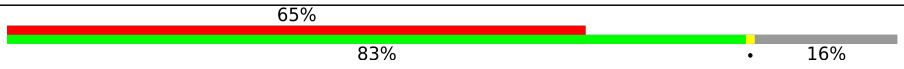
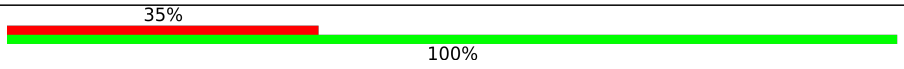
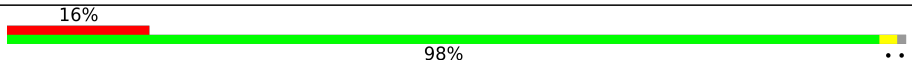
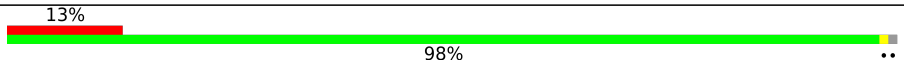
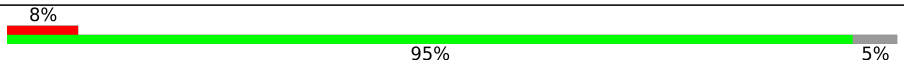
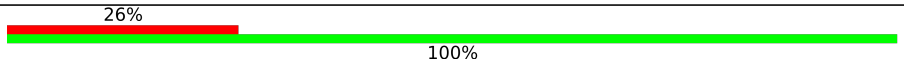
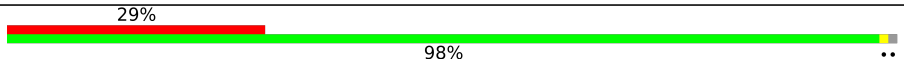
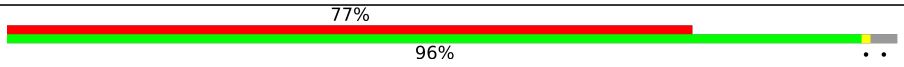
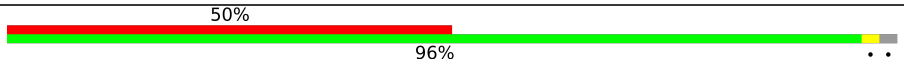
Metric	Whole archive (#Entries)	EM structures (#Entries)
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826
RNA backbone	4643	859

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

Mol	Chain	Length	Quality of chain
1	2	1870	9% 67% 23% • 9%
2	B	295	23% 73% • 26%
3	C	264	17% 80% • 19%
4	D	221	27% 100%
5	E	281	60% 80% • 19%
6	F	263	16% 100%
7	G	204	64% 94% 6%
8	H	249	24% 94% • 5%

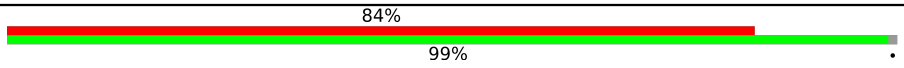
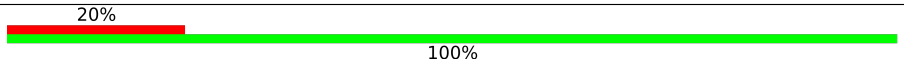
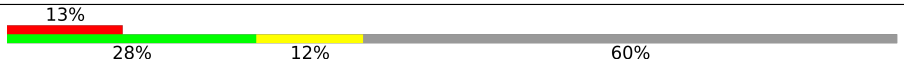
Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
9	I	432	
10	J	208	
11	K	194	
12	L	149	
13	M	158	
14	N	132	
15	O	151	
16	P	168	
17	Q	145	
18	R	172	
19	S	135	
20	T	152	
21	U	145	
22	V	119	
23	W	83	
24	X	130	
25	Y	143	
26	Z	131	
27	a	124	
28	b	101	
29	c	84	
30	d	69	
31	e	56	
32	f	133	
33	g	188	

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
34	h	317	
35	n	25	
36	z	400	

2 Entry composition

There are 37 unique types of molecules in this entry. The entry contains 78761 atoms, of which 0 are hydrogens and 0 are deuteriums.

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

- Molecule 1 is a RNA chain called 18S rRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	P		
1	2	1697	36227	16170	6504	11857	1696	0	0

- Molecule 2 is a protein called 40S ribosomal protein SA.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
2	B	217	1710	1086	300	316	8	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	114	THR	ALA	conflict	UNP G1TLT8

- Molecule 3 is a protein called 40S ribosomal protein S3a.

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

- Molecule 4 is a protein called uS5 (S2).

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
4	D	221	1716	1111	295	301	9	0	0

- Molecule 5 is a protein called uS3.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
5	E	228	1768	1126	318	316	8	0	0

- Molecule 6 is a protein called 40S ribosomal protein S4.

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

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
F	25	GLY	SER	conflict	UNP G1TK17
F	51	ARG	LYS	conflict	UNP G1TK17
F	78	THR	ALA	conflict	UNP G1TK17
F	156	VAL	MET	conflict	UNP G1TK17

- Molecule 7 is a protein called uS7.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	G	191	Total	C	N	O	S	0	0
			1509	943	286	273	7		

- Molecule 8 is a protein called eS6.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	H	237	Total	C	N	O	S	0	0
			1923	1200	387	329	7		

- Molecule 9 is a protein called 40S ribosomal protein S7.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	I	185	Total	C	N	O	S	0	0
			1488	952	271	264	1		

- Molecule 10 is a protein called eS8.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	J	206	Total	C	N	O	S	0	0
			1686	1058	332	291	5		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
J	47	ARG	GLY	conflict	UNP G1TJW1

- Molecule 11 is a protein called uS4.

Mol	Chain	Residues	Atoms					AltConf	Trace
11	K	185	Total	C	N	O	S	0	0
			1525	969	306	248	2		

- Molecule 12 is a protein called eS10.

Mol	Chain	Residues	Atoms					AltConf	Trace
12	L	96	Total	C	N	O	S	0	0
			810	530	143	131	6		

- Molecule 13 is a protein called uS17.

Mol	Chain	Residues	Atoms					AltConf	Trace
13	M	151	Total	C	N	O	S	0	0
			1233	785	231	211	6		

- Molecule 14 is a protein called eS12.

Mol	Chain	Residues	Atoms					AltConf	Trace
14	N	117	Total	C	N	O	S	0	0
			908	570	161	169	8		

- Molecule 15 is a protein called uS15.

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

- Molecule 16 is a protein called uS11.

Mol	Chain	Residues	Atoms					AltConf	Trace
16	P	136	Total	C	N	O	S	0	0
			1016	621	199	190	6		

- Molecule 17 is a protein called uS19.

Mol	Chain	Residues	Atoms					AltConf	Trace
17	Q	115	Total	C	N	O	S	0	0
			956	610	176	163	7		

- Molecule 18 is a protein called uS9.

Mol	Chain	Residues	Atoms					AltConf	Trace
18	R	142	Total	C	N	O	S	0	0
			1128	717	213	195	3		

- Molecule 19 is a protein called eS17.

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

- Molecule 20 is a protein called uS13.

Mol	Chain	Residues	Atoms					AltConf	Trace
20	T	144	Total	C	N	O	S	0	0
			1190	746	241	202	1		

- Molecule 21 is a protein called eS19.

Mol	Chain	Residues	Atoms					AltConf	Trace
21	U	141	Total	C	N	O	S	0	0
			1097	688	211	195	3		

- Molecule 22 is a protein called uS10.

Mol	Chain	Residues	Atoms					AltConf	Trace
22	V	100	Total	C	N	O	S	0	0
			795	498	152	141	4		

- Molecule 23 is a protein called 40S ribosomal protein S21.

Mol	Chain	Residues	Atoms					AltConf	Trace
23	W	83	Total	C	N	O	S	0	0
			636	393	117	121	5		

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
W	3	ASN	SER	conflict	UNP G1TM82
W	4	ASP	ASN	conflict	UNP G1TM82
W	33	GLN	PRO	conflict	UNP G1TM82
W	50	PHE	SER	conflict	UNP G1TM82
W	75	ALA	SER	conflict	UNP G1TM82
W	76	ASP	HIS	conflict	UNP G1TM82

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
W	81	LYS	GLN	conflict	UNP G1TM82

- Molecule 24 is a protein called uS8.

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

- Molecule 25 is a protein called uS12.

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

- Molecule 26 is a protein called 40S ribosomal protein S24.

Mol	Chain	Residues	Atoms					AltConf	Trace
26	Z	124	Total	C	N	O	S	0	0
			1011	640	198	168	5		

- Molecule 27 is a protein called 40S ribosomal protein S25.

Mol	Chain	Residues	Atoms					AltConf	Trace
27	a	77	Total	C	N	O	S	0	0
			614	393	114	106	1		

- Molecule 28 is a protein called 40S ribosomal protein S26.

Mol	Chain	Residues	Atoms					AltConf	Trace
28	b	101	Total	C	N	O	S	0	0
			814	507	170	132	5		

- Molecule 29 is a protein called eS27.

Mol	Chain	Residues	Atoms					AltConf	Trace
29	c	83	Total	C	N	O	S	0	0
			651	408	121	115	7		

- Molecule 30 is a protein called eS28.

Mol	Chain	Residues	Atoms					AltConf	Trace
30	d	67	Total	C	N	O	S	0	0
			530	321	108	99	2		

- Molecule 31 is a protein called eS29.

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

- Molecule 32 is a protein called eS30.

Mol	Chain	Residues	Atoms					AltConf	Trace
32	f	57	Total	C	N	O	S	0	0
			457	282	101	73	1		

- Molecule 33 is a protein called 40S ribosomal protein S27a.

Mol	Chain	Residues	Atoms					AltConf	Trace
33	g	68	Total	C	N	O	S	0	0
			555	351	103	94	7		

- Molecule 34 is a protein called RACK1.

Mol	Chain	Residues	Atoms					AltConf	Trace
34	h	313	Total	C	N	O	S	0	0
			2436	1535	424	465	12		

- Molecule 35 is a protein called 60s ribosomal protein l41.

Mol	Chain	Residues	Atoms					AltConf	Trace
35	n	25	Total	C	N	O	S	0	0
			239	145	64	27	3		

- Molecule 36 is a RNA chain called HCV IRES.

Mol	Chain	Residues	Atoms					AltConf	Trace
36	z	162	Total	C	N	O	P	0	0
			3465	1542	621	1140	162		

- Molecule 37 is ZINC ION (three-letter code: ZN) (formula: Zn).

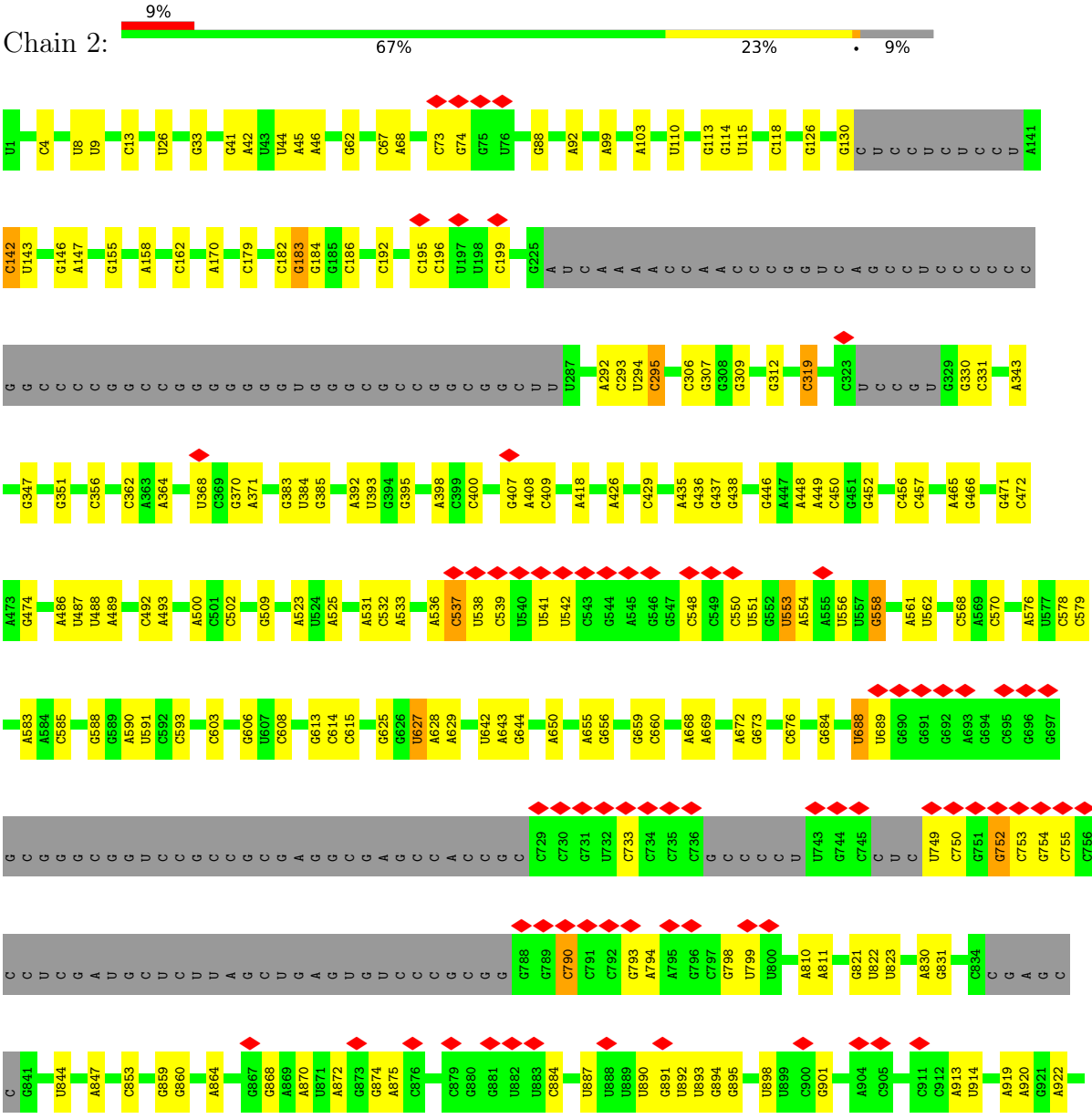
Mol	Chain	Residues	Atoms	AltConf
37	b	1	Total Zn 1 1	0
37	g	1	Total Zn 1 1	0

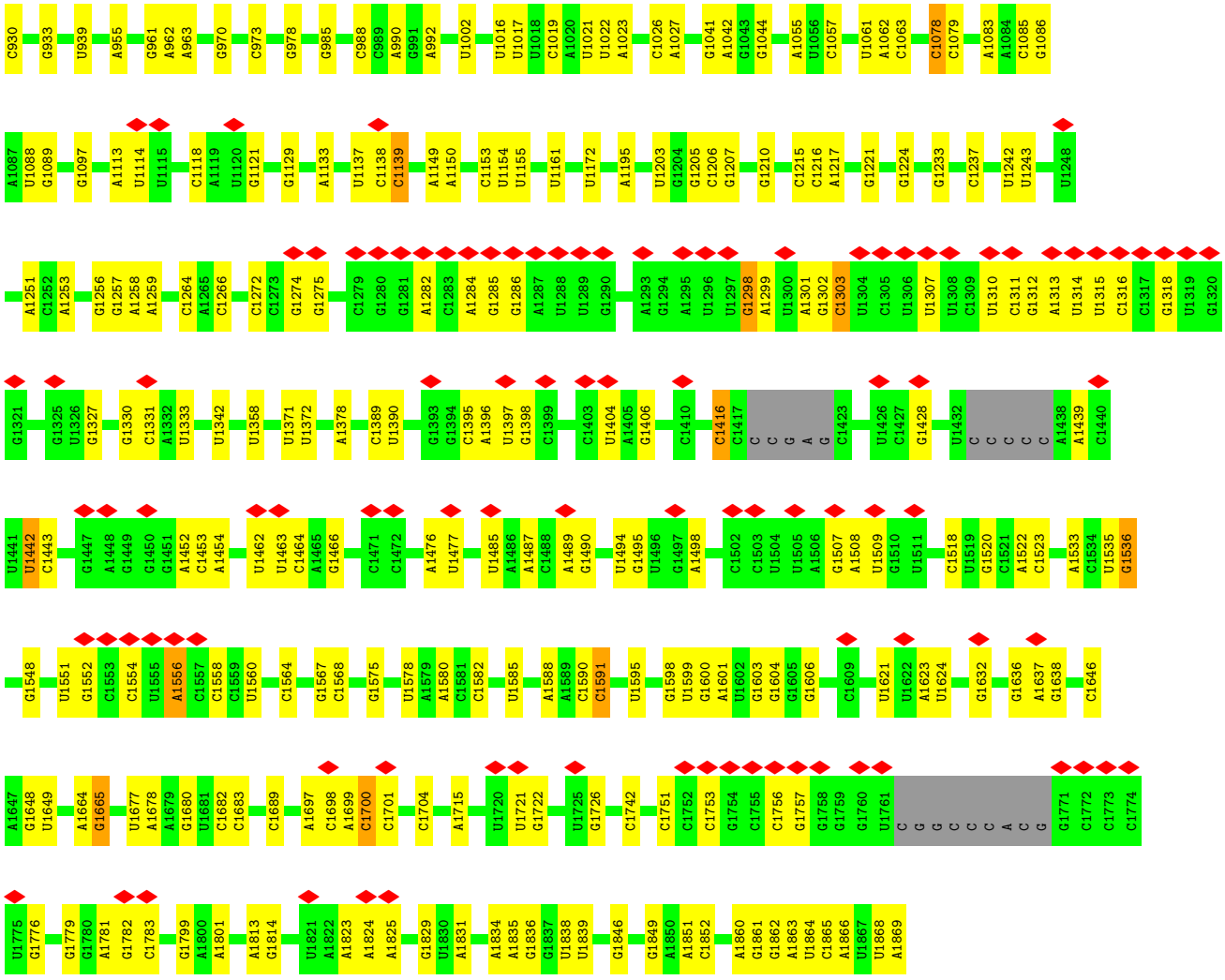
3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion in map density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red diamond above a residue indicates a poor fit to the EM map for this residue (all-atom inclusion < 40%). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: 18S rRNA

