



wwPDB EM Validation Summary Report ⓘ

Jun 18, 2026 – 11:11 pm BST

PDB ID : 6ZXG / pdb_00006zxg
EMDB ID : EMD-11520
Title : Cryo-EM structure of a late human pre-40S ribosomal subunit - State H1
Authors : Ameismeier, M.; Zemp, I.; van den Heuvel, J.; Thoms, M.; Berninghausen, O.;
Kutay, U.; Beckmann, R.
Deposited on : 2020-07-29
Resolution : 2.60 Å(reported)

This is a wwPDB EM Validation Summary 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:

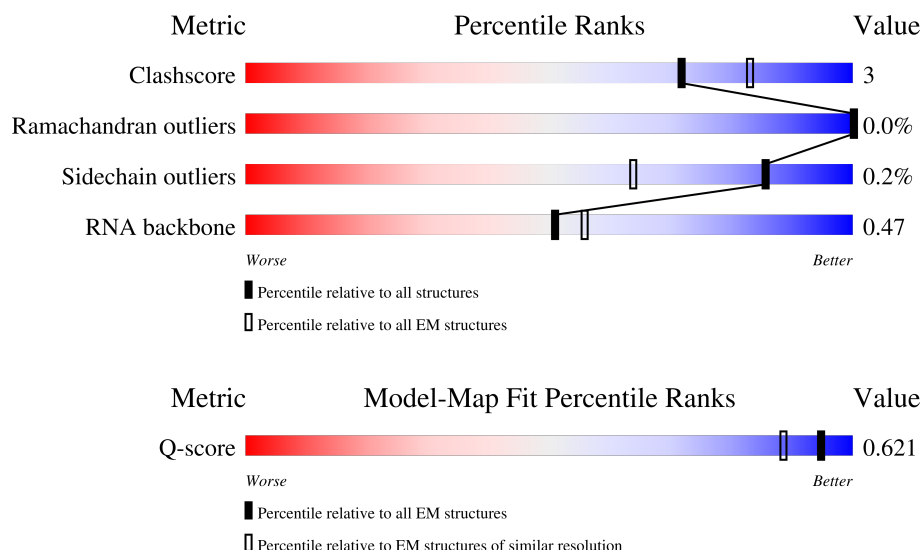
EMDB validation analysis : 0.0.1.dev132
Mogul : 1.8.4, CSD as541be (2020)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

1 Overall quality at a glance

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

The reported resolution of this entry is 2.60 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)	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	8728 (2.10 - 3.10)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion $< 40\%$). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	2	1869	
2	A	295	
3	B	264	

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
4	C	293	
5	h	115	
6	E	263	
7	D	243	
8	G	249	
9	H	194	
10	I	208	
11	J	194	
12	F	204	
13	L	158	
14	K	165	
15	N	151	
16	O	151	
17	M	132	
18	P	145	
19	R	135	
20	Q	146	
21	S	152	
22	T	145	
23	V	83	
24	W	130	
25	X	143	
26	Y	133	
27	U	119	
28	Z	125	

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
29	b	84	
30	c	69	
31	d	56	
32	e	59	
33	f	156	
34	g	317	
35	j	165	
36	k	583	
37	z	568	

2 Entry composition

There are 41 unique types of molecules in this entry. The entry contains 79865 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 pre-18S ribosomal RNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	2	1651	Total	C	N	O	P	0	0
			35310	15787	6344	11528	1651		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
2	A	216	Total	C	N	O	S	0	0
			1705	1083	299	315	8		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
3	B	211	Total	C	N	O	S	0	0
			1715	1088	307	306	14		

- Molecule 4 is a protein called 40S ribosomal protein S2.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	C	216	Total	C	N	O	S	0	0
			1674	1085	287	292	10		

- Molecule 5 is a protein called 40S ribosomal protein S26.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	h	99	Total	C	N	O	S	0	0
			792	492	165	130	5		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
6	E	255	Total	C	N	O	S	0	0
			2031	1299	377	347	8		

- Molecule 7 is a protein called 40S ribosomal protein S3.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	D	224	Total	C	N	O	S	0	0
			1745	1112	314	312	7		

- Molecule 8 is a protein called 40S ribosomal protein S6.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	G	223	Total	C	N	O	S	0	0
			1802	1128	358	309	7		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
9	H	179	Total	C	N	O	S	0	0
			1446	927	264	254	1		

- Molecule 10 is a protein called 40S ribosomal protein S8.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	I	199	Total	C	N	O	S	0	0
			1638	1027	322	284	5		

- Molecule 11 is a protein called 40S ribosomal protein S9.

Mol	Chain	Residues	Atoms					AltConf	Trace
11	J	176	Total	C	N	O	S	0	0
			1465	934	295	234	2		

- Molecule 12 is a protein called 40S ribosomal protein S5.

Mol	Chain	Residues	Atoms					AltConf	Trace
12	F	184	Total	C	N	O	S	0	0
			1461	914	276	264	7		

- Molecule 13 is a protein called 40S ribosomal protein S11.

Mol	Chain	Residues	Atoms					AltConf	Trace
13	L	136	Total	C	N	O	S	0	0
			1127	719	212	190	6		

- Molecule 14 is a protein called 40S ribosomal protein S10.

Mol	Chain	Residues	Atoms					AltConf	Trace
14	K	95	Total	C	N	O	S	0	0
			799	524	139	130	6		

- Molecule 15 is a protein called 40S ribosomal protein S13.

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

- Molecule 16 is a protein called 40S ribosomal protein S14.

Mol	Chain	Residues	Atoms					AltConf	Trace
16	O	126	Total	C	N	O	S	0	0
			947	580	188	173	6		

- Molecule 17 is a protein called 40S ribosomal protein S12.

Mol	Chain	Residues	Atoms					AltConf	Trace
17	M	108	Total	C	N	O	S	0	0
			837	530	147	153	7		

- Molecule 18 is a protein called 40S ribosomal protein S15.

Mol	Chain	Residues	Atoms					AltConf	Trace
18	P	132	Total	C	N	O	S	0	0
			1077	685	202	183	7		

- Molecule 19 is a protein called 40S ribosomal protein S17.

Mol	Chain	Residues	Atoms					AltConf	Trace
19	R	132	Total	C	N	O	S	0	0
			1066	669	199	194	4		

- Molecule 20 is a protein called 40S ribosomal protein S16.

Mol	Chain	Residues	Atoms					AltConf	Trace
20	Q	140	Total	C	N	O	S	0	0
			1116	710	211	192	3		

- Molecule 21 is a protein called 40S ribosomal protein S18.

Mol	Chain	Residues	Atoms					AltConf	Trace
21	S	141	Total	C	N	O	S	0	0
			1162	731	232	198	1		

- Molecule 22 is a protein called 40S ribosomal protein S19.

Mol	Chain	Residues	Atoms					AltConf	Trace
22	T	141	Total	C	N	O	S	0	0
			1094	685	210	196	3		

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

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

- Molecule 24 is a protein called 40S ribosomal protein S15a.

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

- Molecule 25 is a protein called 40S ribosomal protein S23.

Mol	Chain	Residues	Atoms					AltConf	Trace
25	X	139	Total	C	N	O	S	0	0
			1080	682	214	181	3		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
26	Y	122	Total	C	N	O	S	0	0
			987	624	193	165	5		

- Molecule 27 is a protein called 40S ribosomal protein S20.

Mol	Chain	Residues	Atoms					AltConf	Trace
27	U	98	Total	C	N	O	S	0	0
			774	486	145	139	4		

- Molecule 28 is a protein called 40S ribosomal protein S25.

Mol	Chain	Residues	Atoms					AltConf	Trace
28	Z	82	Total	C	N	O	S	0	0
			638	410	117	110	1		

- Molecule 29 is a protein called 40S ribosomal protein S27.

Mol	Chain	Residues	Atoms					AltConf	Trace
29	b	77	Total	C	N	O	S	0	0
			611	386	113	107	5		

- Molecule 30 is a protein called 40S ribosomal protein S28.

Mol	Chain	Residues	Atoms					AltConf	Trace
30	c	61	Total	C	N	O	S	0	0
			479	292	95	90	2		

- Molecule 31 is a protein called 40S ribosomal protein S29.

Mol	Chain	Residues	Atoms					AltConf	Trace
31	d	54	Total	C	N	O	S	0	0
			450	282	93	70	5		

- Molecule 32 is a protein called 40S ribosomal protein S30.

Mol	Chain	Residues	Atoms					AltConf	Trace
32	e	50	Total	C	N	O	S	0	0
			398	244	90	63	1		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
33	f	57	Total	C	N	O	S	0	0
			465	295	89	74	7		

- Molecule 34 is a protein called Receptor of activated protein C kinase 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
34	g	295	Total	C	N	O	S	0	0
			2306	1460	403	431	12		

- Molecule 35 is a protein called Probable RNA-binding protein EIF1AD.

Mol	Chain	Residues	Atoms					AltConf	Trace
35	j	116	Total	C	N	O	S	0	0
			934	593	167	172	2		

- Molecule 36 is a protein called Leucine-rich repeat-containing protein 47.

Mol	Chain	Residues	Atoms					AltConf	Trace
36	k	430	Total	C	N	O	S	0	0
			2886	1788	555	536	7		

- Molecule 37 is a protein called Serine/threonine-protein kinase RIO1.

Mol	Chain	Residues	Atoms					AltConf	Trace
37	z	349	Total	C	N	O	S	0	0
			2803	1757	515	511	20		

- Molecule 38 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		AltConf
38	2	123	Total	Mg	0
			123	123	
38	h	1	Total	Mg	0
			1	1	
38	G	1	Total	Mg	0
			1	1	
38	F	1	Total	Mg	0
			1	1	
38	O	1	Total	Mg	0
			1	1	
38	X	1	Total	Mg	0
			1	1	
38	d	1	Total	Mg	0
			1	1	
38	z	2	Total	Mg	0
			2	2	

- Molecule 39 is ZINC ION (CCD ID: ZN) (formula: Zn).

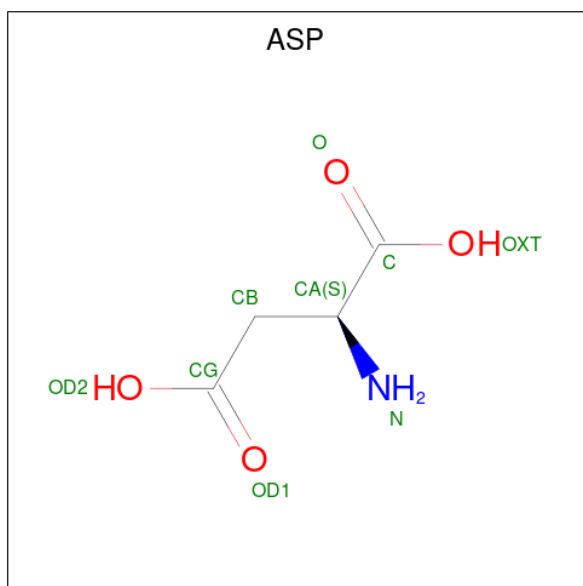
Mol	Chain	Residues	Atoms		AltConf
39	h	1	Total	Zn	0
			1	1	
39	d	1	Total	Zn	0
			1	1	

Continued on next page...