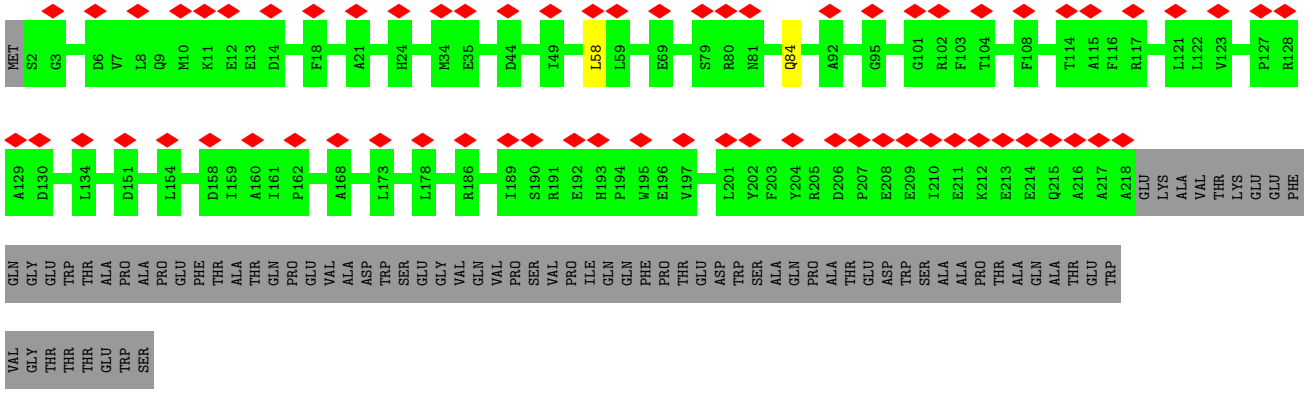
Chain 2:





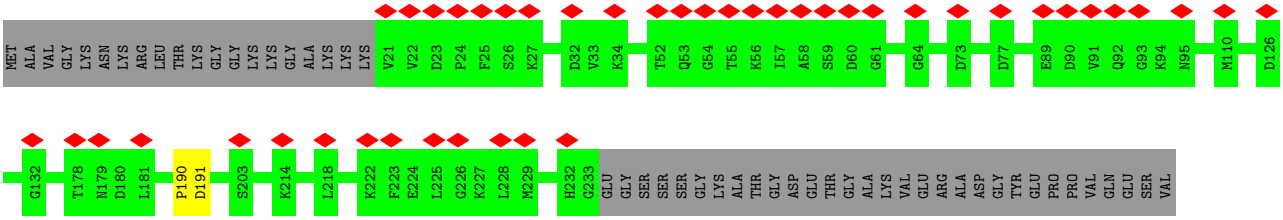
• Molecule 2: 40S ribosomal protein SA

Chain B:

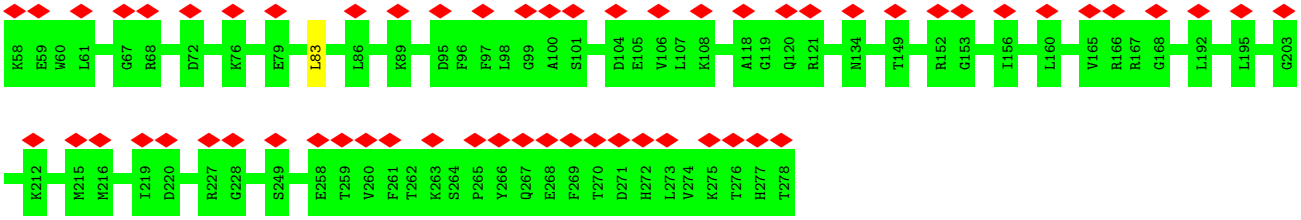


• Molecule 3: 40S ribosomal protein S3a

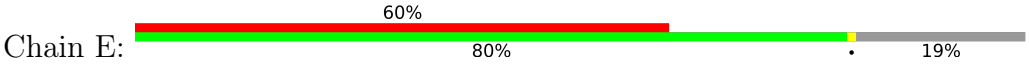
Chain C:

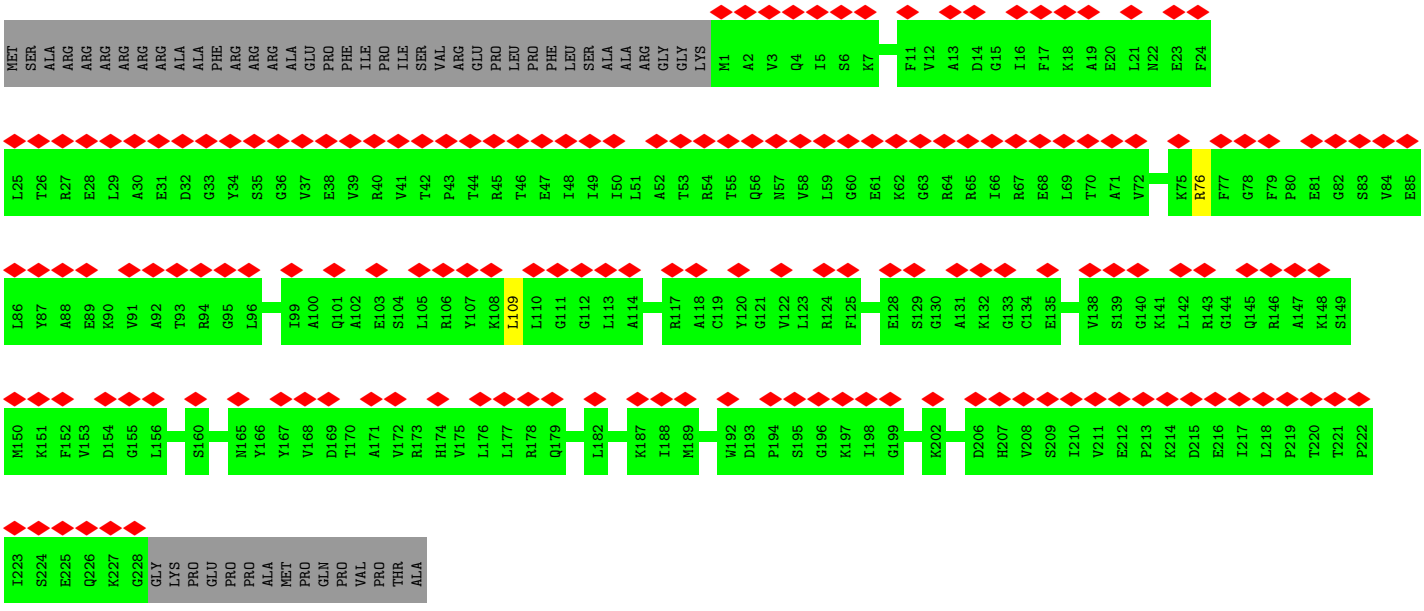


• Molecule 4: uS5 (S2)

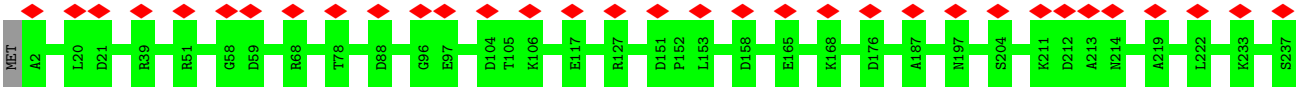


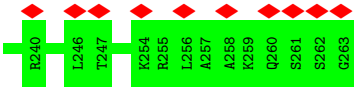
• Molecule 5: uS3





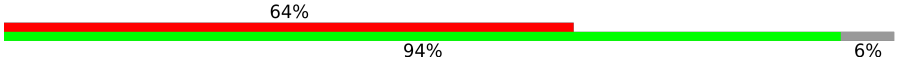
• Molecule 6: 40S ribosomal protein S4

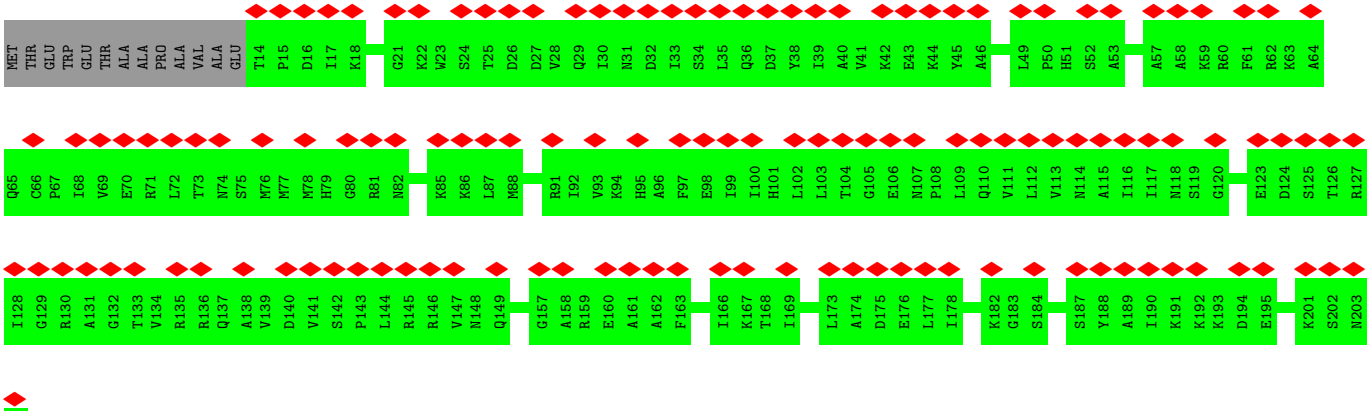




• Molecule 7: uS7

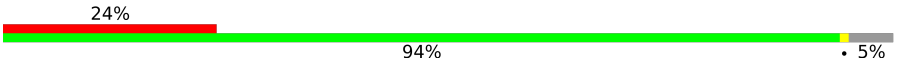
Chain G:

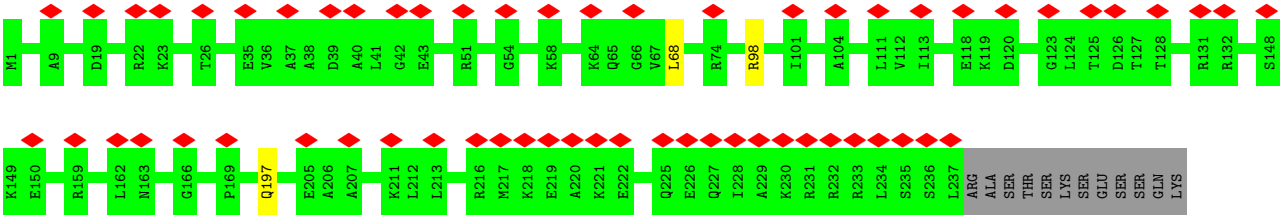




• Molecule 8: eS6

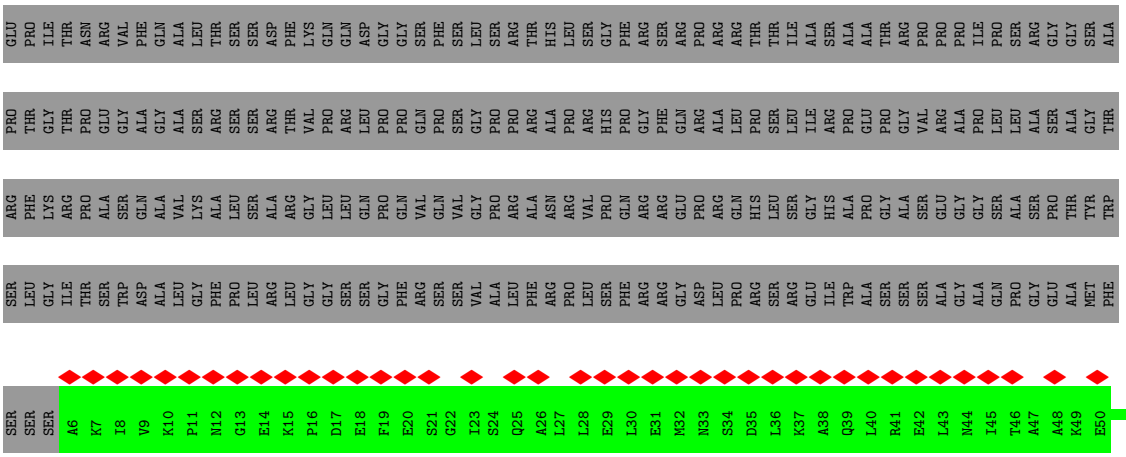
Chain H:

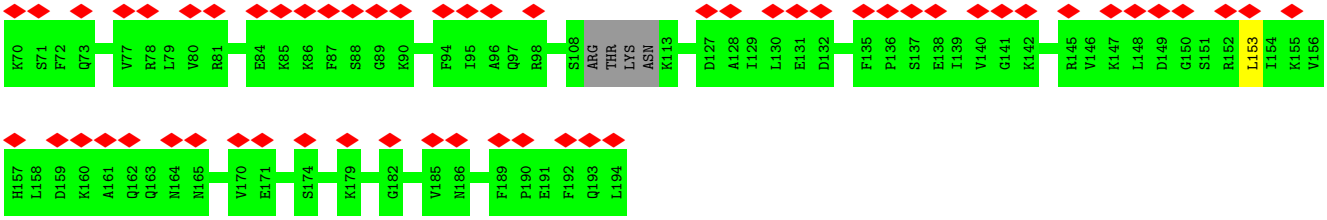




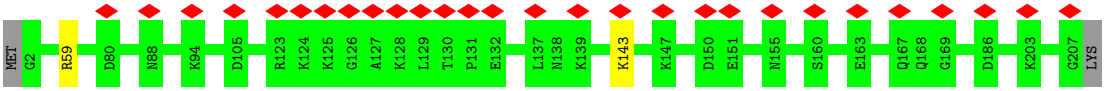
• Molecule 9: 40S ribosomal protein S7

Chain I:

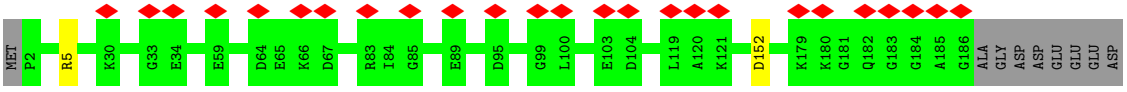




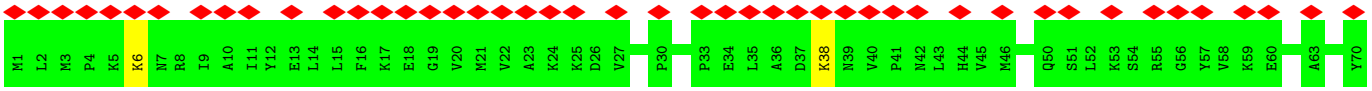
• Molecule 10: eS8

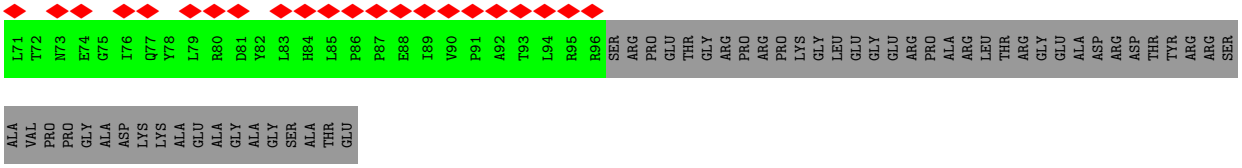


• Molecule 11: uS4

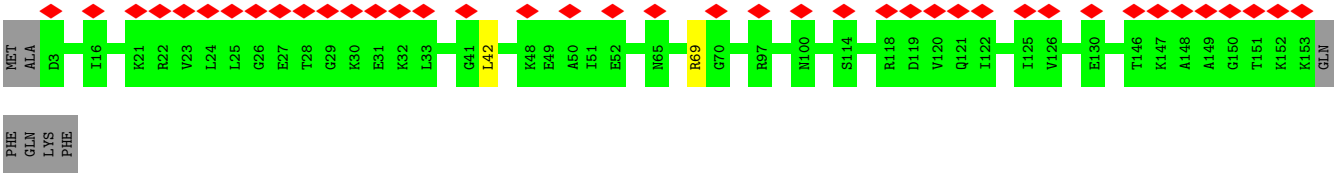


• Molecule 12: eS10

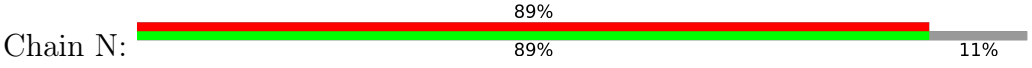


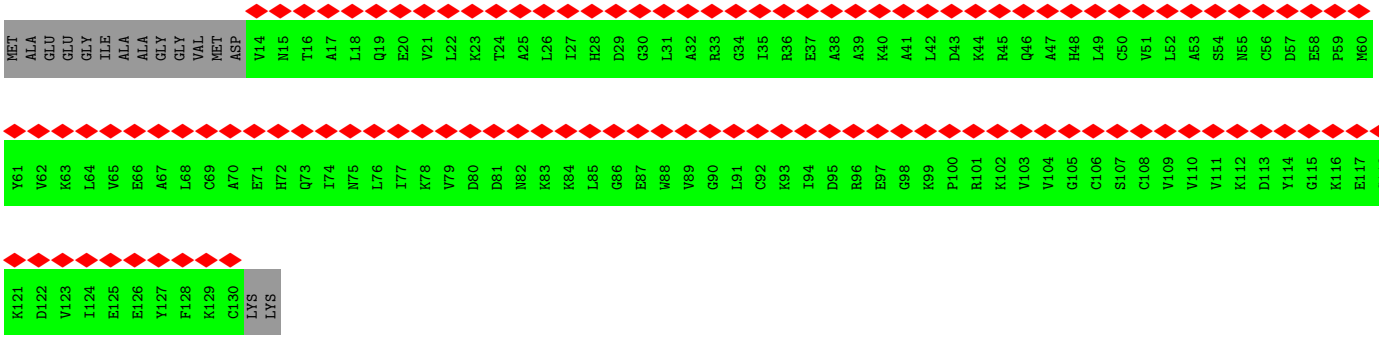


• Molecule 13: uS17

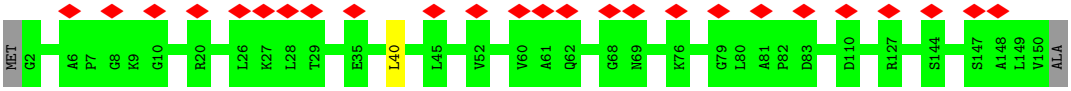


• Molecule 14: eS12

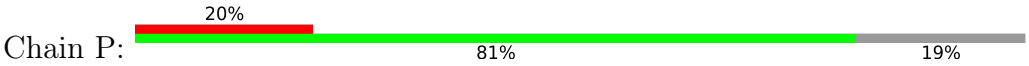


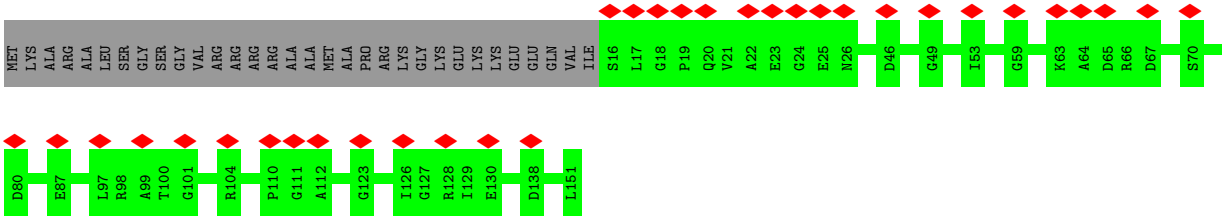


• Molecule 15: uS15

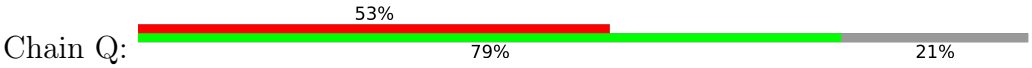


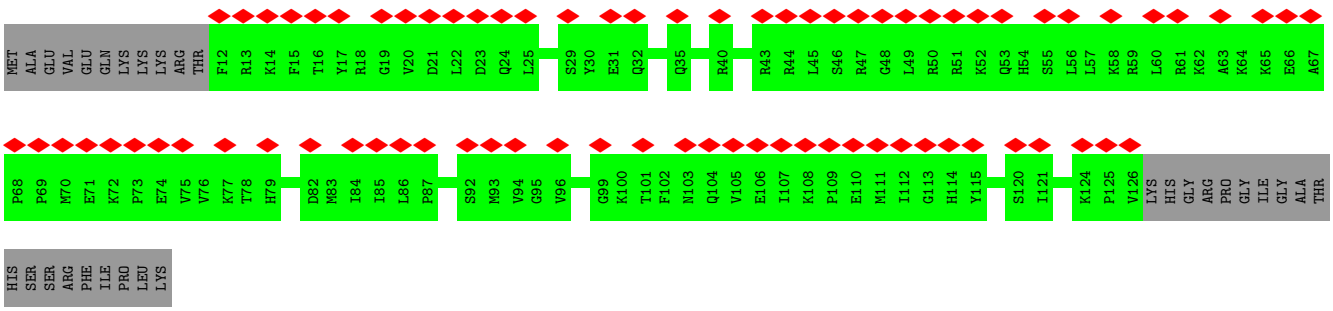
• Molecule 16: uS11



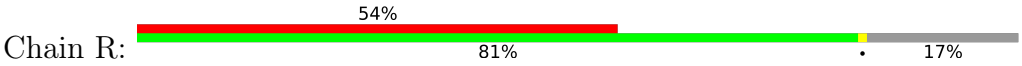


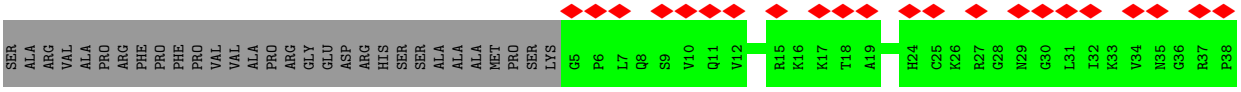
• Molecule 17: uS19

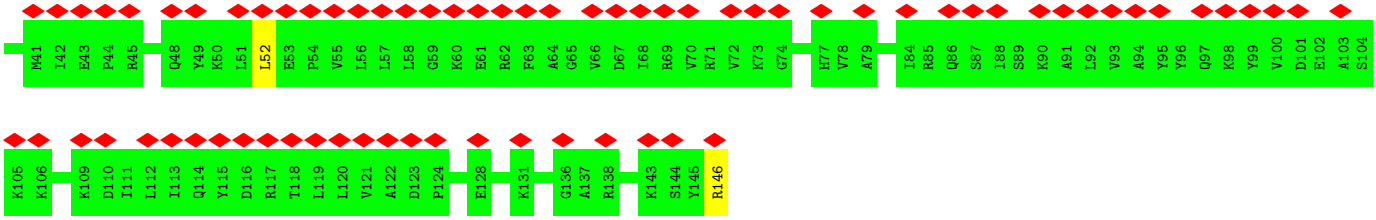




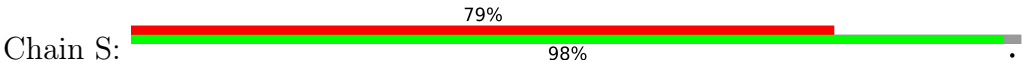
• Molecule 18: uS9

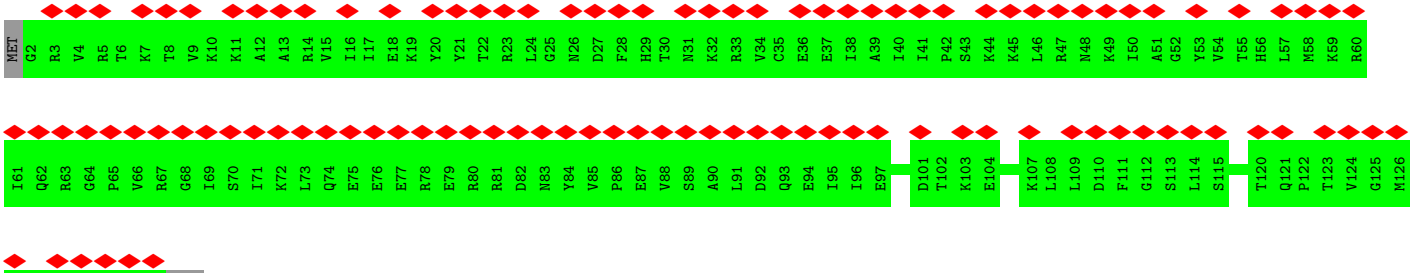




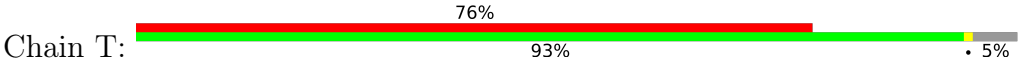


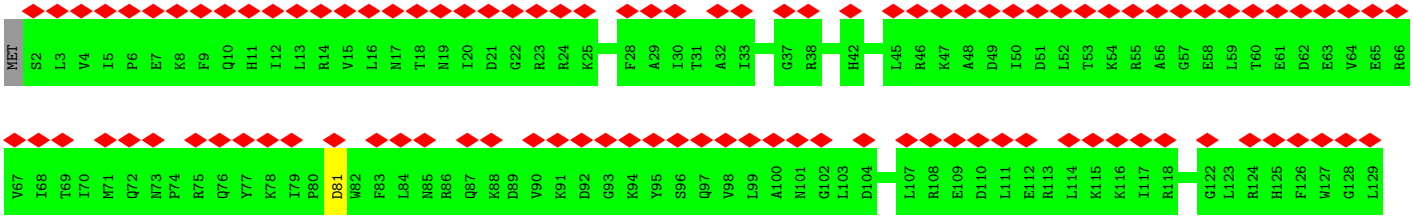
• Molecule 19: eS17



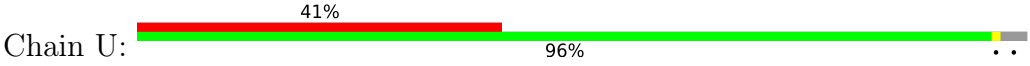


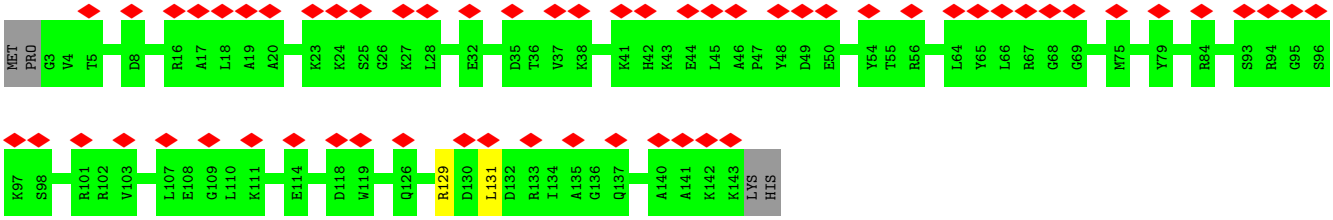
• Molecule 20: uS13



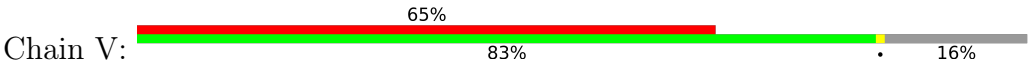


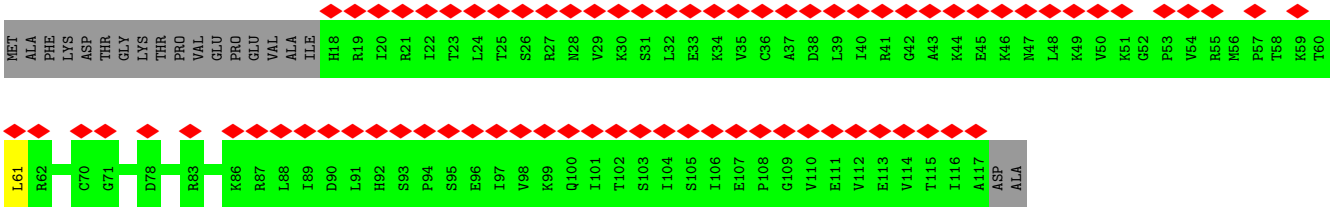
• Molecule 21: eS19



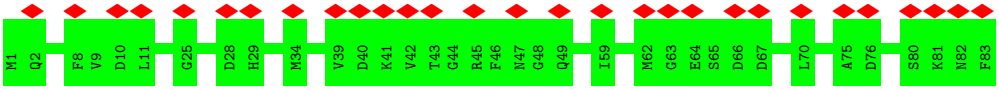


• Molecule 22: uS10

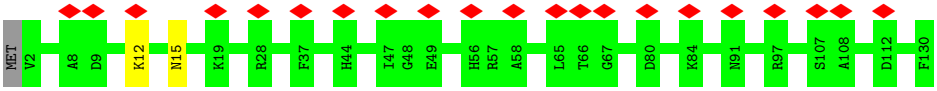




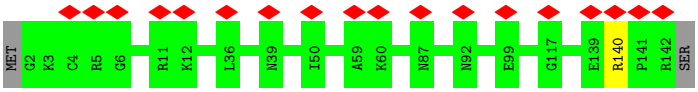
• Molecule 23: 40S ribosomal protein S21



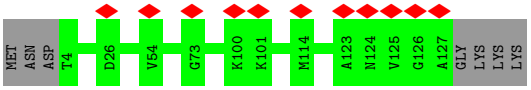
• Molecule 24: uS8



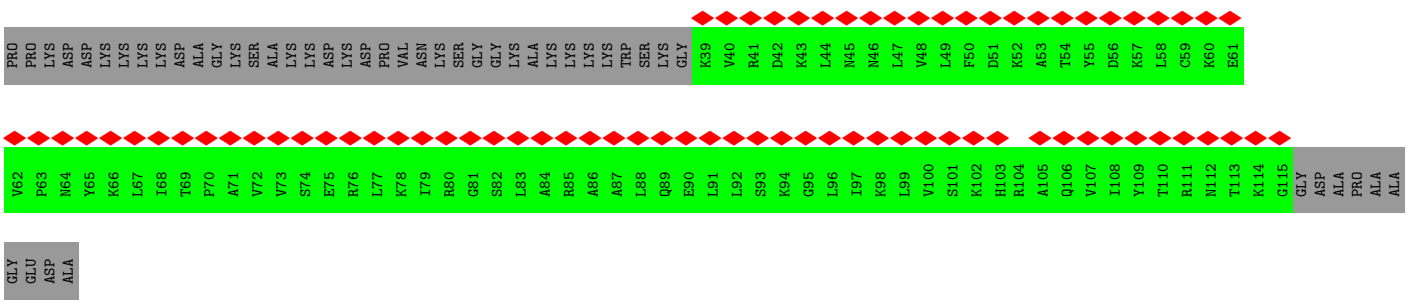
• Molecule 25: uS12



• Molecule 26: 40S ribosomal protein S24

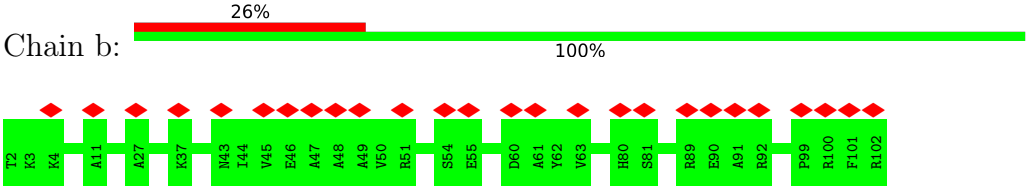


• Molecule 27: 40S ribosomal protein S25



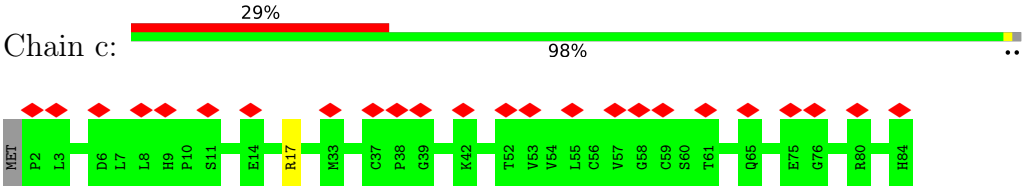
• Molecule 28: 40S ribosomal protein S26