Continued from previous page...

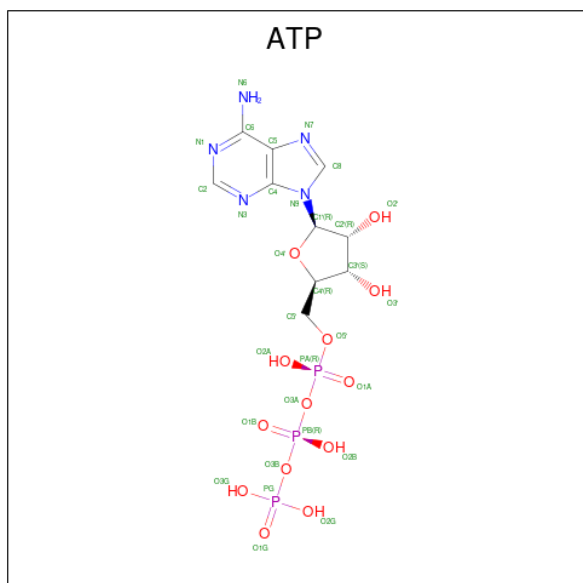
Mol	Chain	Residues	Atoms		AltConf
39	f	1	Total	Zn	0
			1	1	

- Molecule 40 is ASPARTIC ACID (CCD ID: ASP) (formula: $C_4H_7NO_4$).



Mol	Chain	Residues	Atoms				AltConf
40	U	1	Total	C	N	O	0
			8	4	1	3	

- Molecule 41 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$).

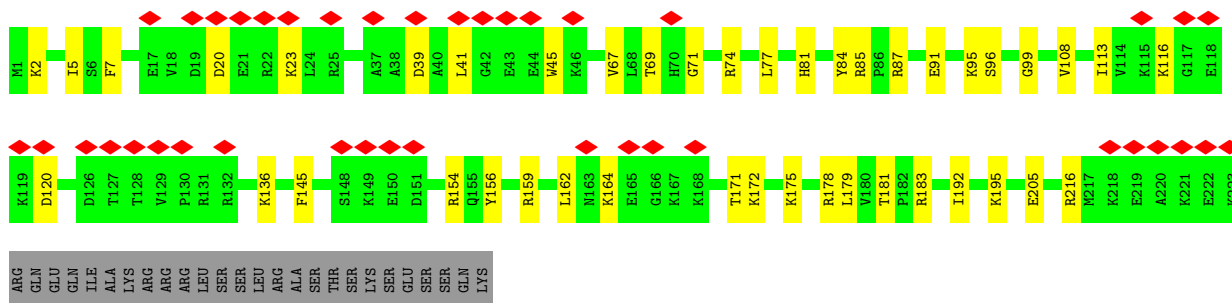


Mol	Chain	Residues	Atoms					AltConf
			Total	C	N	O	P	
41	z	1	31	10	5	13	3	0

ALA
MET
PRO
GLN
PRO
VAL
PRO
THR
ALA

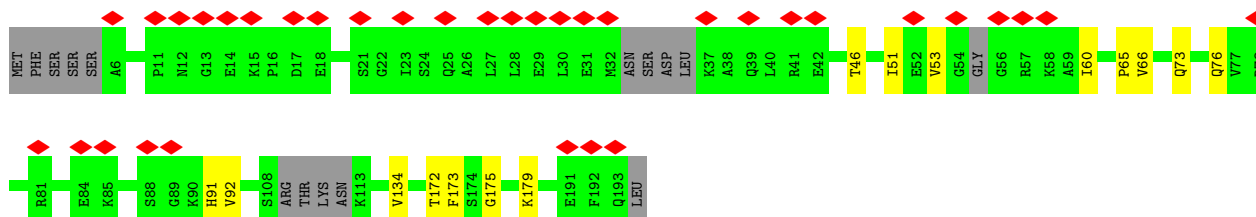
• Molecule 8: 40S ribosomal protein S6

Chain G: 16% 72% 17% 10%



• Molecule 9: 40S ribosomal protein S7

Chain H: 18% 85% 8% 8%



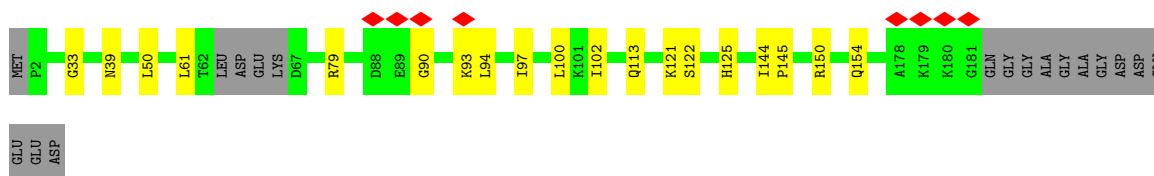
• Molecule 10: 40S ribosomal protein S8

Chain I: 11% 82% 13% .

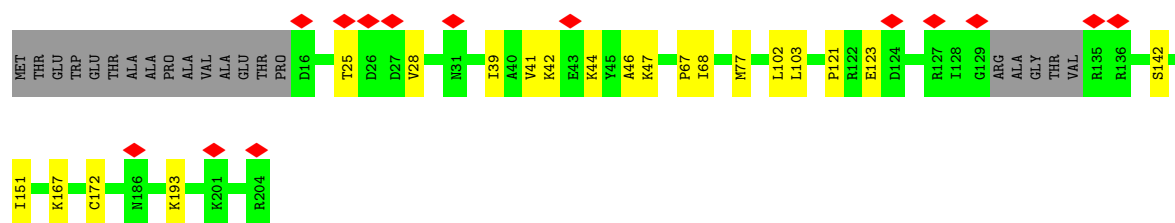
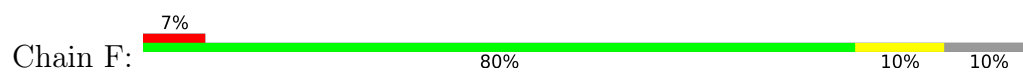


• Molecule 11: 40S ribosomal protein S9

Chain J: . 81% 10% 9%



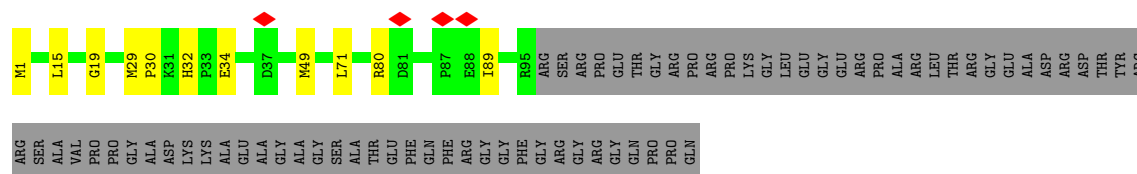
• Molecule 12: 40S ribosomal protein S5



- Molecule 13: 40S ribosomal protein S11



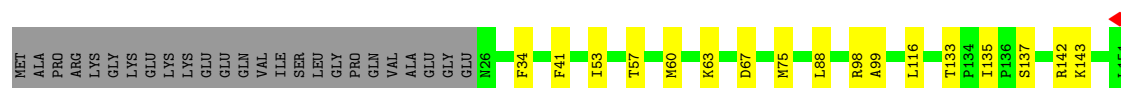
- Molecule 14: 40S ribosomal protein S10



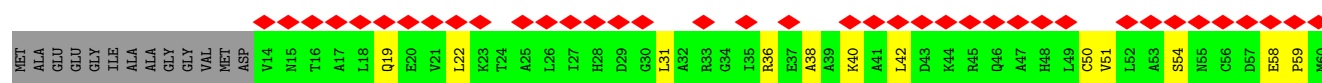
- Molecule 15: 40S ribosomal protein S13



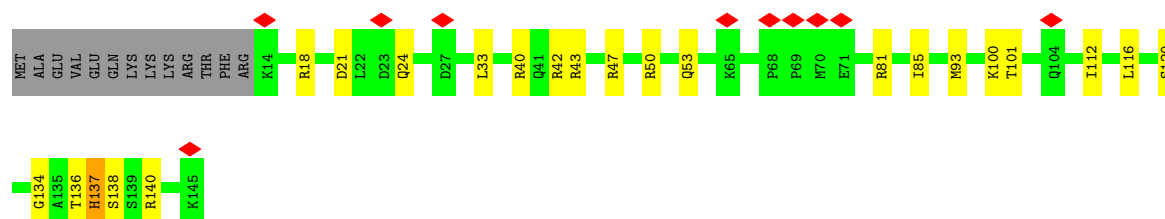
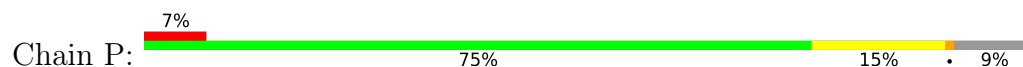
- Molecule 16: 40S ribosomal protein S14



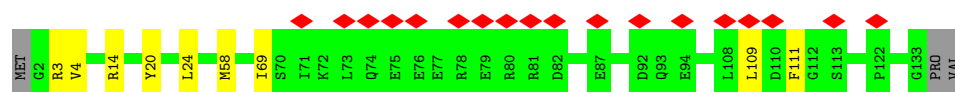
- Molecule 17: 40S ribosomal protein S12



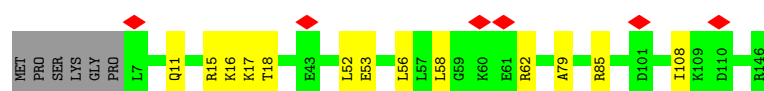
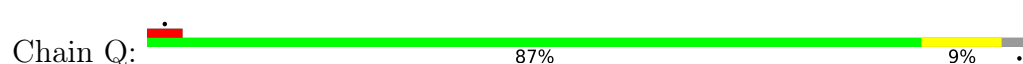
- Molecule 18: 40S ribosomal protein S15



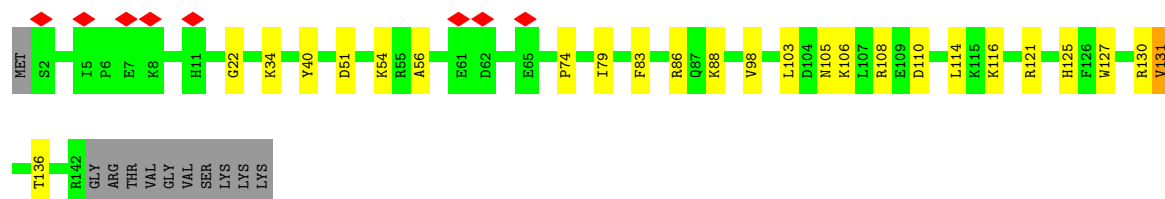
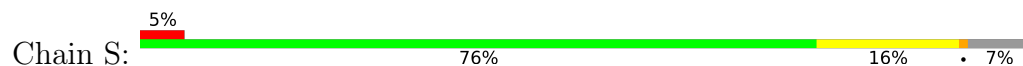
- Molecule 19: 40S ribosomal protein S17