Chain b:



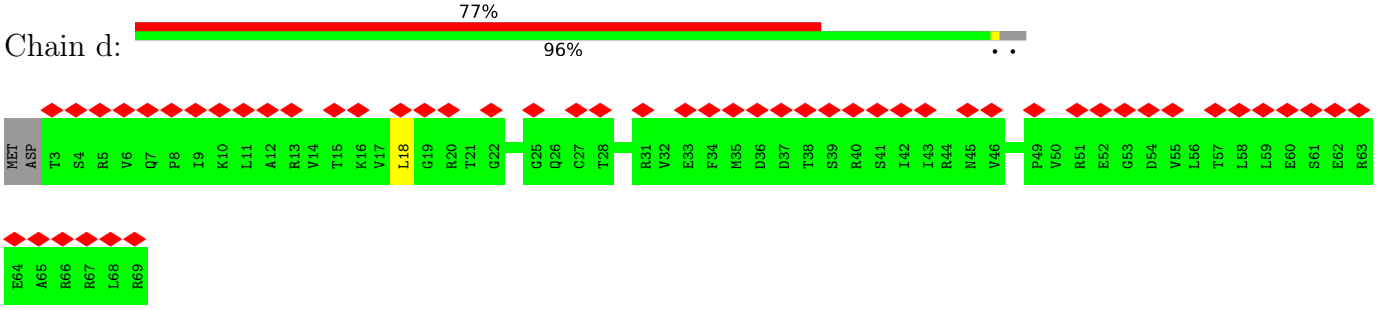
• Molecule 29: eS27

Chain c:



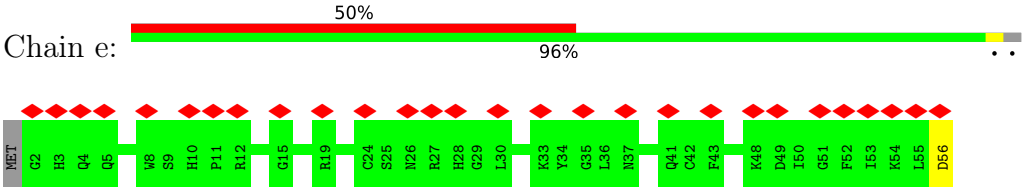
• Molecule 30: eS28

Chain d:



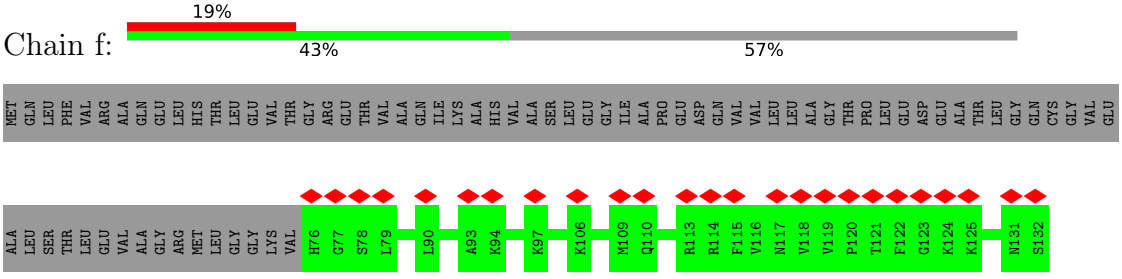
• Molecule 31: eS29

Chain e:



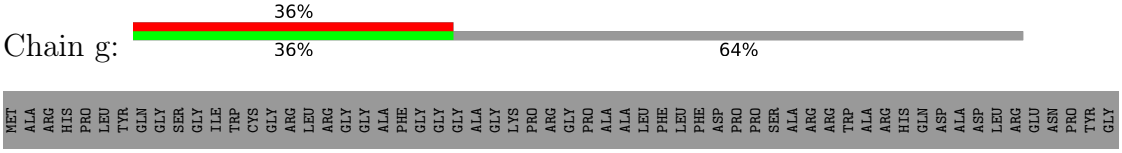
• Molecule 32: eS30

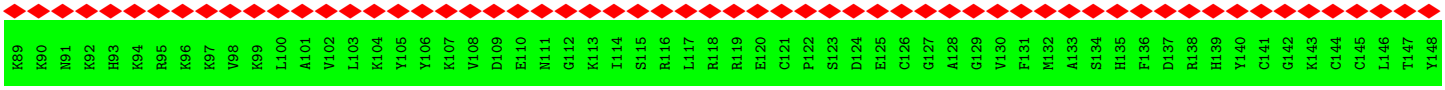
Chain f:

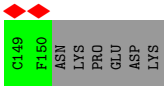


• Molecule 33: 40S ribosomal protein S27a

Chain g:

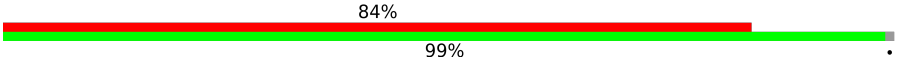


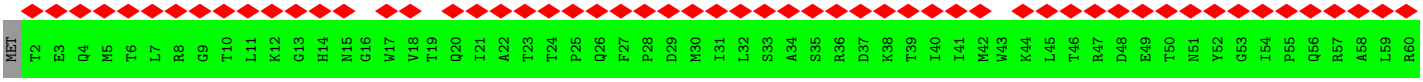


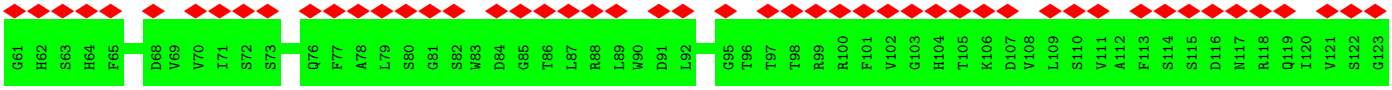


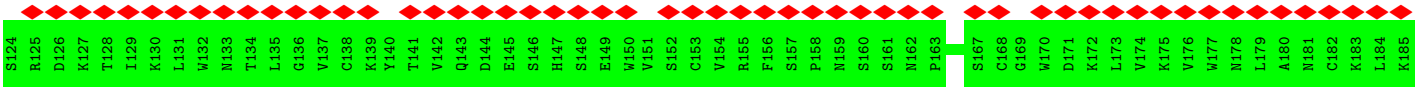
• Molecule 34: RACK1

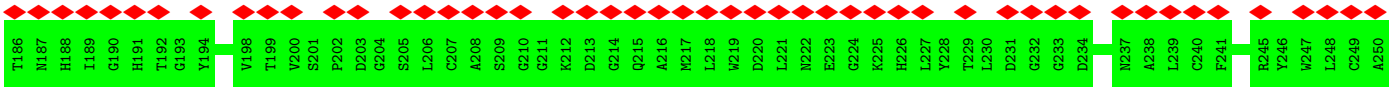
Chain h:

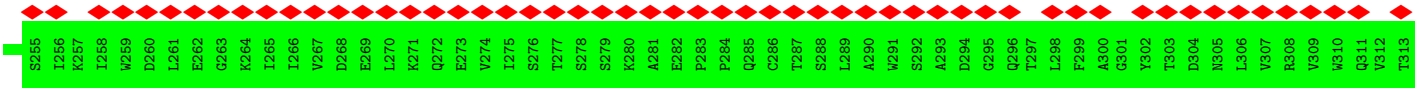






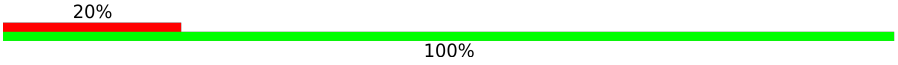


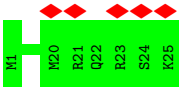




• Molecule 35: 60s ribosomal protein l41

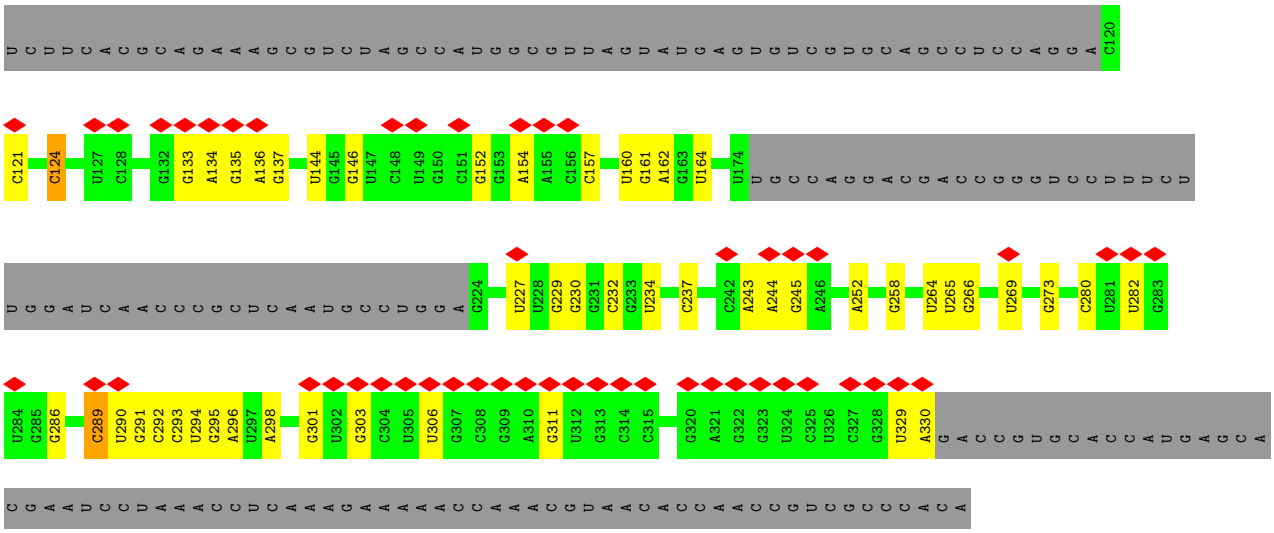
Chain n:





• Molecule 36: HCV IRES

Chain z:


U C U U C A C G C A G A A A G C G U C U U A G C C C A U G G C G U U A G U A U G A G U G U C G U G C A G C C U C C C A G G A C120
C121 C124 U127 C128 G132 G133 A134 G135 A136 G137 U144 G145 G146 U147 C148 U149 G150 C151 G152 G153 A154 A155 C156 C157 U160 G161 A162 G163 U164 U174
U G C C A G G A C G A C C G G G U C C U U U C U
U G G A U C A A C C C G C U C A A U G C C C U G G A G224
U227 U228 G229 G230 G231 C232 G233 U234 C237 C242 A243 A244 G245 A246 A252 G258 U264 U265 G266 U269 G273 C280 U281 U282 G283
U284 G285 G286 C289 U290 G291 C292 C293 U294 G295 A296 U297 A298 G301 U302 G303 C304 U305 U306 G307 C308 G309 A310 G311 U312 G313 C314 C315 G320 A321 G322 G323 U324 C325 U326 C327 G328 U329 A330
G A C C G U G C A C C A A U G A G C A
C G A A U C C U A A A C C U C A A A A G A A A A A A C C C A C G U A A C C A A C C C G U C G C C C A C A

4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	28684	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TECNAI F30	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/Å^2$)	70.9	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	52000	Depositor
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.036	Depositor
Minimum map value	-0.020	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.002	Depositor
Recommended contour level	0.00883	Depositor
Map size (Å)	380.0, 380.0, 380.0	wwPDB
Map dimensions	400, 400, 400	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.95, 0.95, 0.95	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZN

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	#$\|Z\|>5$	RMSZ	#$\|Z\|>5$
1	2	0.27	0/40506	1.01	166/63123 (0.3%)
2	B	0.28	0/1747	0.65	1/2374 (0.0%)
3	C	0.27	0/1756	0.60	0/2350
4	D	0.27	0/1753	0.62	1/2369 (0.0%)
5	E	0.29	0/1796	0.62	1/2417 (0.0%)
6	F	0.27	0/2118	0.58	0/2849
7	G	0.26	0/1531	0.60	0/2059
8	H	0.26	0/1946	0.64	2/2590 (0.1%)
9	I	0.28	0/1510	0.63	1/2022 (0.0%)
10	J	0.25	0/1715	0.58	0/2287
11	K	0.29	0/1550	0.65	1/2069 (0.0%)
12	L	0.26	0/834	0.57	0/1125
13	M	0.28	0/1254	0.64	1/1677 (0.1%)
14	N	0.31	0/918	0.64	0/1233
15	O	0.26	0/1226	0.60	1/1649 (0.1%)
16	P	0.29	0/1029	0.65	0/1380
17	Q	0.27	0/974	0.67	0/1301
18	R	0.26	0/1146	0.68	1/1534 (0.1%)
19	S	0.30	0/1082	0.69	0/1452
20	T	0.27	0/1208	0.69	1/1618 (0.1%)
21	U	0.25	0/1115	0.59	1/1493 (0.1%)
22	V	0.26	0/805	0.69	1/1081 (0.1%)
23	W	0.27	0/643	0.59	0/860
24	X	0.30	0/1051	0.73	0/1406
25	Y	0.26	0/1116	0.58	0/1490
26	Z	0.28	0/1028	0.61	0/1366
27	a	0.27	0/620	0.62	0/831
28	b	0.27	0/828	0.64	0/1109
29	c	0.27	0/665	0.61	0/891
30	d	0.25	0/532	0.74	1/712 (0.1%)
31	e	0.27	0/470	0.63	1/623 (0.2%)
32	f	0.25	0/462	0.60	0/607

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\#\|Z\| > 5$	RMSZ	$\#\|Z\| > 5$
33	g	0.24	0/567	0.50	0/753
34	h	0.26	0/2493	0.60	0/3394
35	n	0.25	0/240	0.74	0/305
36	z	0.27	0/3871	1.00	12/6034 (0.2%)
All	All	0.27	0/84105	0.86	193/122433 (0.2%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
3	C	0	1
21	U	0	1
All	All	0	2

There are no bond length outliers.

All (193) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	2	1453	C	N1-C2-O2	12.46	126.38	118.90
1	2	1453	C	C2-N1-C1'	11.43	131.37	118.80
1	2	1078	C	N1-C2-O2	10.66	125.30	118.90
1	2	1078	C	N3-C2-O2	-10.63	114.46	121.90
1	2	1453	C	N3-C2-O2	-9.83	115.02	121.90
1	2	853	C	C2-N1-C1'	9.22	128.94	118.80
1	2	853	C	N1-C2-O2	9.11	124.37	118.90
1	2	537	C	N1-C2-O2	8.89	124.24	118.90
1	2	457	C	N3-C2-O2	-8.81	115.73	121.90
1	2	1314	U	C2-N1-C1'	8.78	128.23	117.70
1	2	538	U	N1-C2-O2	8.65	128.85	122.80
20	T	81	ASP	CB-CG-OD2	8.38	125.84	118.30
1	2	539	C	N1-C2-O2	8.34	123.90	118.90
1	2	1314	U	N1-C2-O2	8.18	128.52	122.80
1	2	1700	C	C6-N1-C2	-8.11	117.06	120.30
1	2	1591	C	N1-C2-O2	8.07	123.74	118.90
1	2	1689	C	N1-C2-O2	8.07	123.74	118.90
36	z	292	C	N1-C2-O2	8.06	123.74	118.90
1	2	538	U	N3-C2-O2	-8.05	116.56	122.20
1	2	1078	C	C6-N1-C2	-7.85	117.16	120.30
1	2	1453	C	C6-N1-C1'	-7.85	111.39	120.80
1	2	1314	U	N3-C2-O2	-7.82	116.73	122.20

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	2	1453	C	C6-N1-C2	-7.67	117.23	120.30
1	2	1139	C	N1-C2-O2	7.65	123.49	118.90
1	2	1624	U	C2-N1-C1'	7.63	126.86	117.70
1	2	1442	U	N1-C2-O2	7.61	128.13	122.80
1	2	537	C	N3-C2-O2	-7.59	116.59	121.90
1	2	1700	C	N1-C2-O2	7.57	123.44	118.90
1	2	1442	U	C2-N1-C1'	7.55	126.76	117.70
1	2	853	C	N3-C2-O2	-7.39	116.72	121.90
1	2	823	U	C2-N1-C1'	7.32	126.48	117.70
1	2	1442	U	N3-C2-O2	-7.29	117.10	122.20
1	2	578	C	N1-C2-O2	7.26	123.26	118.90
36	z	124	C	N1-C2-O2	7.25	123.25	118.90
1	2	537	C	C2-N1-C1'	7.20	126.72	118.80
1	2	1303	C	N1-C2-O2	7.20	123.22	118.90
8	H	68	LEU	CA-CB-CG	7.18	131.82	115.30
1	2	1019	C	N1-C2-O2	7.15	123.19	118.90
18	R	52	LEU	CA-CB-CG	7.15	131.74	115.30
1	2	1019	C	N3-C2-O2	-7.08	116.94	121.90
1	2	1303	C	C2-N1-C1'	6.98	126.48	118.80
11	K	152	ASP	CB-CG-OD1	6.98	124.58	118.30
1	2	1139	C	C6-N1-C2	-6.98	117.51	120.30
1	2	823	U	N3-C2-O2	-6.97	117.32	122.20
1	2	1700	C	N3-C2-O2	-6.94	117.04	121.90
1	2	1298	G	N3-C4-N9	6.90	130.14	126.00
1	2	538	U	C2-N1-C1'	6.90	125.98	117.70
1	2	1689	C	N3-C2-O2	-6.86	117.10	121.90
1	2	1298	G	N3-C4-C5	-6.86	125.17	128.60
1	2	823	U	N1-C2-O2	6.79	127.55	122.80
1	2	539	C	N3-C2-O2	-6.70	117.21	121.90
1	2	1753	C	N3-C2-O2	-6.67	117.23	121.90
1	2	195	C	C2-N1-C1'	6.64	126.11	118.80
1	2	1139	C	N3-C2-O2	-6.62	117.27	121.90
1	2	539	C	C2-N1-C1'	6.60	126.06	118.80
1	2	752	G	P-O3'-C3'	6.60	127.62	119.70
1	2	585	C	N1-C2-O2	6.58	122.85	118.90
36	z	292	C	N3-C2-O2	-6.57	117.30	121.90
1	2	1665	G	O5'-P-OP1	-6.56	99.79	105.70
1	2	183	G	N3-C4-C5	-6.55	125.33	128.60
1	2	659	G	C4-N9-C1'	6.53	134.99	126.50
1	2	859	G	N3-C4-C5	-6.52	125.34	128.60
1	2	558	G	C4-N9-C1'	-6.50	118.06	126.50
1	2	553	U	P-O3'-C3'	6.49	127.49	119.70

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	2	749	U	N3-C2-O2	-6.47	117.67	122.20
1	2	790	C	N1-C2-O2	6.47	122.78	118.90
1	2	1063	C	N1-C2-O2	6.46	122.78	118.90
1	2	688	U	P-O3'-C3'	6.45	127.44	119.70
1	2	1298	G	C4-N9-C1'	6.45	134.88	126.50
1	2	457	C	C6-N1-C2	-6.43	117.73	120.30
1	2	558	G	C8-N9-C1'	6.43	135.36	127.00
1	2	1443	C	N1-C2-O2	6.43	122.76	118.90
1	2	1453	C	C5-C6-N1	6.42	124.21	121.00
1	2	1591	C	N3-C2-O2	-6.41	117.41	121.90
1	2	568	C	N1-C2-O2	6.40	122.74	118.90
1	2	853	C	C6-N1-C1'	-6.39	113.13	120.80
1	2	179	C	N1-C2-O2	6.38	122.73	118.90
1	2	558	G	N3-C4-N9	-6.35	122.19	126.00
1	2	1591	C	C6-N1-C2	-6.32	117.77	120.30
1	2	183	G	C4-N9-C1'	6.18	134.54	126.50
1	2	1624	U	N1-C2-O2	6.12	127.09	122.80
1	2	859	G	C4-N9-C1'	6.11	134.45	126.50
1	2	1078	C	C2-N1-C1'	6.10	125.51	118.80
1	2	578	C	N3-C2-O2	-6.09	117.64	121.90
1	2	539	C	C6-N1-C2	-6.07	117.87	120.30
1	2	1139	C	C2-N1-C1'	6.07	125.47	118.80
1	2	195	C	C6-N1-C2	-6.06	117.88	120.30
1	2	1057	C	C2-N1-C1'	6.05	125.46	118.80
1	2	1689	C	C2-N1-C1'	6.05	125.46	118.80
1	2	859	G	N3-C4-N9	6.03	129.62	126.00
1	2	1689	C	C6-N1-C2	-6.01	117.89	120.30
36	z	124	C	C2-N1-C1'	6.01	125.41	118.80
1	2	356	C	C2-N1-C1'	6.00	125.39	118.80
1	2	1079	C	N1-C2-O2	5.99	122.49	118.90
1	2	1518	C	C2-N1-C1'	5.98	125.38	118.80
1	2	537	C	C6-N1-C2	-5.95	117.92	120.30
1	2	853	C	C6-N1-C2	-5.92	117.93	120.30
1	2	749	U	N1-C2-O2	5.91	126.94	122.80
1	2	1591	C	C2-N1-C1'	5.86	125.25	118.80
1	2	1314	U	C6-N1-C1'	-5.85	113.00	121.20
1	2	356	C	N1-C2-O2	5.85	122.41	118.90
1	2	1624	U	N3-C2-O2	-5.85	118.11	122.20
1	2	894	G	C4-N9-C1'	5.85	134.10	126.50
1	2	183	G	N3-C4-N9	5.84	129.50	126.00
1	2	293	C	N1-C2-O2	5.83	122.39	118.90
36	z	124	C	N3-C2-O2	-5.82	117.82	121.90

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	2	579	C	N1-C2-O2	5.81	122.39	118.90
36	z	280	C	N1-C2-O2	5.79	122.38	118.90
36	z	292	C	C2-N1-C1'	5.79	125.17	118.80
1	2	1016	U	C2-N1-C1'	5.78	124.64	117.70
1	2	13	C	C2-N1-C1'	5.77	125.15	118.80
1	2	627	U	P-O3'-C3'	5.76	126.61	119.70
1	2	183	G	C2-N3-C4	5.74	114.77	111.90
1	2	1443	C	C2-N1-C1'	5.73	125.11	118.80
9	I	153	LEU	CA-CB-CG	5.71	128.44	115.30
1	2	1485	U	N1-C2-O2	5.71	126.80	122.80
30	d	18	LEU	CA-CB-CG	5.69	128.39	115.30
1	2	1700	C	C5-C6-N1	5.68	123.84	121.00
1	2	659	G	C8-N9-C1'	-5.68	119.61	127.00
1	2	1303	C	N3-C2-O2	-5.68	117.93	121.90
1	2	1389	C	N1-C2-O2	5.67	122.30	118.90
1	2	1016	U	N1-C2-O2	5.66	126.76	122.80
1	2	457	C	N1-C2-O2	5.64	122.28	118.90
1	2	118	C	N1-C2-O2	5.63	122.28	118.90
1	2	196	C	N1-C2-O2	5.63	122.28	118.90
1	2	1742	C	N1-C2-O2	5.62	122.27	118.90
1	2	570	C	N1-C2-O2	5.62	122.27	118.90
36	z	124	C	C6-N1-C2	-5.60	118.06	120.30
1	2	755	C	N1-C2-O2	5.57	122.24	118.90
1	2	1523	C	N1-C2-O2	5.56	122.24	118.90
15	O	40	LEU	CA-CB-CG	5.56	128.08	115.30
1	2	1442	U	C5-C6-N1	5.54	125.47	122.70
2	B	58	LEU	CA-CB-CG	5.52	128.00	115.30
22	V	61	LEU	CA-CB-CG	5.50	127.96	115.30
1	2	1057	C	N1-C2-O2	5.50	122.20	118.90
1	2	437	G	N1-C2-N2	-5.50	111.25	116.20
1	2	319	C	O5'-P-OP1	5.46	117.25	110.70
1	2	1485	U	N3-C2-O2	-5.45	118.39	122.20
1	2	585	C	N3-C2-O2	-5.45	118.09	121.90
1	2	1753	C	N1-C2-O2	5.44	122.17	118.90
1	2	1063	C	C6-N1-C2	-5.44	118.12	120.30
36	z	289	C	N1-C2-O2	5.41	122.14	118.90
1	2	659	G	N3-C4-N9	5.40	129.24	126.00
1	2	1518	C	N1-C2-O2	5.39	122.14	118.90
1	2	1016	U	N3-C2-O2	-5.39	118.42	122.20
1	2	1298	G	C8-N9-C1'	-5.39	120.00	127.00
1	2	1564	C	C2-N1-C1'	5.38	124.71	118.80
36	z	289	C	N3-C2-O2	-5.34	118.16	121.90

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	2	195	C	N1-C2-O2	5.34	122.10	118.90
1	2	1389	C	N3-C2-O2	-5.33	118.17	121.90
1	2	939	U	N1-C2-O2	5.30	126.51	122.80
1	2	456	C	N1-C2-O2	5.30	122.08	118.90
1	2	142	C	N1-C2-O2	5.30	122.08	118.90
1	2	1298	G	C2-N3-C4	5.29	114.55	111.90
13	M	42	LEU	CA-CB-CG	5.29	127.46	115.30
1	2	894	G	N3-C4-C5	-5.28	125.96	128.60
1	2	790	C	C5-C6-N1	5.28	123.64	121.00
1	2	1063	C	N3-C2-O2	-5.27	118.21	121.90
1	2	1331	C	N1-C2-O2	5.26	122.06	118.90
1	2	195	C	C5-C6-N1	5.25	123.62	121.00
1	2	1443	C	C6-N1-C2	-5.24	118.20	120.30
1	2	295	C	N1-C2-O2	5.24	122.04	118.90
1	2	1624	U	O4'-C1'-N1	5.23	112.39	108.20
1	2	568	C	N3-C2-O2	-5.23	118.24	121.90
1	2	894	G	N3-C4-N9	5.22	129.13	126.00
1	2	1591	C	C5-C6-N1	5.21	123.61	121.00
1	2	615	C	N1-C2-O2	5.20	122.02	118.90
1	2	466	G	C4-N9-C1'	5.18	133.23	126.50
1	2	750	C	N1-C2-O2	5.18	122.01	118.90
1	2	973	C	N1-C2-O2	5.18	122.01	118.90
4	D	83	LEU	CA-CB-CG	5.17	127.20	115.30
8	H	197	GLN	N-CA-CB	5.17	119.91	110.60
1	2	1556	A	C2-N3-C4	5.15	113.17	110.60
1	2	676	C	N1-C2-O2	5.14	121.98	118.90
1	2	1139	C	C5-C6-N1	5.13	123.57	121.00
5	E	109	LEU	CA-CB-CG	5.13	127.09	115.30
1	2	179	C	N3-C2-O2	-5.11	118.32	121.90
1	2	1700	C	C2-N1-C1'	5.11	124.42	118.80
1	2	790	C	N3-C2-O2	-5.09	118.33	121.90
1	2	1416	C	C5-C6-N1	5.09	123.55	121.00
1	2	1590	C	N1-C2-O2	5.09	121.95	118.90
1	2	1172	U	N1-C2-O2	5.09	126.36	122.80
1	2	1536	G	N3-C4-C5	-5.08	126.06	128.60
1	2	930	C	N1-C2-O2	5.07	121.94	118.90
1	2	961	G	C4-N9-C1'	5.07	133.09	126.50
36	z	160	U	P-O3'-C3'	5.06	125.78	119.70
1	2	196	C	N3-C2-O2	-5.06	118.36	121.90
21	U	131	LEU	CA-CB-CG	5.05	126.92	115.30
36	z	280	C	N3-C2-O2	-5.04	118.37	121.90
1	2	466	G	N3-C4-C5	-5.04	126.08	128.60

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	2	1389	C	C6-N1-C2	-5.02	118.29	120.30
31	e	56	ASP	CB-CG-OD1	5.01	122.81	118.30
1	2	578	C	C6-N1-C2	-5.01	118.30	120.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	C	191	ASP	Peptide
21	U	129	ARG	Sidechain

5.2 Too-close contacts (i)

Due to software issues we are unable to calculate clashes - this section is therefore empty.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	215/295 (73%)	202 (94%)	13 (6%)	0	100	100
3	C	211/264 (80%)	198 (94%)	12 (6%)	1 (0%)	29	68
4	D	219/221 (99%)	208 (95%)	11 (5%)	0	100	100
5	E	226/281 (80%)	214 (95%)	12 (5%)	0	100	100
6	F	260/263 (99%)	247 (95%)	13 (5%)	0	100	100
7	G	189/204 (93%)	178 (94%)	11 (6%)	0	100	100
8	H	235/249 (94%)	231 (98%)	4 (2%)	0	100	100
9	I	181/432 (42%)	173 (96%)	8 (4%)	0	100	100
10	J	204/208 (98%)	188 (92%)	16 (8%)	0	100	100
11	K	183/194 (94%)	179 (98%)	4 (2%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
12	L	94/149 (63%)	87 (93%)	7 (7%)	0	100	100
13	M	149/158 (94%)	132 (89%)	17 (11%)	0	100	100
14	N	115/132 (87%)	104 (90%)	11 (10%)	0	100	100
15	O	147/151 (97%)	144 (98%)	3 (2%)	0	100	100
16	P	134/168 (80%)	127 (95%)	7 (5%)	0	100	100
17	Q	113/145 (78%)	101 (89%)	12 (11%)	0	100	100
18	R	140/172 (81%)	132 (94%)	8 (6%)	0	100	100
19	S	130/135 (96%)	117 (90%)	13 (10%)	0	100	100
20	T	142/152 (93%)	135 (95%)	7 (5%)	0	100	100
21	U	139/145 (96%)	134 (96%)	5 (4%)	0	100	100
22	V	98/119 (82%)	95 (97%)	3 (3%)	0	100	100
23	W	81/83 (98%)	77 (95%)	4 (5%)	0	100	100
24	X	127/130 (98%)	121 (95%)	6 (5%)	0	100	100
25	Y	139/143 (97%)	133 (96%)	6 (4%)	0	100	100
26	Z	122/131 (93%)	116 (95%)	6 (5%)	0	100	100
27	a	75/124 (60%)	73 (97%)	2 (3%)	0	100	100
28	b	99/101 (98%)	94 (95%)	5 (5%)	0	100	100
29	c	81/84 (96%)	77 (95%)	4 (5%)	0	100	100
30	d	65/69 (94%)	63 (97%)	2 (3%)	0	100	100
31	e	53/56 (95%)	51 (96%)	2 (4%)	0	100	100
32	f	55/133 (41%)	51 (93%)	4 (7%)	0	100	100
33	g	66/188 (35%)	61 (92%)	5 (8%)	0	100	100
34	h	311/317 (98%)	277 (89%)	34 (11%)	0	100	100
35	n	23/25 (92%)	23 (100%)	0	0	100	100
All	All	4821/5821 (83%)	4543 (94%)	277 (6%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	C	190	PRO

5.3.2 Protein sidechains (i)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
2	B	180/245 (74%)	179 (99%)	1 (1%)	86 92
3	C	194/231 (84%)	194 (100%)	0	100 100
4	D	187/187 (100%)	187 (100%)	0	100 100
5	E	190/232 (82%)	189 (100%)	1 (0%)	88 93
6	F	224/225 (100%)	224 (100%)	0	100 100
7	G	161/170 (95%)	161 (100%)	0	100 100
8	H	207/218 (95%)	206 (100%)	1 (0%)	88 93
9	I	165/360 (46%)	165 (100%)	0	100 100
10	J	178/180 (99%)	176 (99%)	2 (1%)	73 85
11	K	161/168 (96%)	160 (99%)	1 (1%)	86 92
12	L	87/125 (70%)	85 (98%)	2 (2%)	50 70
13	M	136/142 (96%)	135 (99%)	1 (1%)	84 90
14	N	99/108 (92%)	99 (100%)	0	100 100
15	O	130/131 (99%)	130 (100%)	0	100 100
16	P	106/130 (82%)	106 (100%)	0	100 100
17	Q	105/130 (81%)	105 (100%)	0	100 100
18	R	117/140 (84%)	116 (99%)	1 (1%)	78 87
19	S	119/121 (98%)	119 (100%)	0	100 100
20	T	125/132 (95%)	124 (99%)	1 (1%)	81 89
21	U	111/116 (96%)	111 (100%)	0	100 100
22	V	92/107 (86%)	92 (100%)	0	100 100
23	W	67/67 (100%)	67 (100%)	0	100 100
24	X	112/113 (99%)	110 (98%)	2 (2%)	59 77
25	Y	113/115 (98%)	112 (99%)	1 (1%)	78 87
26	Z	107/113 (95%)	107 (100%)	0	100 100
27	a	68/102 (67%)	68 (100%)	0	100 100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
28	b	88/88 (100%)	88 (100%)	0	100 100
29	c	75/76 (99%)	74 (99%)	1 (1%)	69 82
30	d	60/62 (97%)	60 (100%)	0	100 100
31	e	48/49 (98%)	48 (100%)	0	100 100
32	f	47/106 (44%)	47 (100%)	0	100 100
33	g	61/154 (40%)	61 (100%)	0	100 100
34	h	272/275 (99%)	272 (100%)	0	100 100
35	n	24/24 (100%)	24 (100%)	0	100 100
All	All	4216/4942 (85%)	4201 (100%)	15 (0%)	91 94