- Molecule 20: 40S ribosomal protein S16



- Molecule 21: 40S ribosomal protein S18

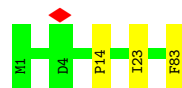


- Molecule 22: 40S ribosomal protein S19





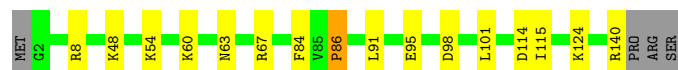
- Molecule 23: 40S ribosomal protein S21



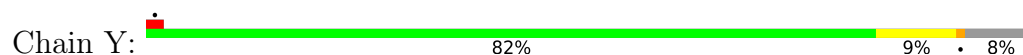
- Molecule 24: 40S ribosomal protein S15a



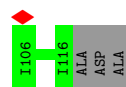
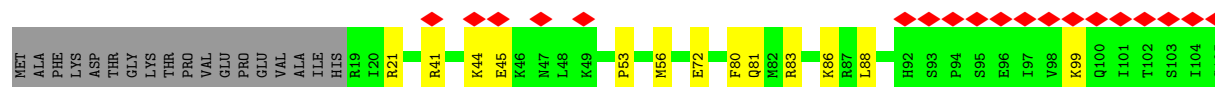
- Molecule 25: 40S ribosomal protein S23



- Molecule 26: 40S ribosomal protein S24



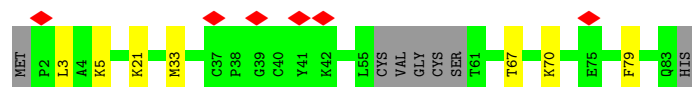
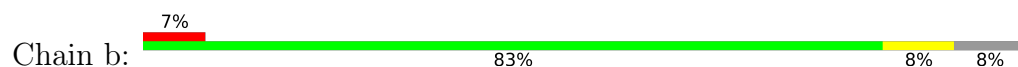
- Molecule 27: 40S ribosomal protein S20



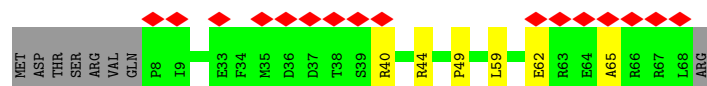
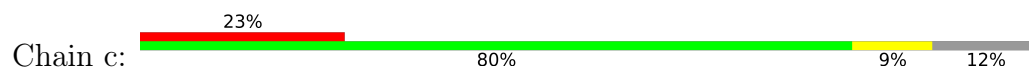
- Molecule 28: 40S ribosomal protein S25



- Molecule 29: 40S ribosomal protein S27



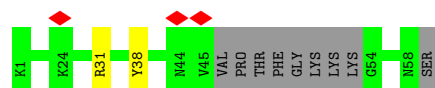
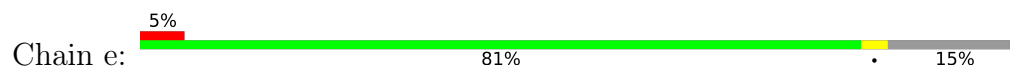
- Molecule 30: 40S ribosomal protein S28



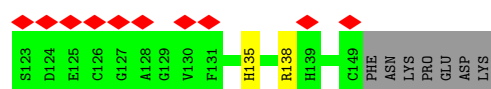
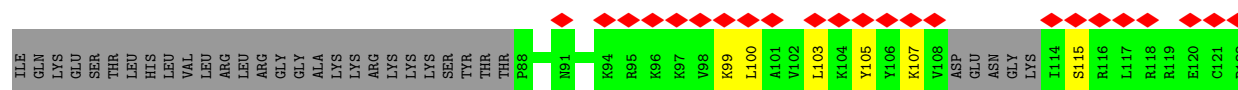
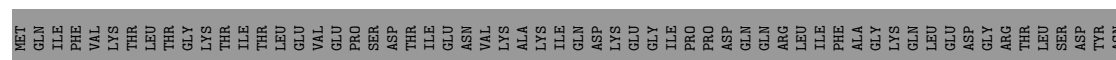
- Molecule 31: 40S ribosomal protein S29



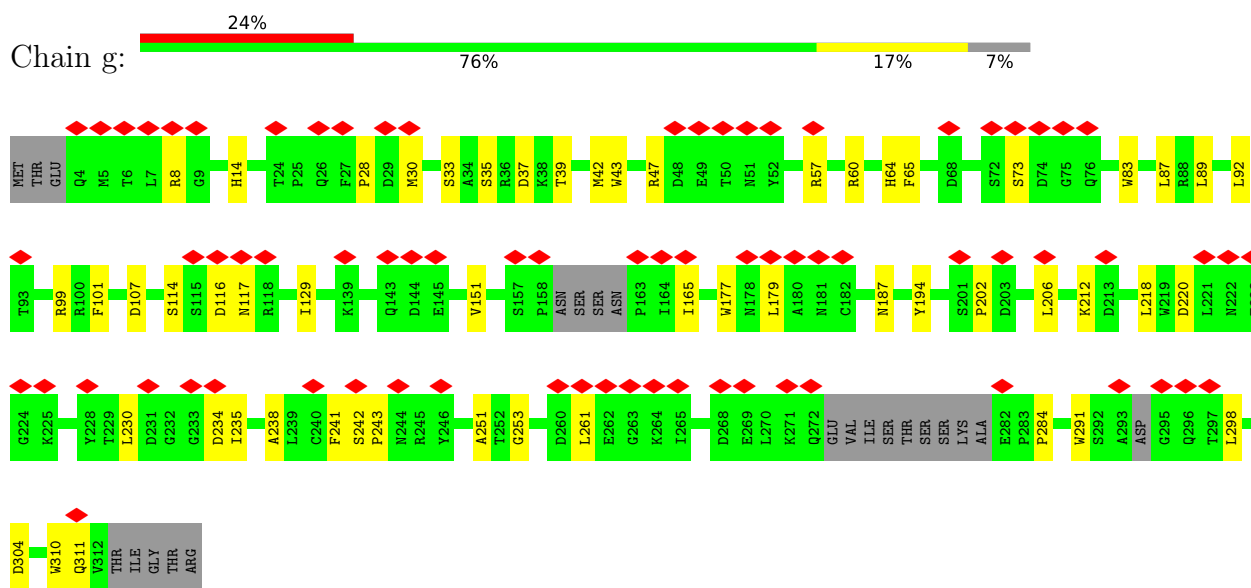
- Molecule 32: 40S ribosomal protein S30



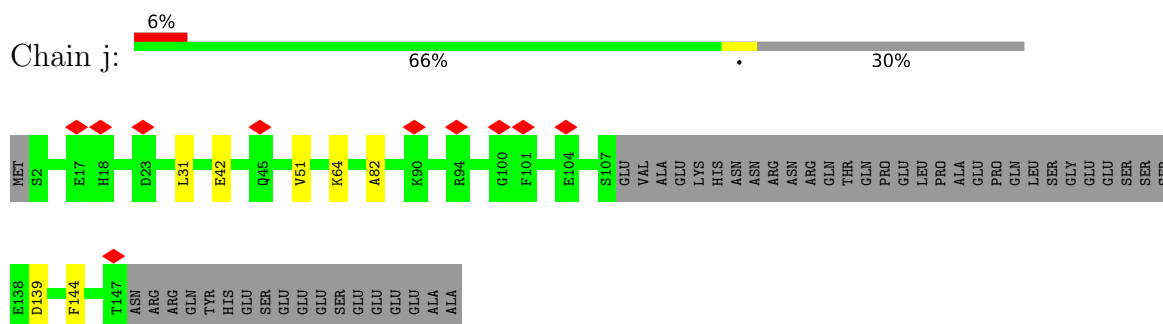
- Molecule 33: Ubiquitin-40S ribosomal protein S27a



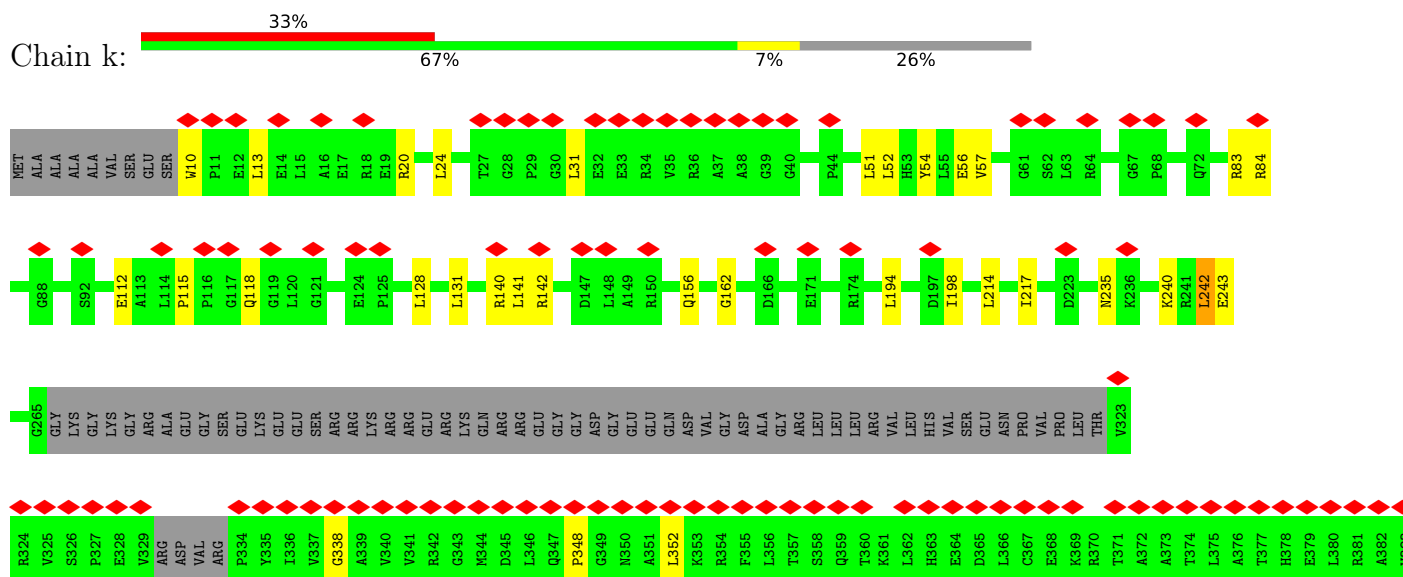
- Molecule 34: Receptor of activated protein C kinase 1