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	84	GLN
5	E	76	ARG
8	H	98	ARG
10	J	59	ARG
10	J	143	LYS
11	K	5	ARG
12	L	6	LYS
12	L	38	LYS
13	M	69	ARG
18	R	146	ARG
20	T	142	ARG
24	X	12	LYS
24	X	15	ASN
25	Y	140	ARG
29	c	17	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	33	GLN
12	L	7	ASN
24	X	15	ASN
24	X	24	GLN
27	a	106	GLN

5.3.3 RNA (i)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	2	1685/1870 (90%)	388 (23%)	7 (0%)
36	z	160/400 (40%)	47 (29%)	0
All	All	1845/2270 (81%)	435 (23%)	7 (0%)

All (435) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	2	4	C
1	2	8	U
1	2	9	U
1	2	26	U
1	2	33	G
1	2	41	G
1	2	42	A
1	2	44	U
1	2	45	A
1	2	46	A
1	2	62	G
1	2	67	C
1	2	68	A
1	2	73	C
1	2	74	G
1	2	88	G
1	2	92	A
1	2	99	A
1	2	103	A
1	2	110	U
1	2	113	G
1	2	114	G
1	2	115	U
1	2	126	G
1	2	130	G
1	2	142	C
1	2	143	U
1	2	146	G
1	2	147	A
1	2	155	G
1	2	158	A
1	2	162	C
1	2	170	A
1	2	182	C

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	2	183	G
1	2	184	G
1	2	186	C
1	2	192	C
1	2	199	C
1	2	292	A
1	2	294	U
1	2	295	C
1	2	306	C
1	2	307	G
1	2	309	G
1	2	312	G
1	2	319	C
1	2	330	G
1	2	331	C
1	2	343	A
1	2	347	G
1	2	351	G
1	2	362	C
1	2	364	A
1	2	368	U
1	2	370	G
1	2	371	A
1	2	383	G
1	2	384	U
1	2	385	G
1	2	392	A
1	2	393	U
1	2	395	G
1	2	398	A
1	2	400	C
1	2	407	G
1	2	408	A
1	2	409	C
1	2	418	A
1	2	426	A
1	2	429	C
1	2	435	A
1	2	436	G
1	2	438	G
1	2	446	G
1	2	448	A

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	2	449	A
1	2	450	C
1	2	452	G
1	2	465	A
1	2	471	G
1	2	472	C
1	2	474	G
1	2	486	A
1	2	487	U
1	2	488	U
1	2	489	A
1	2	492	C
1	2	493	A
1	2	500	A
1	2	502	C
1	2	509	G
1	2	523	A
1	2	525	A
1	2	531	A
1	2	532	C
1	2	533	A
1	2	536	A
1	2	537	C
1	2	541	U
1	2	542	U
1	2	548	C
1	2	550	C
1	2	551	U
1	2	554	A
1	2	556	U
1	2	558	G
1	2	561	A
1	2	562	U
1	2	576	A
1	2	583	A
1	2	588	G
1	2	590	A
1	2	591	U
1	2	593	C
1	2	603	C
1	2	606	G
1	2	608	C

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	2	613	G
1	2	614	C
1	2	625	G
1	2	627	U
1	2	628	A
1	2	629	A
1	2	642	U
1	2	643	A
1	2	644	G
1	2	650	A
1	2	655	A
1	2	656	G
1	2	660	C
1	2	668	A
1	2	669	A
1	2	672	A
1	2	673	G
1	2	684	G
1	2	689	U
1	2	733	C
1	2	752	G
1	2	753	C
1	2	754	G
1	2	790	C
1	2	793	G
1	2	794	A
1	2	798	G
1	2	799	U
1	2	810	A
1	2	811	A
1	2	821	G
1	2	822	U
1	2	830	A
1	2	831	G
1	2	844	U
1	2	847	A
1	2	860	G
1	2	864	A
1	2	868	G
1	2	870	A
1	2	872	A
1	2	874	G

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	2	875	A
1	2	884	C
1	2	887	U
1	2	890	U
1	2	891	G
1	2	892	U
1	2	893	U
1	2	895	G
1	2	898	U
1	2	901	G
1	2	913	A
1	2	914	U
1	2	919	A
1	2	920	A
1	2	922	A
1	2	933	G
1	2	955	A
1	2	962	A
1	2	963	A
1	2	970	G
1	2	978	G
1	2	985	G
1	2	988	C
1	2	990	A
1	2	992	A
1	2	1002	U
1	2	1017	U
1	2	1021	U
1	2	1022	U
1	2	1023	A
1	2	1026	C
1	2	1027	A
1	2	1041	G
1	2	1042	A
1	2	1044	G
1	2	1055	A
1	2	1061	U
1	2	1062	A
1	2	1078	C
1	2	1083	A
1	2	1085	C
1	2	1086	G

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	2	1088	U
1	2	1089	G
1	2	1097	G
1	2	1113	A
1	2	1114	U
1	2	1118	C
1	2	1121	G
1	2	1129	G
1	2	1133	A
1	2	1138	C
1	2	1139	C
1	2	1149	A
1	2	1150	A
1	2	1153	C
1	2	1154	U
1	2	1155	U
1	2	1161	U
1	2	1195	A
1	2	1203	U
1	2	1205	G
1	2	1206	C
1	2	1207	G
1	2	1210	G
1	2	1215	C
1	2	1216	C
1	2	1217	A
1	2	1221	G
1	2	1224	G
1	2	1233	G
1	2	1237	C
1	2	1242	U
1	2	1243	U
1	2	1251	A
1	2	1253	A
1	2	1256	G
1	2	1257	G
1	2	1258	A
1	2	1259	A
1	2	1264	C
1	2	1266	C
1	2	1272	C
1	2	1274	G

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	2	1275	G
1	2	1282	A
1	2	1284	A
1	2	1285	G
1	2	1286	G
1	2	1298	G
1	2	1299	A
1	2	1301	A
1	2	1302	G
1	2	1303	C
1	2	1307	U
1	2	1310	U
1	2	1311	C
1	2	1312	G
1	2	1313	A
1	2	1315	U
1	2	1316	C
1	2	1318	G
1	2	1327	G
1	2	1330	G
1	2	1333	U
1	2	1342	U
1	2	1358	U
1	2	1371	U
1	2	1372	U
1	2	1378	A
1	2	1390	U
1	2	1395	C
1	2	1396	A
1	2	1397	U
1	2	1398	G
1	2	1404	U
1	2	1406	G
1	2	1416	C
1	2	1428	G
1	2	1439	A
1	2	1442	U
1	2	1452	A
1	2	1454	A
1	2	1462	U
1	2	1463	U
1	2	1464	C

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	2	1466	G
1	2	1476	A
1	2	1477	U
1	2	1487	A
1	2	1489	A
1	2	1490	G
1	2	1494	U
1	2	1495	G
1	2	1498	A
1	2	1507	G
1	2	1508	A
1	2	1509	U
1	2	1520	G
1	2	1522	A
1	2	1533	A
1	2	1535	U
1	2	1536	G
1	2	1548	G
1	2	1551	U
1	2	1552	G
1	2	1554	C
1	2	1556	A
1	2	1558	C
1	2	1560	U
1	2	1567	G
1	2	1568	C
1	2	1575	G
1	2	1578	U
1	2	1580	A
1	2	1582	C
1	2	1585	U
1	2	1588	A
1	2	1591	C
1	2	1595	U
1	2	1598	G
1	2	1599	U
1	2	1600	G
1	2	1601	A
1	2	1603	G
1	2	1604	G
1	2	1606	G
1	2	1621	U

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	2	1623	A
1	2	1632	G
1	2	1636	G
1	2	1637	A
1	2	1638	G
1	2	1646	C
1	2	1648	G
1	2	1649	U
1	2	1664	A
1	2	1665	G
1	2	1677	U
1	2	1678	A
1	2	1680	G
1	2	1682	C
1	2	1683	C
1	2	1697	A
1	2	1698	C
1	2	1699	A
1	2	1700	C
1	2	1701	C
1	2	1704	C
1	2	1715	A
1	2	1721	U
1	2	1722	G
1	2	1726	G
1	2	1751	C
1	2	1756	C
1	2	1757	G
1	2	1776	G
1	2	1779	G
1	2	1781	A
1	2	1782	G
1	2	1783	C
1	2	1799	G
1	2	1801	A
1	2	1813	A
1	2	1814	G
1	2	1823	A
1	2	1824	A
1	2	1825	A
1	2	1829	G
1	2	1831	A

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	2	1834	A
1	2	1835	A
1	2	1836	G
1	2	1838	U
1	2	1839	U
1	2	1846	G
1	2	1849	G
1	2	1851	A
1	2	1852	C
1	2	1860	A
1	2	1861	G
1	2	1862	G
1	2	1863	A
1	2	1864	U
1	2	1865	C
1	2	1866	A
1	2	1868	U
1	2	1869	A
36	z	121	C
36	z	124	C
36	z	133	G
36	z	134	A
36	z	135	G
36	z	136	A
36	z	137	G
36	z	144	U
36	z	146	G
36	z	152	G
36	z	154	A
36	z	157	C
36	z	161	G
36	z	162	A
36	z	164	U
36	z	227	U
36	z	229	G
36	z	230	G
36	z	232	C
36	z	234	U
36	z	237	C
36	z	243	A
36	z	244	A
36	z	245	G

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
36	z	252	A
36	z	258	G
36	z	264	U
36	z	265	U
36	z	266	G
36	z	269	U
36	z	273	G
36	z	282	U
36	z	286	G
36	z	289	C
36	z	290	U
36	z	291	G
36	z	293	C
36	z	294	U
36	z	295	G
36	z	296	A
36	z	298	A
36	z	301	G
36	z	303	G
36	z	306	U
36	z	311	G
36	z	329	U
36	z	330	A

All (7) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	2	553	U
1	2	561	A
1	2	627	U
1	2	688	U
1	2	752	G
1	2	1137	U
1	2	1664	A

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

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

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 2 ligands modelled in this entry, 2 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.

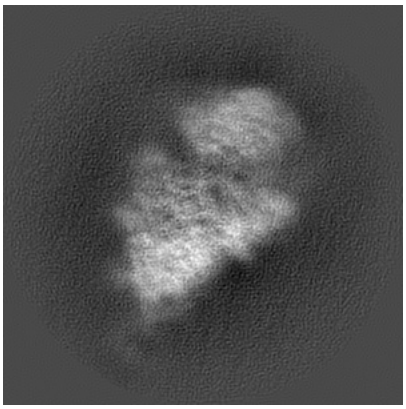
6 Map visualisation

This section contains visualisations of the EMDB entry EMD-25528. These allow visual inspection of the internal detail of the map and identification of artifacts.

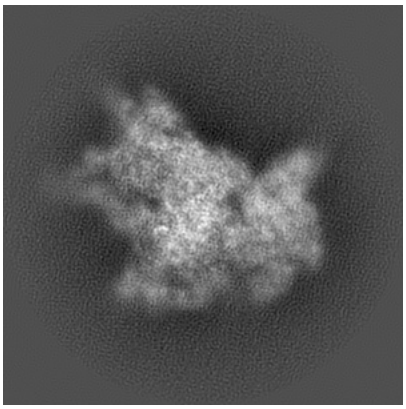
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections

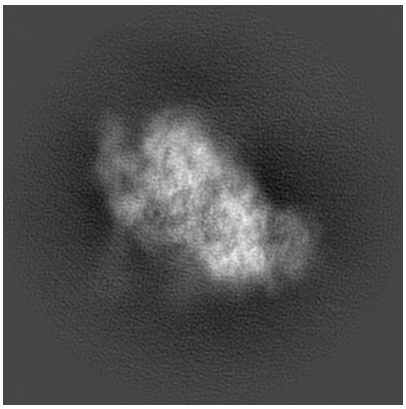
6.1.1 Primary map



X

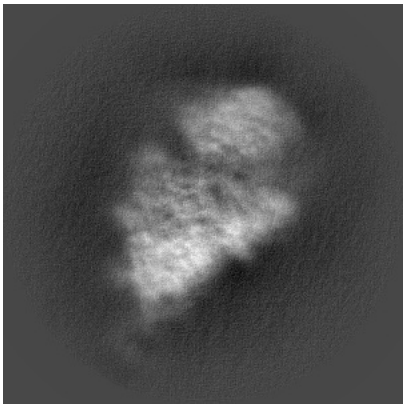


Y

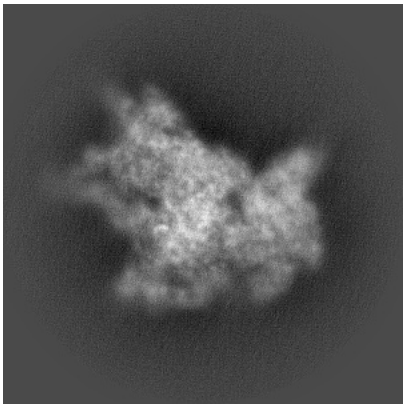


Z

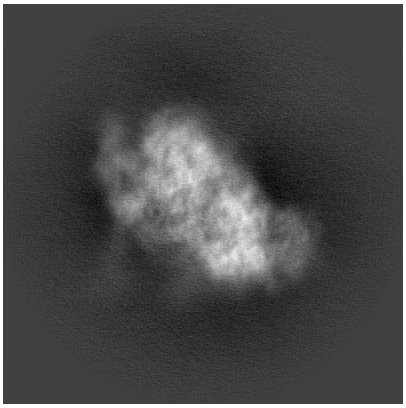
6.1.2 Raw map



X



Y

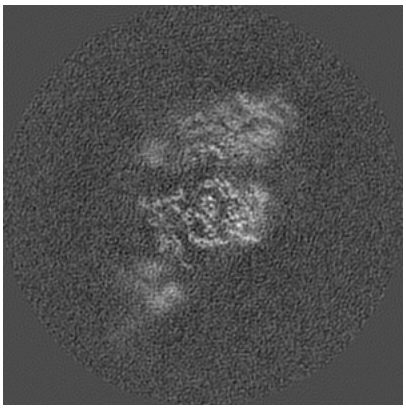


Z

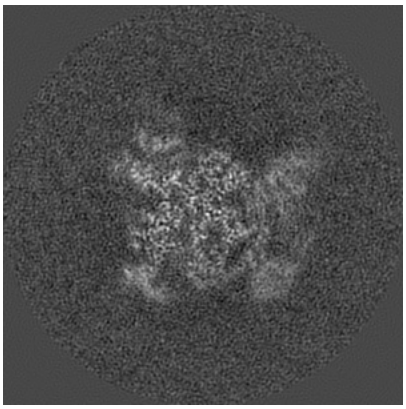
The images above show the map projected in three orthogonal directions.