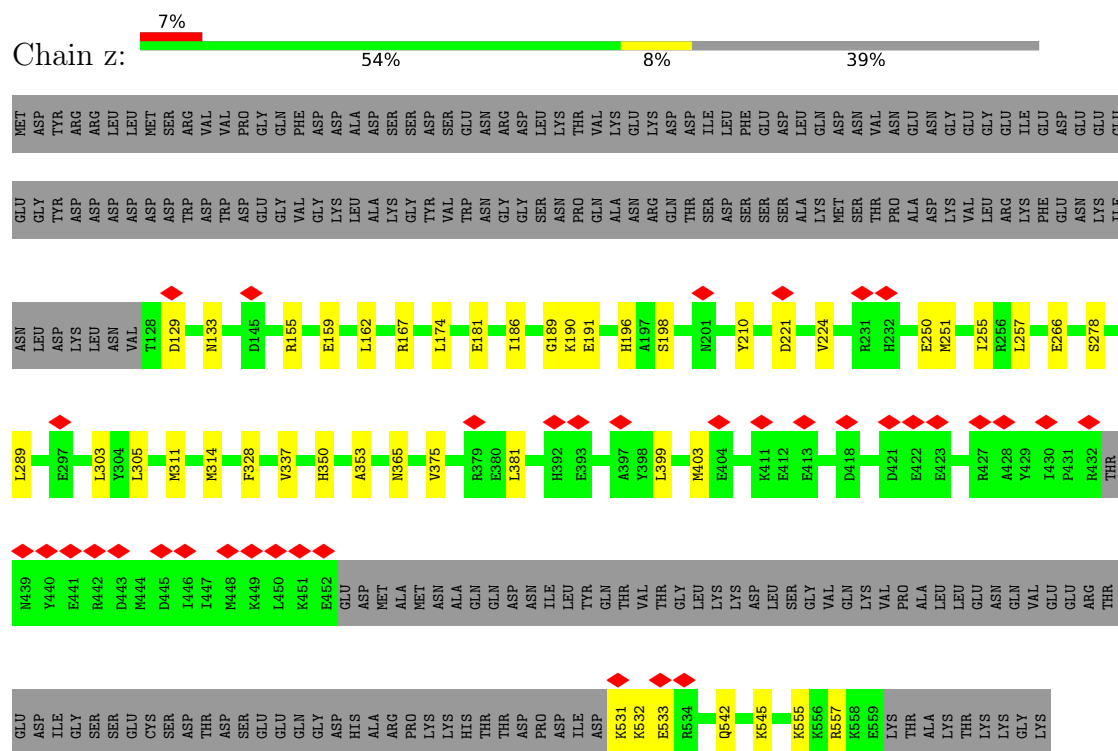
- Molecule 35: Probable RNA-binding protein EIF1AD



- Molecule 36: Leucine-rich repeat-containing protein 47



- Molecule 37: Serine/threonine-protein kinase RIO1



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	121468	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	48	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.856	Depositor
Minimum map value	-0.439	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.016	Depositor
Recommended contour level	0.06	Depositor
Map size (Å)	381.24, 381.24, 381.24	wwPDB
Map dimensions	360, 360, 360	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.059, 1.059, 1.059	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 6MZ, MA6, OMG, OMC, PSU, MG, ZN, 7MG, 4AC, A2M, ATP, B8N, OMU

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	0.40	3/37709 (0.0%)	0.44	2/58755 (0.0%)
2	A	0.32	0/1742	0.50	0/2367
3	B	0.27	0/1742	0.47	0/2330
4	C	0.33	0/1710	0.48	0/2310
5	h	0.39	0/805	0.58	0/1079
6	E	0.30	0/2073	0.49	0/2791
7	D	0.27	0/1773	0.50	0/2387
8	G	0.28	0/1825	0.52	0/2431
9	H	0.25	0/1466	0.48	0/1959
10	I	0.28	0/1666	0.48	0/2223
11	J	0.28	0/1489	0.49	0/1987
12	F	0.27	0/1481	0.53	0/1988
13	L	0.33	0/1147	0.44	0/1535
14	K	0.26	0/823	0.45	0/1111
15	N	0.29	0/1226	0.42	0/1649
16	O	0.30	0/959	0.49	0/1286
17	M	0.21	0/845	0.55	0/1134
18	P	0.28	0/1099	0.57	2/1471 (0.1%)
19	R	0.28	0/1080	0.53	0/1449
20	Q	0.29	0/1133	0.58	0/1517
21	S	0.30	0/1180	0.54	0/1582
22	T	0.31	0/1113	0.47	0/1493
23	V	0.28	0/643	0.42	0/860
24	W	0.36	0/1051	0.55	0/1406
25	X	0.34	0/1097	0.52	0/1464
26	Y	0.26	0/1004	0.47	0/1337
27	U	0.31	0/783	0.53	0/1052
28	Z	0.24	0/646	0.51	0/869
29	b	0.29	0/623	0.50	0/834
30	c	0.22	0/481	0.57	0/643
31	d	0.38	0/461	0.61	0/612
32	e	0.27	0/401	0.51	0/526

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
33	f	0.27	0/474	0.54	0/626
34	g	0.24	0/2360	0.51	0/3207
35	j	0.26	0/952	0.47	0/1282
36	k	0.23	0/2920	0.50	0/3977
37	z	0.28	0/2848	0.57	0/3816
All	All	0.34	3/82830 (0.0%)	0.48	4/119345 (0.0%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
16	O	0	1
18	P	0	2
20	Q	0	1
25	X	0	1
All	All	0	5

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	2	668	A2M	O3'-P	5.38	1.61	1.56
1	2	484	A2M	O3'-P	5.18	1.61	1.56
1	2	1383	A2M	O3'-P	5.12	1.61	1.56

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
18	P	136	THR	CA-C-N	6.74	134.42	121.54
18	P	136	THR	C-N-CA	6.74	134.42	121.54
1	2	1060	A	P-O3'-C3'	5.12	127.87	120.20
1	2	381	C	C2'-C3'-O3'	5.08	121.31	113.70

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
16	O	137	SER	Peptide
18	P	137	HIS	Mainchain
18	P	138	SER	Peptide

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Group
20	Q	15	ARG	Peptide
25	X	60	LYS	Peptide

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	2	35310	0	17883	157	0
2	A	1705	0	1706	14	0
3	B	1715	0	1785	14	0
4	C	1674	0	1764	14	0
5	h	792	0	841	7	0
6	E	2031	0	2133	16	0
7	D	1745	0	1839	8	0
8	G	1802	0	1953	28	0
9	H	1446	0	1540	8	0
10	I	1638	0	1710	19	0
11	J	1465	0	1583	13	0
12	F	1461	0	1511	14	0
13	L	1127	0	1190	7	0
14	K	799	0	823	6	0
15	N	1202	0	1289	9	0
16	O	947	0	976	13	0
17	M	837	0	870	17	0
18	P	1077	0	1123	15	0
19	R	1066	0	1116	7	0
20	Q	1116	0	1185	7	0
21	S	1162	0	1215	15	0
22	T	1094	0	1120	10	0
23	V	636	0	637	2	0
24	W	1034	0	1080	6	0
25	X	1080	0	1147	10	0
26	Y	987	0	1038	9	0
27	U	774	0	839	11	0
28	Z	638	0	679	9	0
29	b	611	0	637	4	0
30	c	479	0	507	4	0
31	d	450	0	444	4	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
32	e	398	0	439	3	0
33	f	465	0	481	5	0
34	g	2306	0	2266	32	0
35	j	934	0	928	5	0
36	k	2886	0	2563	21	0
37	z	2803	0	2799	27	0
38	2	123	0	0	0	0
38	F	1	0	0	0	0
38	G	1	0	0	0	0
38	O	1	0	0	0	0
38	X	1	0	0	0	0
38	d	1	0	0	0	0
38	h	1	0	0	0	0
38	z	2	0	0	0	0
39	d	1	0	0	0	0
39	f	1	0	0	0	0
39	h	1	0	0	0	0
40	U	8	0	3	1	0
41	z	31	0	12	1	0
All	All	79865	0	63654	480	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.

The worst 5 of 480 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:1748:G:H1	1:2:1786:U:H3	1.24	0.82
13:L:104:LYS:HD2	25:X:8:ARG:HH21	1.50	0.77
36:k:217:ILE:HD12	36:k:242:LEU:HD11	1.68	0.75
18:P:137:HIS:HA	18:P:140:ARG:HE	1.51	0.74
1:2:1566:G:N7	22:T:101:ARG:NH2	2.39	0.71

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	A	214/295 (72%)	205 (96%)	9 (4%)	0	100	100
3	B	209/264 (79%)	205 (98%)	4 (2%)	0	100	100
4	C	214/293 (73%)	206 (96%)	8 (4%)	0	100	100
5	h	97/115 (84%)	96 (99%)	1 (1%)	0	100	100
6	E	253/263 (96%)	247 (98%)	6 (2%)	0	100	100
7	D	222/243 (91%)	221 (100%)	1 (0%)	0	100	100
8	G	221/249 (89%)	218 (99%)	3 (1%)	0	100	100
9	H	171/194 (88%)	165 (96%)	6 (4%)	0	100	100
10	I	195/208 (94%)	187 (96%)	8 (4%)	0	100	100
11	J	172/194 (89%)	165 (96%)	7 (4%)	0	100	100
12	F	180/204 (88%)	172 (96%)	8 (4%)	0	100	100
13	L	132/158 (84%)	129 (98%)	3 (2%)	0	100	100
14	K	93/165 (56%)	87 (94%)	6 (6%)	0	100	100
15	N	147/151 (97%)	147 (100%)	0	0	100	100
16	O	124/151 (82%)	121 (98%)	3 (2%)	0	100	100
17	M	102/132 (77%)	98 (96%)	4 (4%)	0	100	100
18	P	130/145 (90%)	124 (95%)	6 (5%)	0	100	100
19	R	130/135 (96%)	125 (96%)	5 (4%)	0	100	100
20	Q	138/146 (94%)	131 (95%)	7 (5%)	0	100	100
21	S	139/152 (91%)	132 (95%)	7 (5%)	0	100	100
22	T	139/145 (96%)	138 (99%)	1 (1%)	0	100	100
23	V	81/83 (98%)	79 (98%)	2 (2%)	0	100	100
24	W	127/130 (98%)	123 (97%)	4 (3%)	0	100	100
25	X	137/143 (96%)	135 (98%)	1 (1%)	1 (1%)	18	38
26	Y	120/133 (90%)	117 (98%)	3 (2%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
27	U	96/119 (81%)	93 (97%)	3 (3%)	0	100	100
28	Z	80/125 (64%)	75 (94%)	5 (6%)	0	100	100
29	b	73/84 (87%)	67 (92%)	6 (8%)	0	100	100
30	c	59/69 (86%)	57 (97%)	2 (3%)	0	100	100
31	d	52/56 (93%)	49 (94%)	3 (6%)	0	100	100
32	e	46/59 (78%)	44 (96%)	2 (4%)	0	100	100
33	f	53/156 (34%)	49 (92%)	4 (8%)	0	100	100
34	g	287/317 (90%)	269 (94%)	18 (6%)	0	100	100
35	j	112/165 (68%)	106 (95%)	6 (5%)	0	100	100
36	k	418/583 (72%)	388 (93%)	30 (7%)	0	100	100
37	z	343/568 (60%)	319 (93%)	24 (7%)	0	100	100
All	All	5506/6792 (81%)	5289 (96%)	216 (4%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
25	X	86	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	
2	A	180/243 (74%)	179 (99%)	1 (1%)	78	91
3	B	192/231 (83%)	192 (100%)	0	100	100
4	C	182/225 (81%)	182 (100%)	0	100	100
5	h	86/98 (88%)	84 (98%)	2 (2%)	44	71
6	E	220/225 (98%)	219 (100%)	1 (0%)	81	92
7	D	188/202 (93%)	187 (100%)	1 (0%)	81	92
8	G	194/218 (89%)	193 (100%)	1 (0%)	81	92

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
9	H	160/174 (92%)	160 (100%)	0	100	100
10	I	174/180 (97%)	174 (100%)	0	100	100
11	J	156/168 (93%)	156 (100%)	0	100	100
12	F	156/170 (92%)	156 (100%)	0	100	100
13	L	126/142 (89%)	126 (100%)	0	100	100
14	K	86/136 (63%)	86 (100%)	0	100	100
15	N	130/131 (99%)	130 (100%)	0	100	100
16	O	99/119 (83%)	99 (100%)	0	100	100
17	M	91/108 (84%)	91 (100%)	0	100	100
18	P	117/130 (90%)	117 (100%)	0	100	100
19	R	118/122 (97%)	118 (100%)	0	100	100
20	Q	116/121 (96%)	115 (99%)	1 (1%)	70	87
21	S	122/132 (92%)	121 (99%)	1 (1%)	73	88
22	T	111/115 (96%)	111 (100%)	0	100	100
23	V	67/67 (100%)	67 (100%)	0	100	100
24	W	112/113 (99%)	112 (100%)	0	100	100
25	X	111/115 (96%)	111 (100%)	0	100	100
26	Y	103/115 (90%)	102 (99%)	1 (1%)	68	86
27	U	90/107 (84%)	90 (100%)	0	100	100
28	Z	67/103 (65%)	67 (100%)	0	100	100
29	b	70/76 (92%)	70 (100%)	0	100	100
30	c	54/62 (87%)	54 (100%)	0	100	100
31	d	47/49 (96%)	46 (98%)	1 (2%)	47	73
32	e	40/48 (83%)	40 (100%)	0	100	100
33	f	51/140 (36%)	51 (100%)	0	100	100
34	g	255/275 (93%)	255 (100%)	0	100	100
35	j	103/150 (69%)	103 (100%)	0	100	100
36	k	231/487 (47%)	230 (100%)	1 (0%)	84	93
37	z	299/512 (58%)	299 (100%)	0	100	100
All	All	4704/5809 (81%)	4693 (100%)	11 (0%)	85	95

5 of 11 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
21	S	131	VAL
26	Y	18	LEU
36	k	242	LEU
31	d	23	VAL
7	D	72	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 98 such sidechains are listed below:

Mol	Chain	Res	Type
21	S	42	HIS
26	Y	63	HIS
21	S	120	HIS
24	W	5	ASN
29	b	9	HIS

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	2	1640/1869 (87%)	353 (21%)	13 (0%)

5 of 353 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	2	2	A
1	2	17	C
1	2	26	U
1	2	33	G
1	2	41	G

5 of 13 RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	2	1060	A
1	2	1065	G
1	2	1814	G
1	2	1440	C
1	2	1605	G

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

74 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$
1	PSU	2	218	1	18,21,22	1.01	1 (5%)	22,30,33	1.75	4 (18%)
1	PSU	2	34	1	18,21,22	1.06	1 (5%)	22,30,33	1.74	4 (18%)
1	OMC	2	1703	1	19,22,23	2.91	8 (42%)	26,31,34	0.90	2 (7%)
1	A2M	2	27	1,38	22,25,26	3.42	10 (45%)	31,36,39	2.47	10 (32%)
1	PSU	2	1232	1	18,21,22	1.14	1 (5%)	22,30,33	1.74	4 (18%)
1	OMC	2	462	1	19,22,23	2.90	8 (42%)	26,31,34	0.79	0
1	OMC	2	174	1	19,22,23	2.88	8 (42%)	26,31,34	0.80	1 (3%)
1	OMU	2	116	1	19,22,23	2.89	7 (36%)	26,31,34	1.76	5 (19%)
1	4AC	2	1842	1,38	21,24,25	3.12	11 (52%)	29,34,37	1.82	6 (20%)
1	OMG	2	601	1	23,26,27	2.23	8 (34%)	33,38,41	1.80	8 (24%)
1	OMU	2	428	1	19,22,23	3.02	8 (42%)	26,31,34	1.93	5 (19%)
1	PSU	2	406	1	18,21,22	1.02	1 (5%)	22,30,33	1.84	4 (18%)
1	A2M	2	484	1	22,25,26	3.38	9 (40%)	31,36,39	2.53	11 (35%)
1	PSU	2	36	1	18,21,22	0.99	1 (5%)	22,30,33	1.90	3 (13%)
1	PSU	2	93	1	18,21,22	1.17	1 (5%)	22,30,33	1.72	4 (18%)
1	PSU	2	1081	1	18,21,22	1.00	2 (11%)	22,30,33	1.76	5 (22%)
1	OMG	2	644	1	23,26,27	2.23	8 (34%)	33,38,41	1.81	7 (21%)
1	PSU	2	1056	1	18,21,22	1.03	1 (5%)	22,30,33	1.86	4 (18%)
1	PSU	2	1174	1,38	18,21,22	1.07	1 (5%)	22,30,33	1.78	4 (18%)
1	PSU	2	119	1	18,21,22	1.07	1 (5%)	22,30,33	1.66	5 (22%)
1	OMU	2	354	1	19,22,23	2.83	7 (36%)	26,31,34	1.74	4 (15%)
1	A2M	2	468	1	22,25,26	3.42	9 (40%)	31,36,39	2.71	12 (38%)
1	PSU	2	1004	1	18,21,22	1.08	1 (5%)	22,30,33	1.72	4 (18%)
1	OMU	2	121	1	19,22,23	2.85	7 (36%)	26,31,34	1.74	5 (19%)
1	PSU	2	1244	1,38	18,21,22	1.11	1 (5%)	22,30,33	1.70	3 (13%)
1	PSU	2	1445	1	18,21,22	1.08	1 (5%)	22,30,33	1.85	4 (18%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	PSU	2	814	1	18,21,22	0.93	1 (5%)	22,30,33	1.75	3 (13%)
1	PSU	2	681	1	18,21,22	1.12	1 (5%)	22,30,33	1.81	4 (18%)
1	PSU	2	822	1	18,21,22	0.99	1 (5%)	22,30,33	1.79	5 (22%)
1	OMC	2	1272	1	19,22,23	2.85	8 (42%)	26,31,34	0.76	0
1	OMG	2	1328	1	23,26,27	2.27	7 (30%)	33,38,41	1.91	8 (24%)
1	PSU	2	1625	1	18,21,22	1.11	1 (5%)	22,30,33	1.70	4 (18%)
1	PSU	2	1238	1	18,21,22	1.12	1 (5%)	22,30,33	1.72	3 (13%)
1	A2M	2	1678	1	22,25,26	3.49	9 (40%)	31,36,39	2.63	11 (35%)
1	A2M	2	1383	1	22,25,26	3.44	10 (45%)	31,36,39	2.56	9 (29%)
1	7MG	2	1639	1	22,26,27	3.59	10 (45%)	29,39,42	2.05	9 (31%)
1	PSU	2	1692	1	18,21,22	1.17	1 (5%)	22,30,33	1.86	5 (22%)
1	OMU	2	172	1	19,22,23	2.88	8 (42%)	26,31,34	1.74	5 (19%)
1	OMC	2	517	1	19,22,23	2.74	7 (36%)	26,31,34	0.76	1 (3%)
1	PSU	2	966	1	18,21,22	0.95	1 (5%)	22,30,33	1.64	3 (13%)
1	6MZ	2	1832	1,38	22,25,26	2.47	5 (22%)	30,36,39	3.16	11 (36%)
1	MA6	2	1850	1	23,26,27	1.43	4 (17%)	34,38,41	3.26	12 (35%)
1	OMG	2	509	1,38	23,26,27	2.29	7 (30%)	33,38,41	1.90	7 (21%)
1	A2M	2	159	1	22,25,26	3.42	9 (40%)	31,36,39	2.52	10 (32%)
1	PSU	2	649	1	18,21,22	1.11	1 (5%)	22,30,33	1.79	4 (18%)
1	PSU	2	105	1	18,21,22	1.05	1 (5%)	22,30,33	1.78	3 (13%)
1	PSU	2	1347	1	18,21,22	1.06	1 (5%)	22,30,33	1.69	4 (18%)
1	OMU	2	1288	1	19,22,23	2.97	8 (42%)	26,31,34	1.69	5 (19%)
1	OMU	2	1442	1,38	19,22,23	2.93	8 (42%)	26,31,34	1.77	4 (15%)
1	OMU	2	1326	1,38	19,22,23	2.99	8 (42%)	26,31,34	1.86	5 (19%)
1	A2M	2	1031	1	22,25,26	3.44	10 (45%)	31,36,39	2.51	8 (25%)
1	A2M	2	166	1	22,25,26	3.45	10 (45%)	31,36,39	2.60	9 (29%)
1	A2M	2	668	1,38	22,25,26	3.42	10 (45%)	31,36,39	2.66	9 (29%)
1	OMG	2	867	1	23,26,27	2.31	9 (39%)	33,38,41	1.78	8 (24%)
1	PSU	2	1367	1	18,21,22	1.04	1 (5%)	22,30,33	1.78	4 (18%)
1	4AC	2	1337	1	21,24,25	3.02	11 (52%)	29,34,37	1.11	2 (6%)
1	PSU	2	686	1	18,21,22	1.03	1 (5%)	22,30,33	1.79	4 (18%)
1	PSU	2	1046	1	18,21,22	1.01	1 (5%)	22,30,33	1.71	4 (18%)
1	PSU	2	109	1	18,21,22	1.03	1 (5%)	22,30,33	1.75	4 (18%)
1	PSU	2	1643	1,38	18,21,22	1.04	1 (5%)	22,30,33	1.60	3 (13%)
1	PSU	2	572	1	18,21,22	1.01	1 (5%)	22,30,33	1.77	3 (13%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	PSU	2	815	1	18,21,22	1.01	1 (5%)	22,30,33	1.75	2 (9%)
1	B8N	2	1248	1	24,29,30	2.98	9 (37%)	29,42,45	1.83	5 (17%)
1	OMC	2	1391	1	19,22,23	2.82	8 (42%)	26,31,34	0.85	1 (3%)
1	PSU	2	1177	1	18,21,22	1.07	1 (5%)	22,30,33	1.64	4 (18%)
1	OMG	2	683	1	23,26,27	2.26	9 (39%)	33,38,41	2.17	14 (42%)
1	PSU	2	1045	1	18,21,22	1.03	1 (5%)	22,30,33	1.69	4 (18%)
1	OMG	2	1490	1,38	23,26,27	2.21	6 (26%)	33,38,41	1.71	7 (21%)
1	PSU	2	651	1	18,21,22	1.03	1 (5%)	22,30,33	1.74	4 (18%)
1	MA6	2	1851	1	23,26,27	1.43	4 (17%)	34,38,41	3.16	12 (35%)
1	A2M	2	576	1	22,25,26	3.44	9 (40%)	31,36,39	2.44	8 (25%)
1	A2M	2	512	1	22,25,26	3.41	10 (45%)	31,36,39	2.69	8 (25%)
1	A2M	2	99	1,38	22,25,26	3.50	10 (45%)	31,36,39	2.43	9 (29%)
1	OMG	2	436	1	23,26,27	2.29	9 (39%)	33,38,41	1.91	8 (24%)