6.2 Central slices (i)

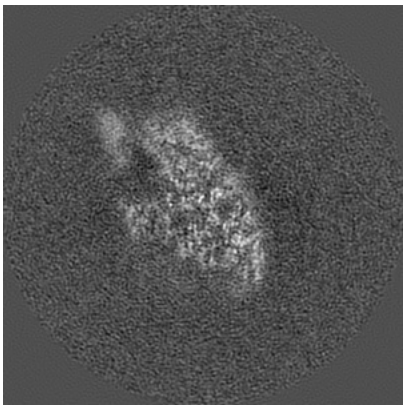
6.2.1 Primary map



X Index: 200

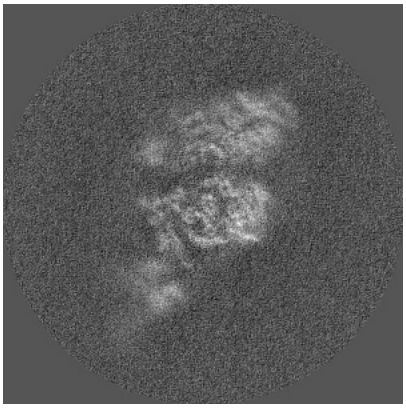


Y Index: 200

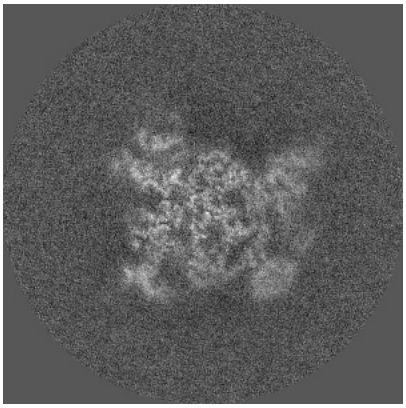


Z Index: 200

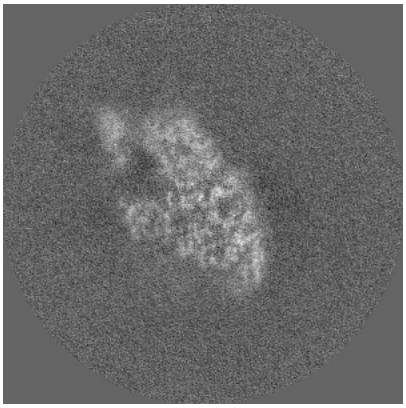
6.2.2 Raw map



X Index: 200



Y Index: 200

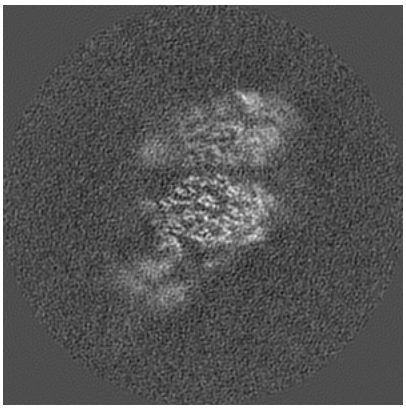


Z Index: 200

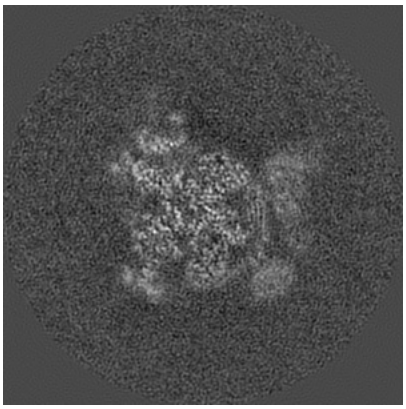
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices (i)

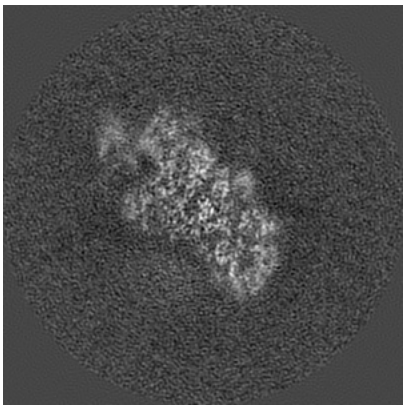
6.3.1 Primary map



X Index: 196

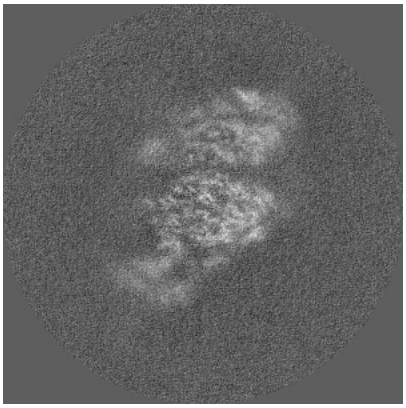


Y Index: 197

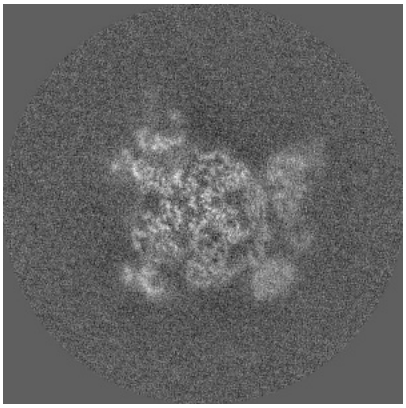


Z Index: 188

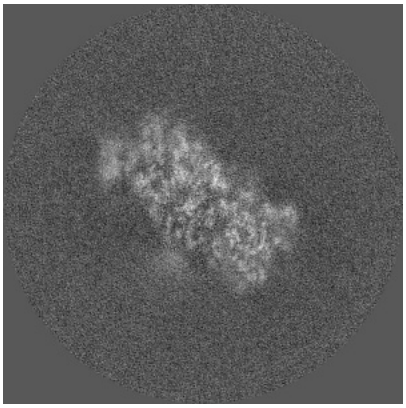
6.3.2 Raw map



X Index: 195



Y Index: 198

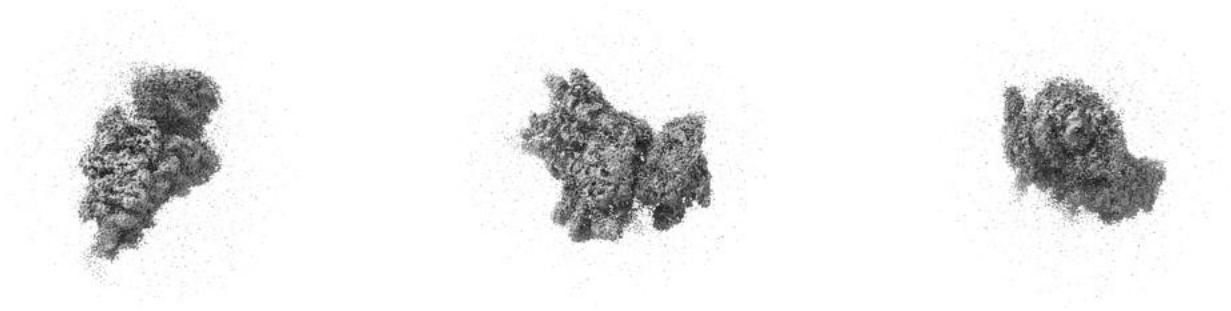


Z Index: 176

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

6.4 Orthogonal surface views (i)

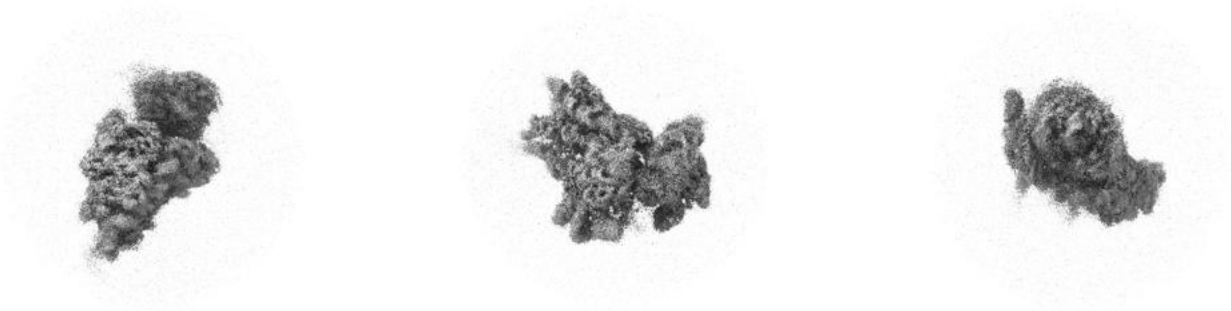
6.4.1 Primary map



X Y Z

The images above show the 3D surface view of the map at the recommended contour level 0.00883. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.4.2 Raw map



X Y Z

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.

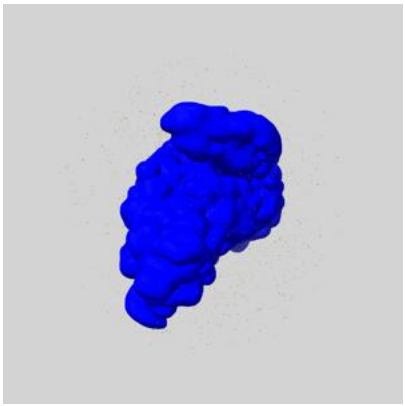
6.5 Mask visualisation (i)

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

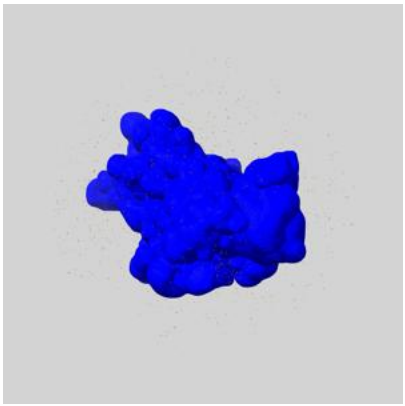
A mask typically either:

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

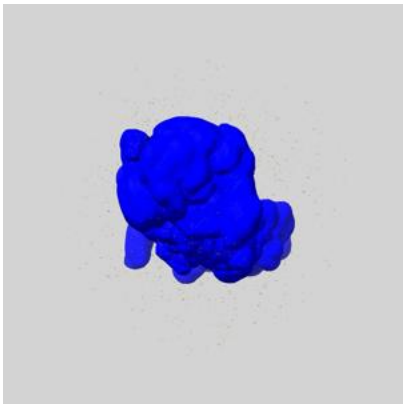
6.5.1 emd_25528_msk_1.map (i)



X



Y

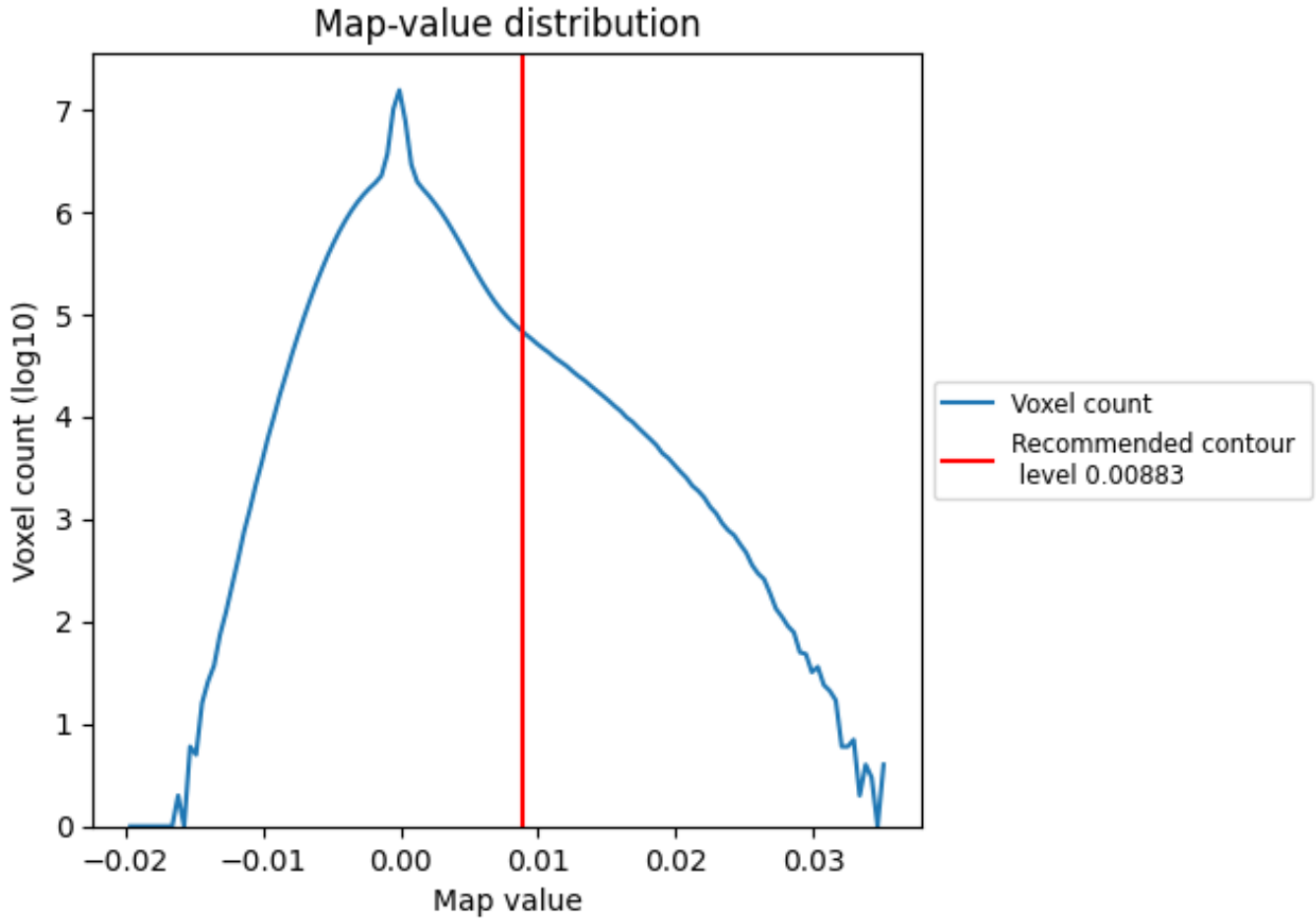


Z

7 Map analysis

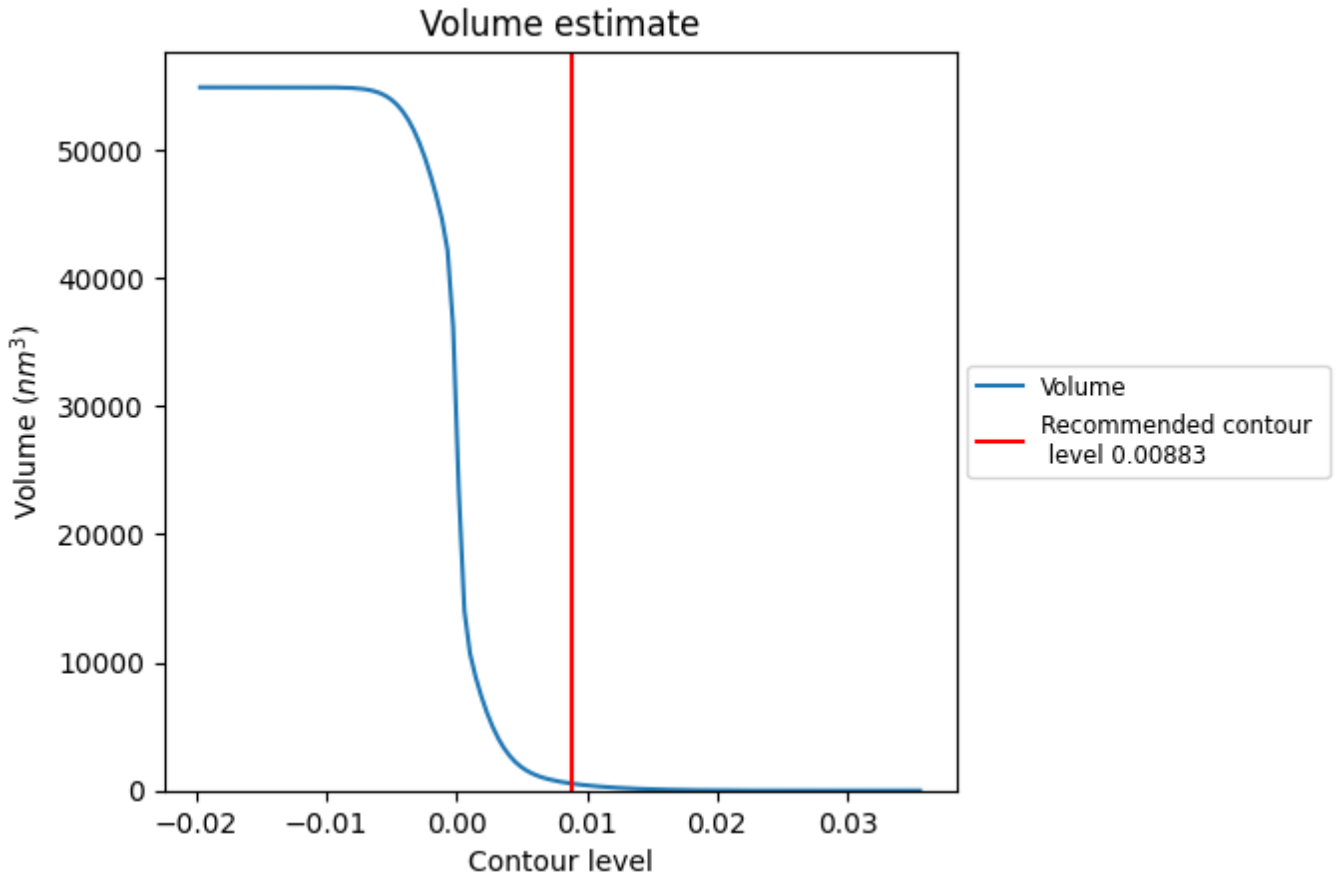
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution



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.