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
1	PSU	2	218	1	-	0/7/25/26	0/2/2/2
1	PSU	2	34	1	-	0/7/25/26	0/2/2/2
1	OMC	2	1703	1	-	0/9/27/28	0/2/2/2
1	A2M	2	27	1,38	-	0/9/27/28	0/3/3/3
1	PSU	2	1232	1	-	0/7/25/26	0/2/2/2
1	OMC	2	462	1	-	0/9/27/28	0/2/2/2
1	OMC	2	174	1	-	0/9/27/28	0/2/2/2
1	OMU	2	116	1	-	0/9/27/28	0/2/2/2
1	4AC	2	1842	1,38	-	2/11/29/30	0/2/2/2
1	OMG	2	601	1	-	0/9/27/28	0/3/3/3
1	OMU	2	428	1	-	6/9/27/28	0/2/2/2
1	PSU	2	406	1	-	0/7/25/26	0/2/2/2
1	A2M	2	484	1	-	0/9/27/28	0/3/3/3
1	PSU	2	36	1	-	0/7/25/26	0/2/2/2
1	PSU	2	93	1	-	0/7/25/26	0/2/2/2
1	PSU	2	1081	1	-	0/7/25/26	0/2/2/2
1	OMG	2	644	1	-	3/9/27/28	0/3/3/3
1	PSU	2	1056	1	-	0/7/25/26	0/2/2/2
1	PSU	2	1174	1,38	-	0/7/25/26	0/2/2/2

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	PSU	2	119	1	-	0/7/25/26	0/2/2/2
1	OMU	2	354	1	-	2/9/27/28	0/2/2/2
1	A2M	2	468	1	-	3/9/27/28	0/3/3/3
1	PSU	2	1004	1	-	0/7/25/26	0/2/2/2
1	OMU	2	121	1	-	0/9/27/28	0/2/2/2
1	PSU	2	1244	1,38	-	0/7/25/26	0/2/2/2
1	PSU	2	1445	1	-	0/7/25/26	0/2/2/2
1	PSU	2	814	1	-	0/7/25/26	0/2/2/2
1	PSU	2	681	1	-	0/7/25/26	0/2/2/2
1	PSU	2	822	1	-	0/7/25/26	0/2/2/2
1	OMC	2	1272	1	-	0/9/27/28	0/2/2/2
1	OMG	2	1328	1	-	0/9/27/28	0/3/3/3
1	PSU	2	1625	1	-	0/7/25/26	0/2/2/2
1	PSU	2	1238	1	-	0/7/25/26	0/2/2/2
1	A2M	2	1678	1	-	1/9/27/28	0/3/3/3
1	A2M	2	1383	1	-	0/9/27/28	0/3/3/3
1	7MG	2	1639	1	-	1/7/37/38	0/3/3/3
1	PSU	2	1692	1	-	2/7/25/26	0/2/2/2
1	OMU	2	172	1	-	3/9/27/28	0/2/2/2
1	OMC	2	517	1	-	0/9/27/28	0/2/2/2
1	PSU	2	966	1	-	0/7/25/26	0/2/2/2
1	6MZ	2	1832	1,38	-	0/9/27/28	0/3/3/3
1	MA6	2	1850	1	-	0/11/29/30	0/3/3/3
1	OMG	2	509	1,38	-	2/9/27/28	0/3/3/3
1	A2M	2	159	1	-	0/9/27/28	0/3/3/3
1	PSU	2	649	1	-	2/7/25/26	0/2/2/2
1	PSU	2	105	1	-	0/7/25/26	0/2/2/2
1	PSU	2	1347	1	-	0/7/25/26	0/2/2/2
1	OMU	2	1288	1	-	0/9/27/28	0/2/2/2
1	OMU	2	1442	1,38	-	2/9/27/28	0/2/2/2
1	OMU	2	1326	1,38	-	0/9/27/28	0/2/2/2
1	A2M	2	1031	1	-	0/9/27/28	0/3/3/3
1	A2M	2	166	1	-	0/9/27/28	0/3/3/3
1	A2M	2	668	1,38	-	3/9/27/28	0/3/3/3
1	OMG	2	867	1	-	1/9/27/28	0/3/3/3
1	PSU	2	1367	1	-	0/7/25/26	0/2/2/2
1	4AC	2	1337	1	-	1/11/29/30	0/2/2/2
1	PSU	2	686	1	-	0/7/25/26	0/2/2/2
1	PSU	2	1046	1	-	0/7/25/26	0/2/2/2
1	PSU	2	109	1	-	0/7/25/26	0/2/2/2

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	PSU	2	1643	1,38	-	0/7/25/26	0/2/2/2
1	PSU	2	572	1	-	0/7/25/26	0/2/2/2
1	PSU	2	815	1	-	0/7/25/26	0/2/2/2
1	B8N	2	1248	1	-	6/16/34/35	0/2/2/2
1	OMC	2	1391	1	-	0/9/27/28	0/2/2/2
1	PSU	2	1177	1	-	0/7/25/26	0/2/2/2
1	OMG	2	683	1	-	2/9/27/28	0/3/3/3
1	PSU	2	1045	1	-	2/7/25/26	0/2/2/2
1	OMG	2	1490	1,38	-	1/9/27/28	0/3/3/3
1	PSU	2	651	1	-	0/7/25/26	0/2/2/2
1	MA6	2	1851	1	-	0/11/29/30	0/3/3/3
1	A2M	2	576	1	-	1/9/27/28	0/3/3/3
1	A2M	2	512	1	-	0/9/27/28	0/3/3/3
1	A2M	2	99	1,38	-	2/9/27/28	0/3/3/3
1	OMG	2	436	1	-	2/9/27/28	0/3/3/3

The worst 5 of 374 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	2	1832	6MZ	C6-N6	9.86	1.44	1.34
1	2	468	A2M	O4'-C1'	8.86	1.63	1.42
1	2	1678	A2M	C2'-C1'	-8.78	1.30	1.53
1	2	99	A2M	C2'-C1'	-8.74	1.30	1.53
1	2	1031	A2M	C2'-C1'	-8.74	1.30	1.53

The worst 5 of 407 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	2	1850	MA6	N1-C6-N6	-12.10	103.85	117.08
1	2	1851	MA6	N1-C6-N6	-12.02	103.94	117.08
1	2	1832	6MZ	C1'-N9-C8	7.60	144.29	127.14
1	2	1832	6MZ	C4-N9-C1'	-7.48	108.77	126.59
1	2	1851	MA6	C5-C6-N6	7.04	137.57	125.30

There are no chirality outliers.

5 of 50 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	2	172	OMU	O4'-C4'-C5'-O5'
1	2	428	OMU	C3'-C4'-C5'-O5'

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms
1	2	428	OMU	O4'-C4'-C5'-O5'
1	2	468	A2M	O4'-C1'-N9-C8
1	2	468	A2M	O4'-C1'-N9-C4

There are no ring outliers.

11 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	2	1232	PSU	2	0
1	2	116	OMU	1	0
1	2	468	A2M	1	0
1	2	814	PSU	1	0
1	2	681	PSU	1	0
1	2	1678	A2M	2	0
1	2	1639	7MG	1	0
1	2	509	OMG	1	0
1	2	1442	OMU	1	0
1	2	1248	B8N	1	0
1	2	1851	MA6	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 136 ligands modelled in this entry, 134 are monoatomic - leaving 2 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
41	ATP	z	602	38	29,33,33	1.32	4 (13%)	44,52,52	1.87	10 (22%)
40	ASP	U	201	-	6,7,8	0.89	0	5,8,10	0.96	0

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
41	ATP	z	602	38	-	3/22/38/38	0/3/3/3
40	ASP	U	201	-	-	3/5/6/8	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
41	z	602	ATP	C5-C4	4.14	1.46	1.39
41	z	602	ATP	C5-N7	-2.53	1.34	1.39
41	z	602	ATP	C5-C6	2.47	1.47	1.41
41	z	602	ATP	C4-N9	-2.07	1.33	1.37

The worst 5 of 10 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
41	z	602	ATP	C5-C4-N3	-6.12	118.76	126.75
41	z	602	ATP	N3-C4-N9	4.93	135.21	127.08
41	z	602	ATP	C2-N3-C4	3.91	121.00	111.75
41	z	602	ATP	N3-C2-N1	-3.28	123.47	128.60
41	z	602	ATP	C4-C5-N7	-2.96	107.01	110.62

There are no chirality outliers.

5 of 6 torsion outliers are listed below:

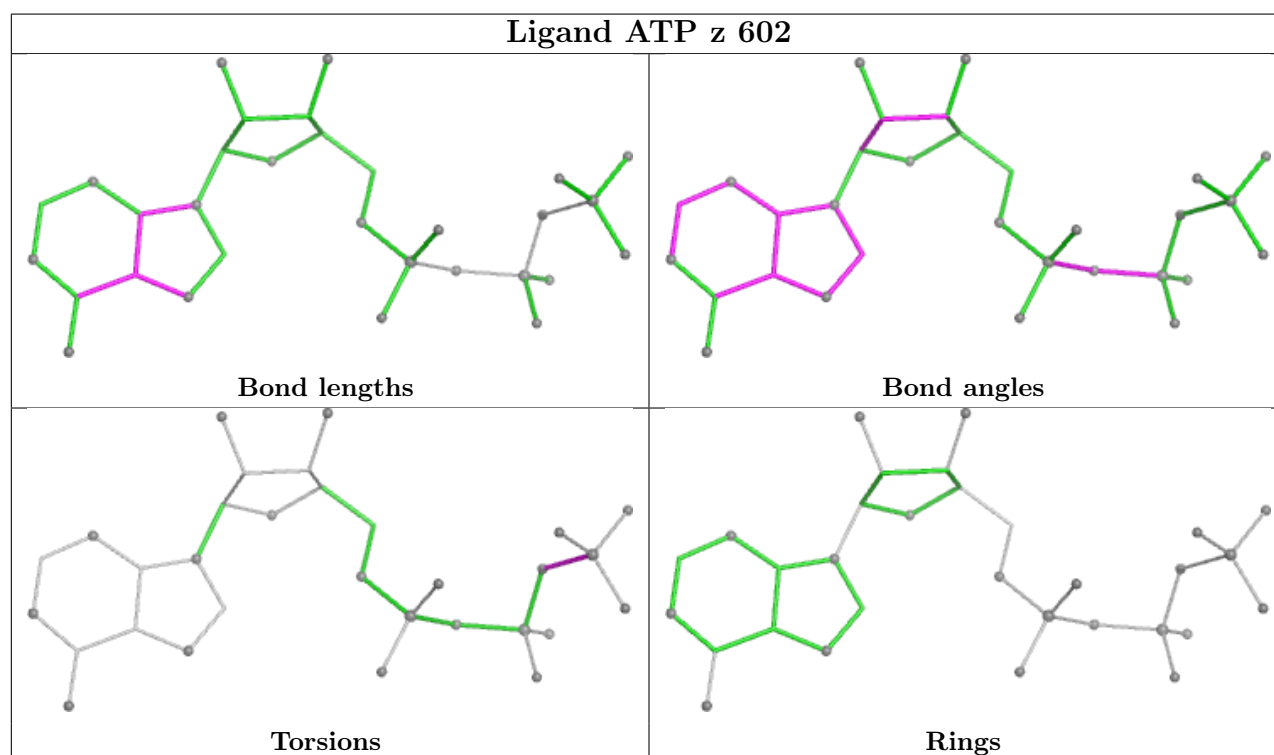
Mol	Chain	Res	Type	Atoms
40	U	201	ASP	O-C-CA-CB
40	U	201	ASP	CA-CB-CG-OD1
40	U	201	ASP	CA-CB-CG-OD2
41	z	602	ATP	PB-O3B-PG-O1G
41	z	602	ATP	PB-O3B-PG-O2G

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
41	z	602	ATP	1	0
40	U	201	ASP	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	2	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	2	563:G	O3'	564:A	P	3.14

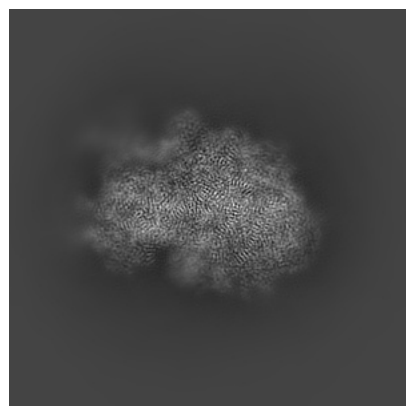
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-11520. These allow visual inspection of the internal detail of the map and identification of artifacts.

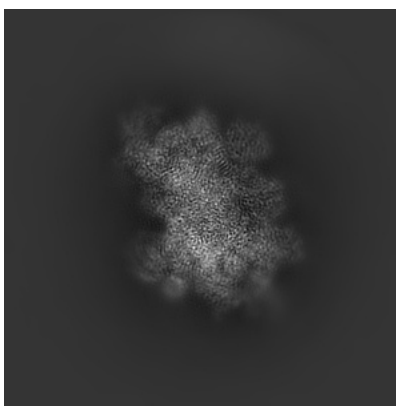
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

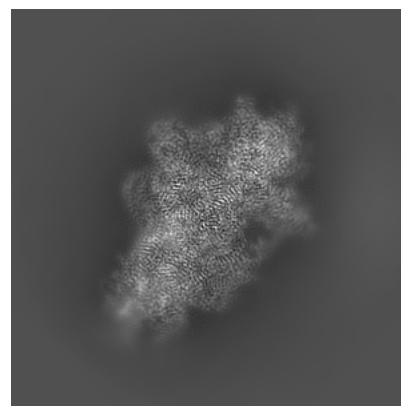
6.1.1 Primary map



X

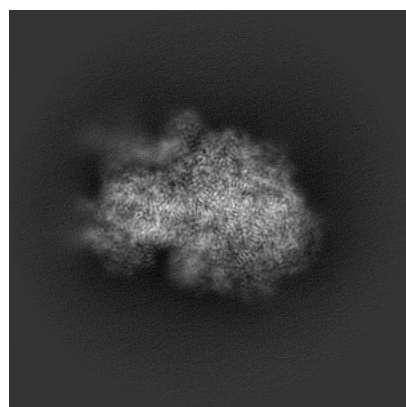


Y

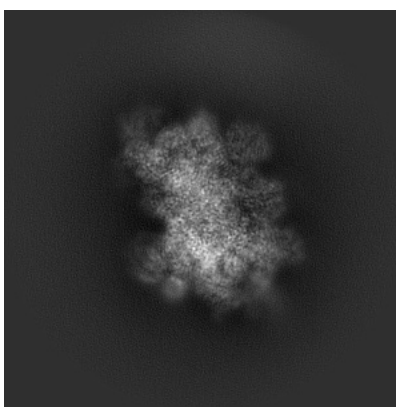


Z

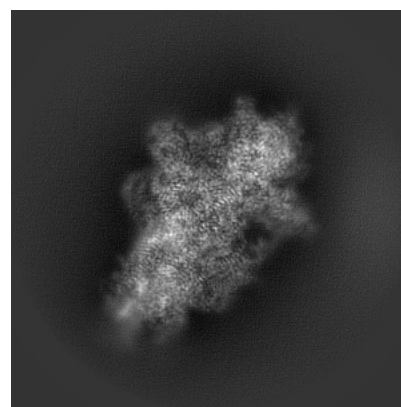
6.1.2 Raw map



X



Y

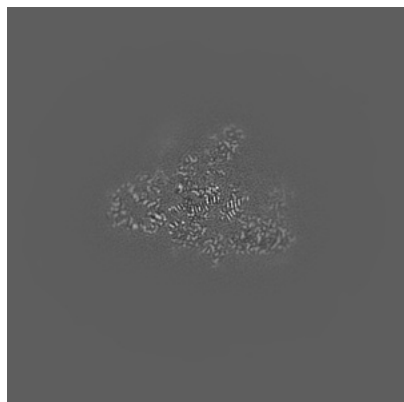


Z

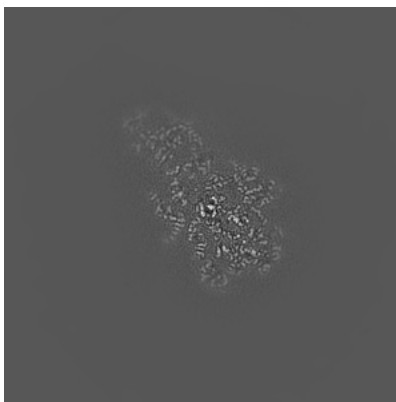
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

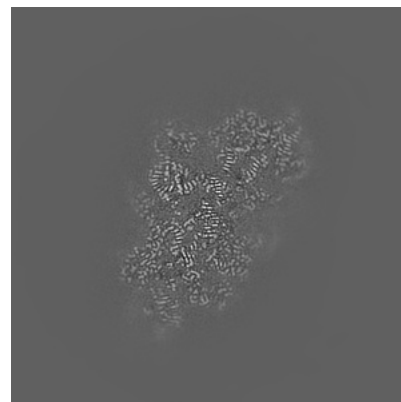
6.2.1 Primary map



X Index: 180

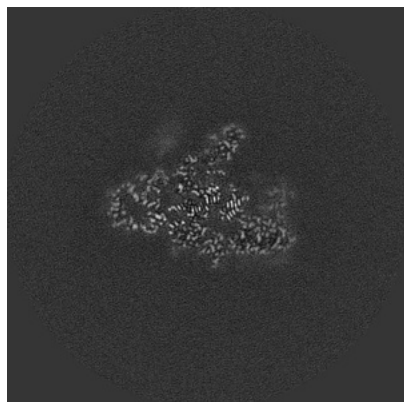


Y Index: 180

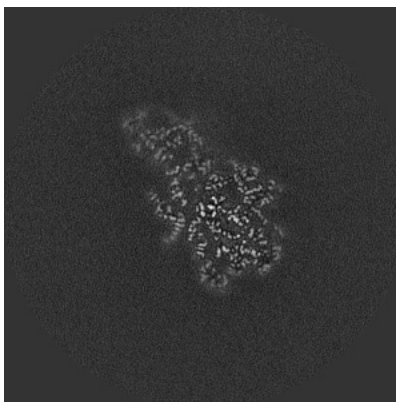


Z Index: 180

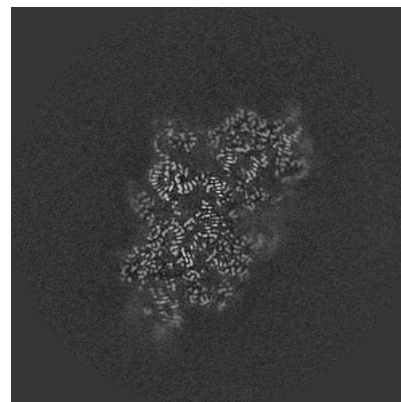
6.2.2 Raw map



X Index: 180



Y Index: 180

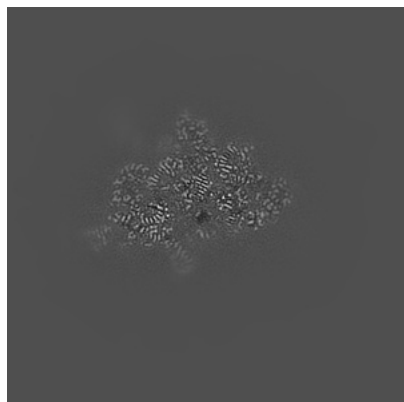


Z Index: 180

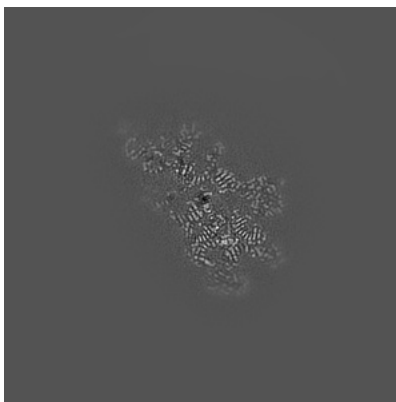
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