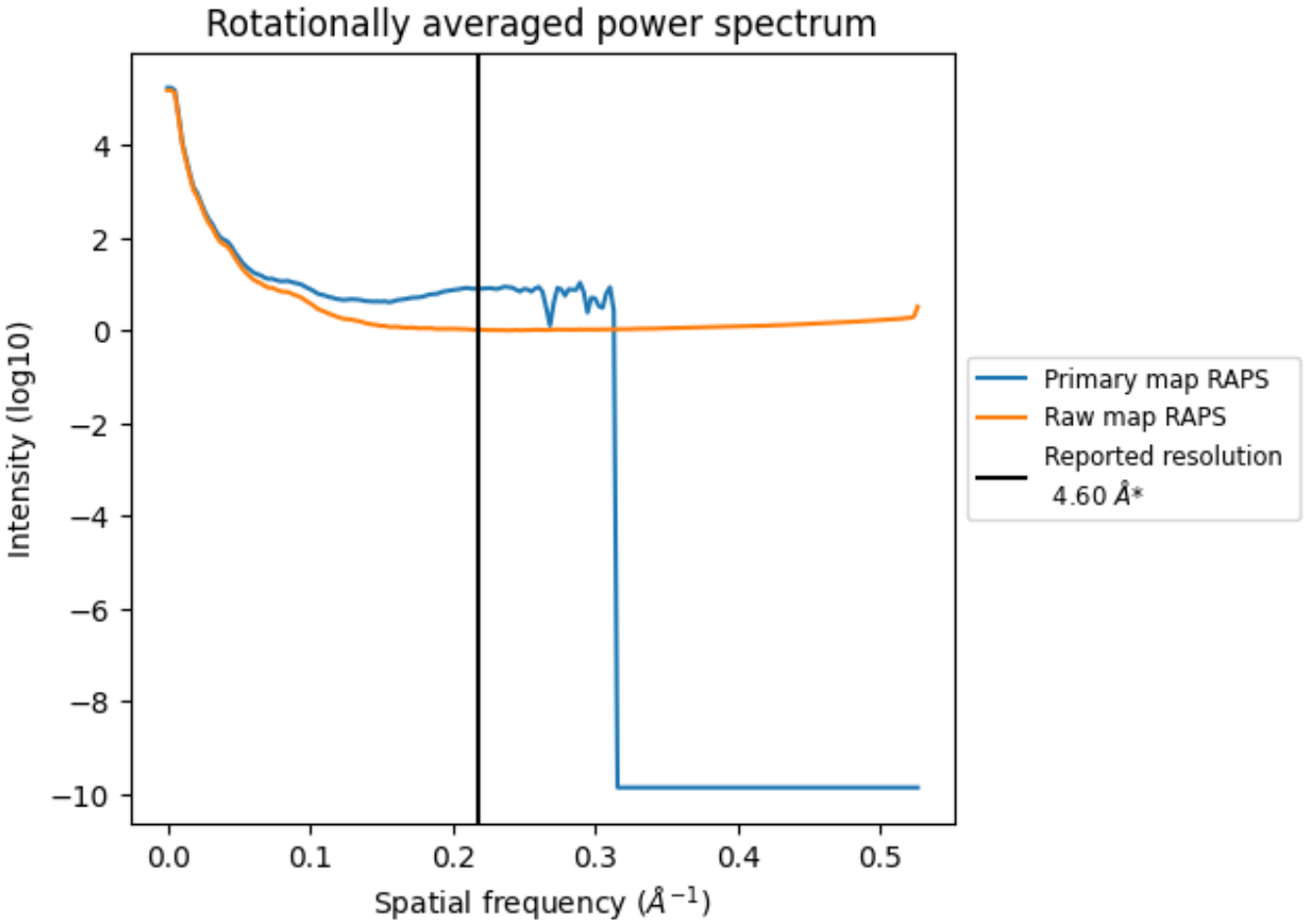
7.2 Volume estimate



The volume at the recommended contour level is 549 nm^3; this corresponds to an approximate mass of 496 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 (i)

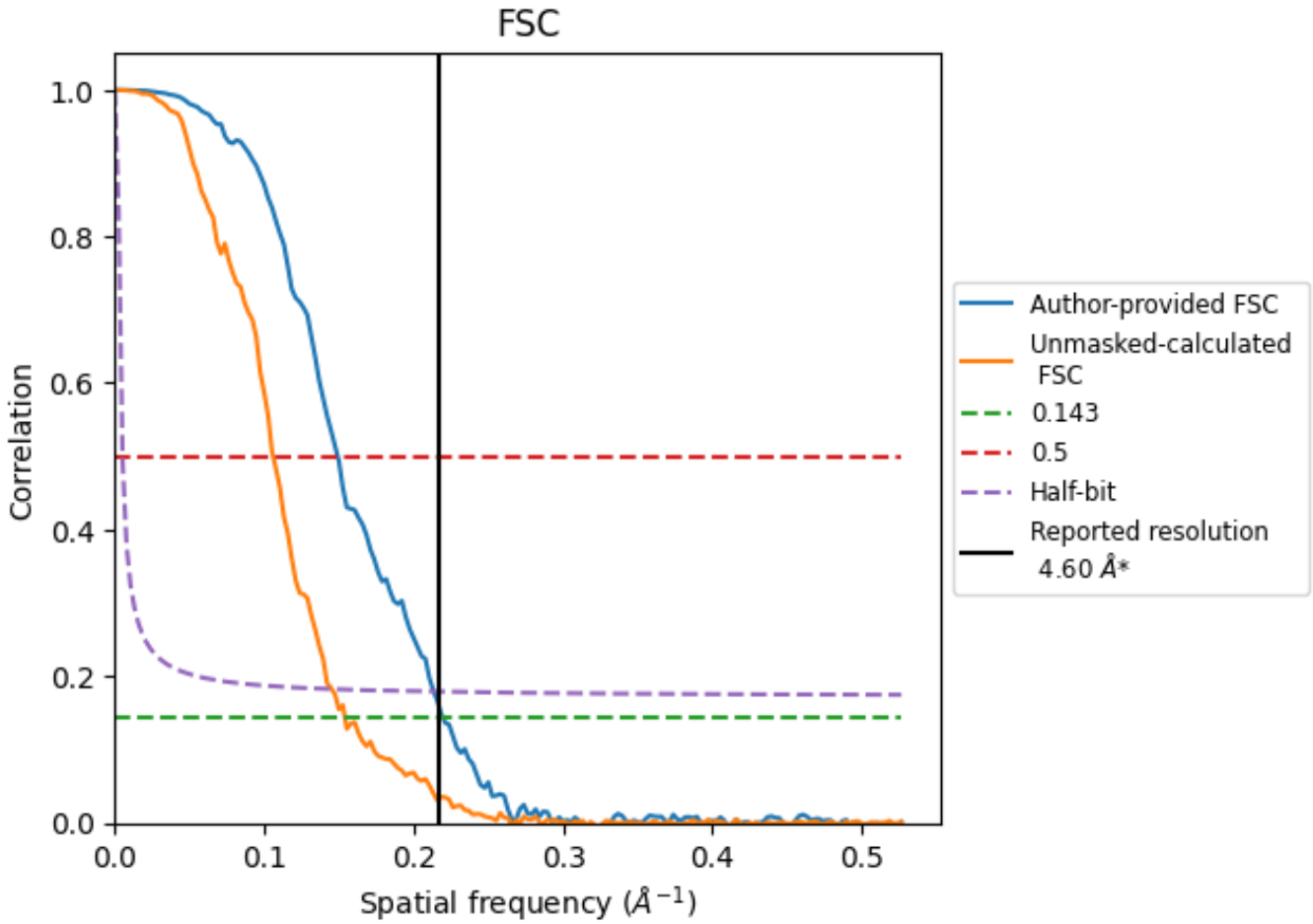


*Reported resolution corresponds to spatial frequency of 0.217 Å^{-1}

8 Fourier-Shell correlation

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



*Reported resolution corresponds to spatial frequency of 0.217 $Å^{-1}$

8.2 Resolution estimates (i)

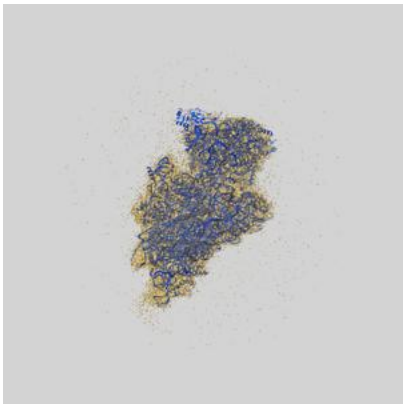
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	4.60	-	-
Author-provided FSC curve	4.55	6.71	4.69
Unmasked-calculated*	6.49	9.44	6.89

*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 6.49 differs from the reported value 4.6 by more than 10 %

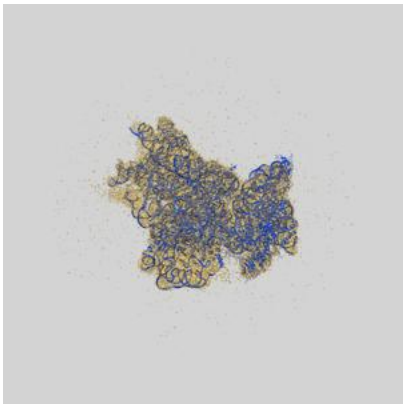
9 Map-model fit

This section contains information regarding the fit between EMDB map EMD-25528 and PDB model 7SYH. Per-residue inclusion information can be found in section 3 on page 12.

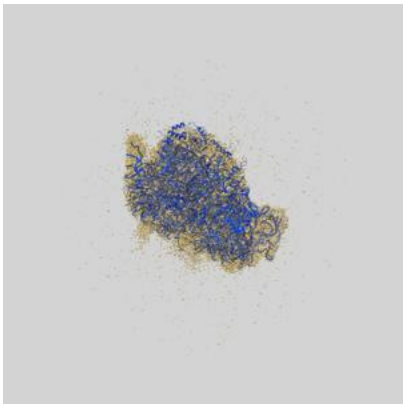
9.1 Map-model overlay



X



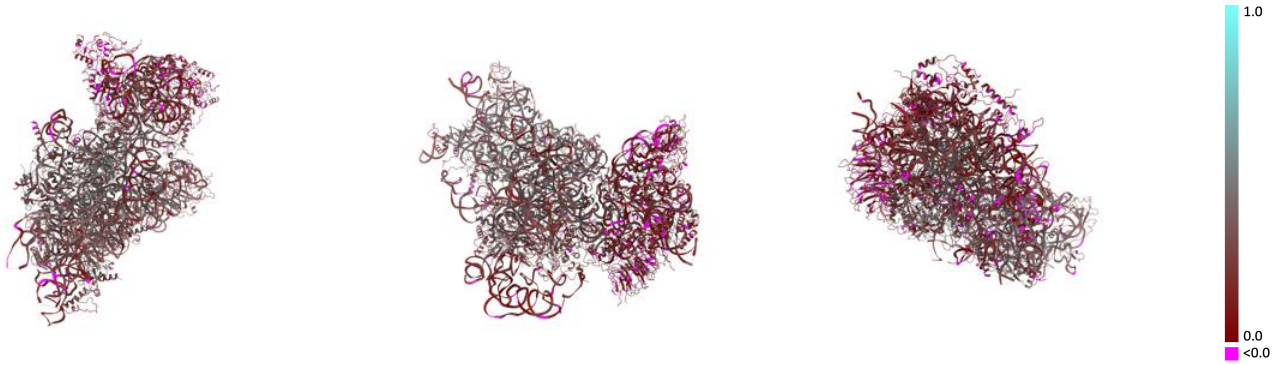
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.00883 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.

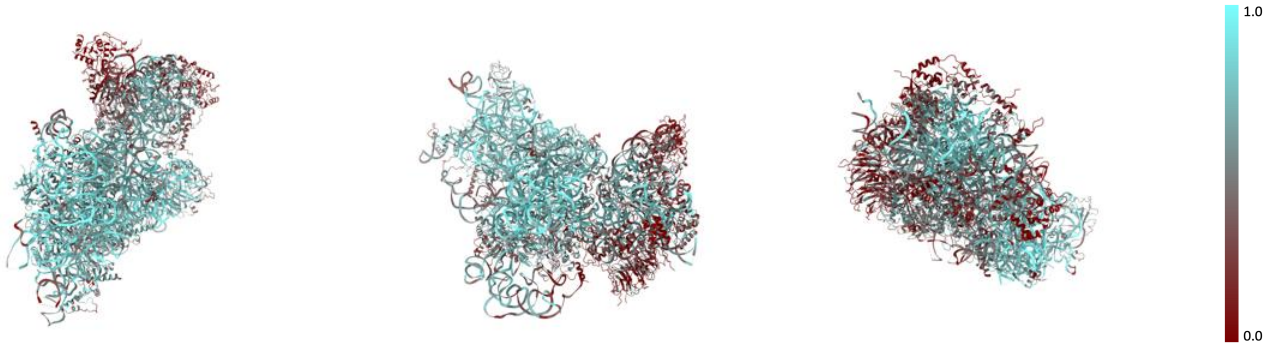
9.2 Q-score mapped to coordinate model (i)



X Y Z

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.

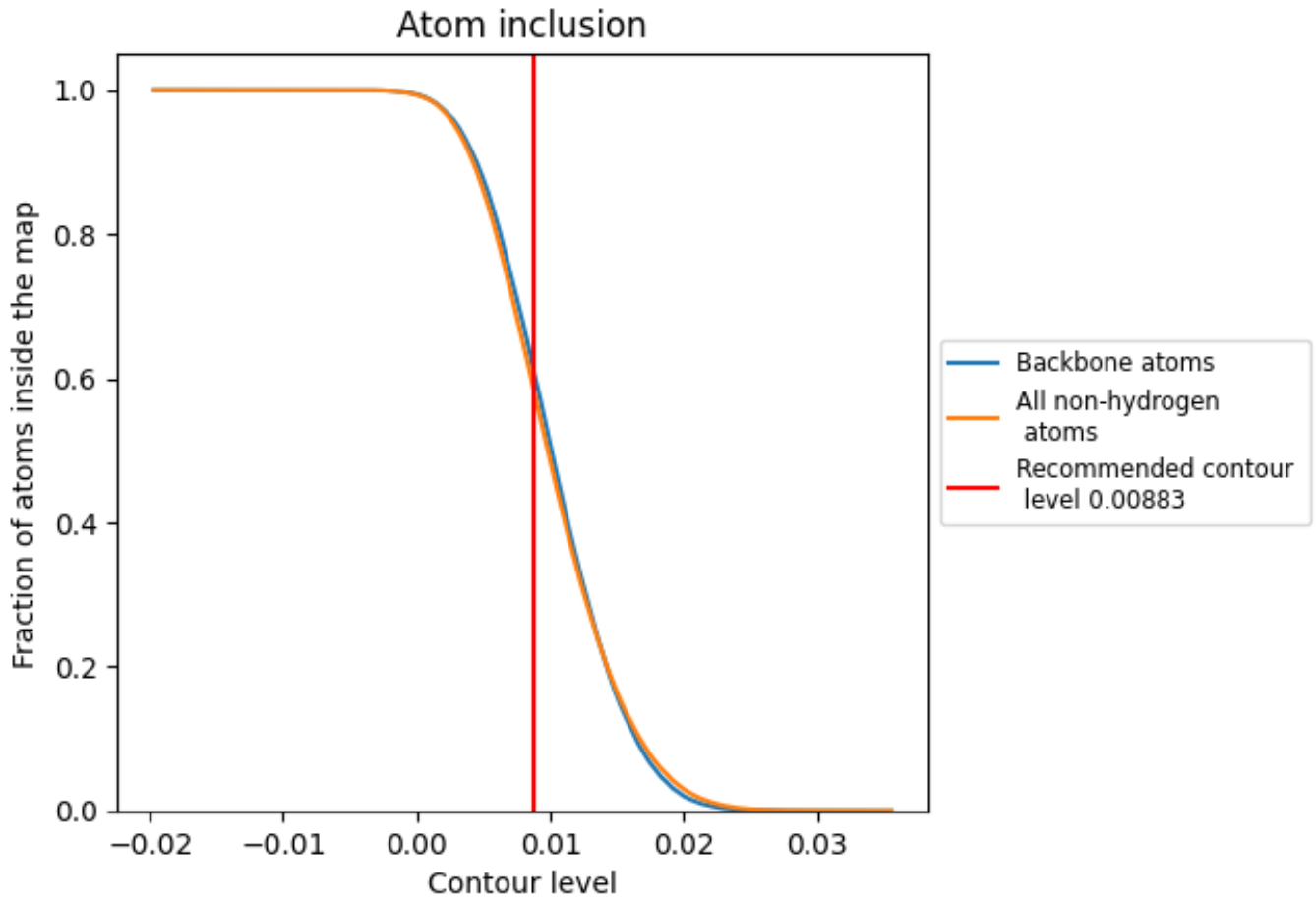
9.3 Atom inclusion mapped to coordinate model (i)



X Y Z

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 (0.00883).

9.4 Atom inclusion (i)



At the recommended contour level, 60% of all backbone atoms, 58% 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 (0.00883) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.5782	0.2710
2	0.7353	0.2910
B	0.5135	0.3310
C	0.5874	0.2940
D	0.5232	0.3480
E	0.2516	0.2010
F	0.6365	0.3670
G	0.3045	0.1740
H	0.5763	0.2820
I	0.3546	0.2820
J	0.6613	0.3300
K	0.6109	0.3480
L	0.2759	0.1920
M	0.5810	0.3810
N	0.0257	0.1030
O	0.6278	0.3440
P	0.6059	0.2760
Q	0.3045	0.1310
R	0.3251	0.1950
S	0.2341	0.2080
T	0.2131	0.1270
U	0.4283	0.2010
V	0.1873	0.2190
W	0.5016	0.3370
X	0.5574	0.3800
Y	0.6153	0.3640
Z	0.6799	0.3380
a	0.0835	0.1250
b	0.5537	0.3420
c	0.4961	0.3430
d	0.2283	0.2090
e	0.3696	0.2140
f	0.4045	0.1970
g	0.0536	0.1190
h	0.1878	0.1290

Continued on next page...

Continued from previous page...

Chain	Atom inclusion	Q-score
n	0.5917	0.2990
z	0.5082	0.1460