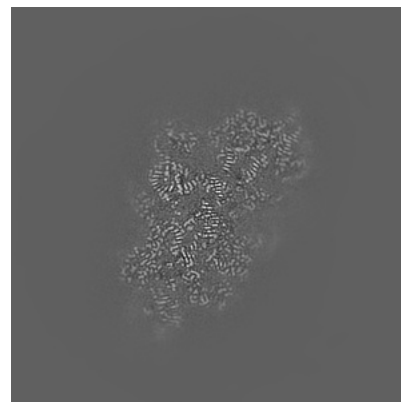
6.3.1 Primary map



X Index: 156

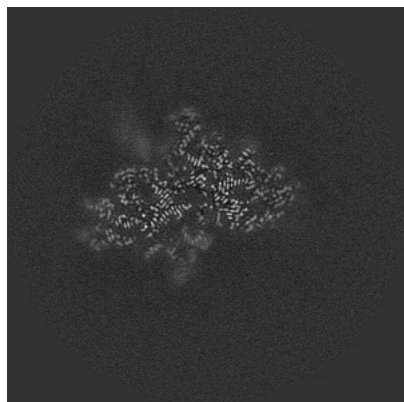


Y Index: 198

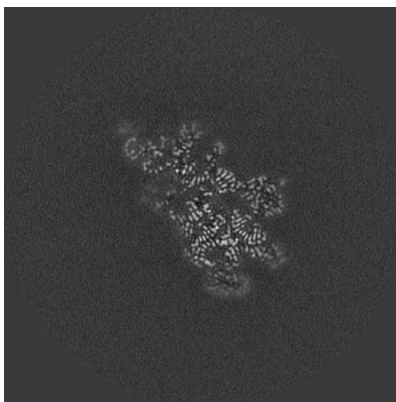


Z Index: 180

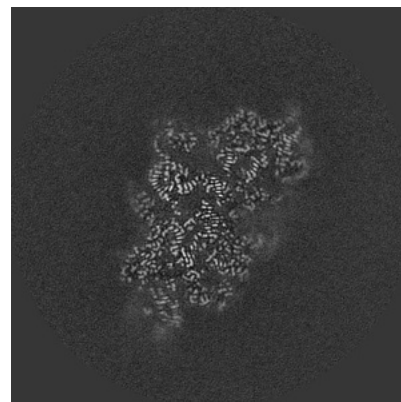
6.3.2 Raw map



X Index: 149



Y Index: 198

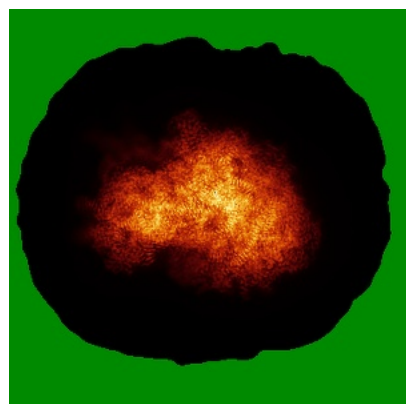


Z Index: 180

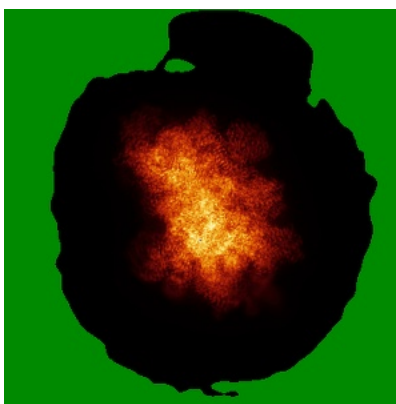
The images above show the largest variance slices of the map in three orthogonal directions.

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

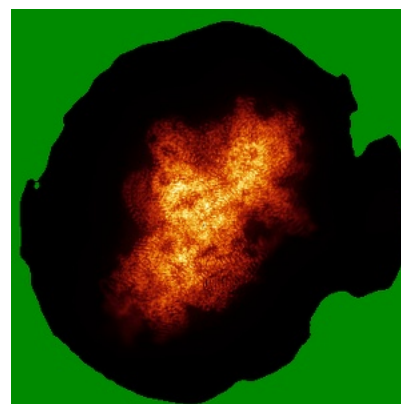
6.4.1 Primary map



X

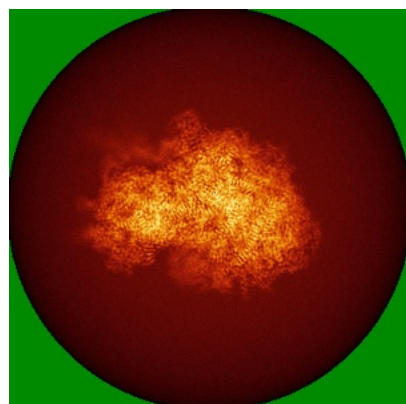


Y

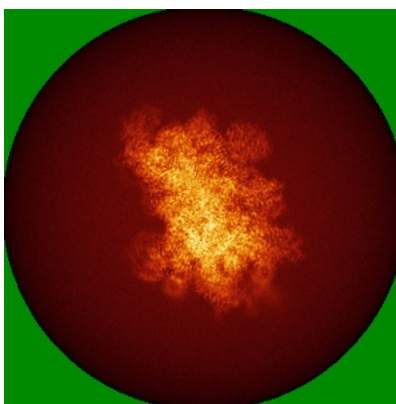


Z

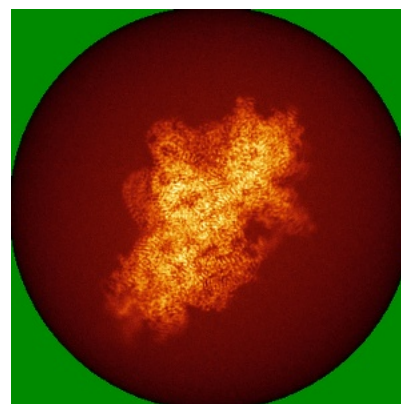
6.4.2 Raw map



X



Y

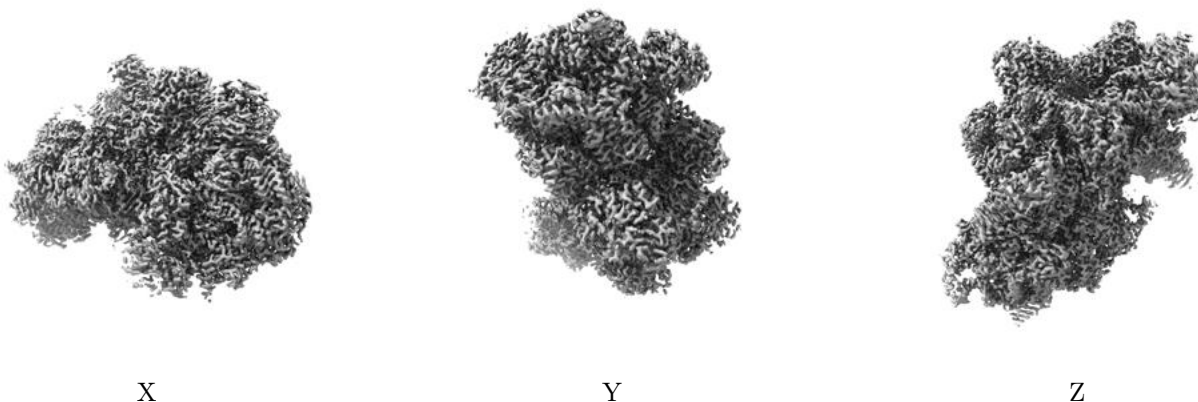


Z

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.

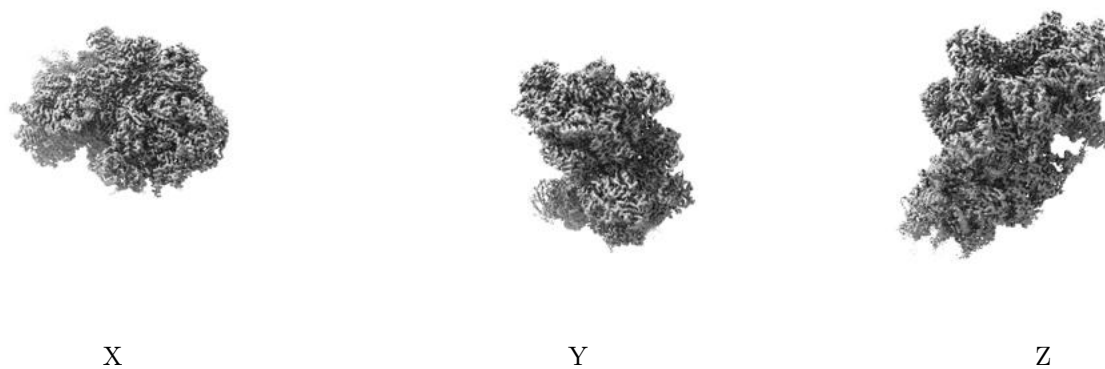
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.06. 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.

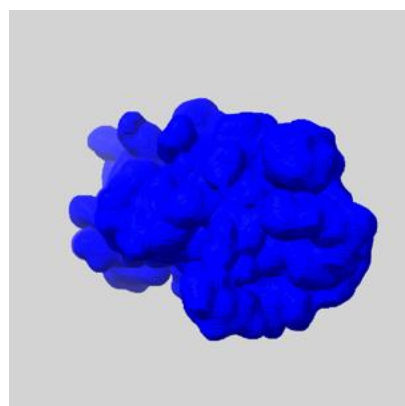
6.6 Mask visualisation [i](#)

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

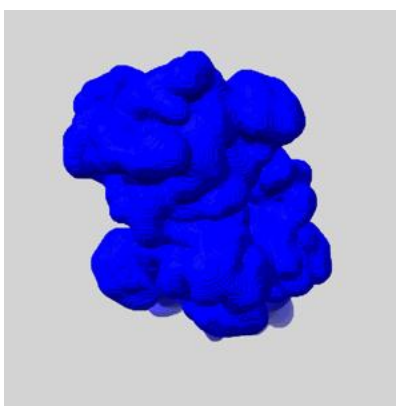
A mask typically either:

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

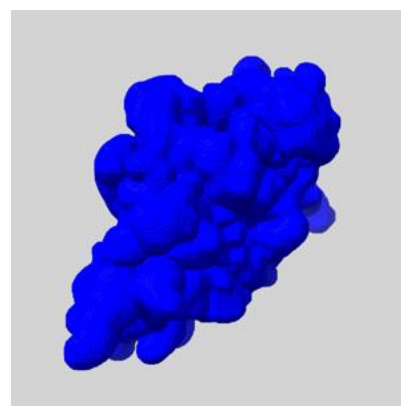
6.6.1 emd_11520_msk_1.map [i](#)



X



Y

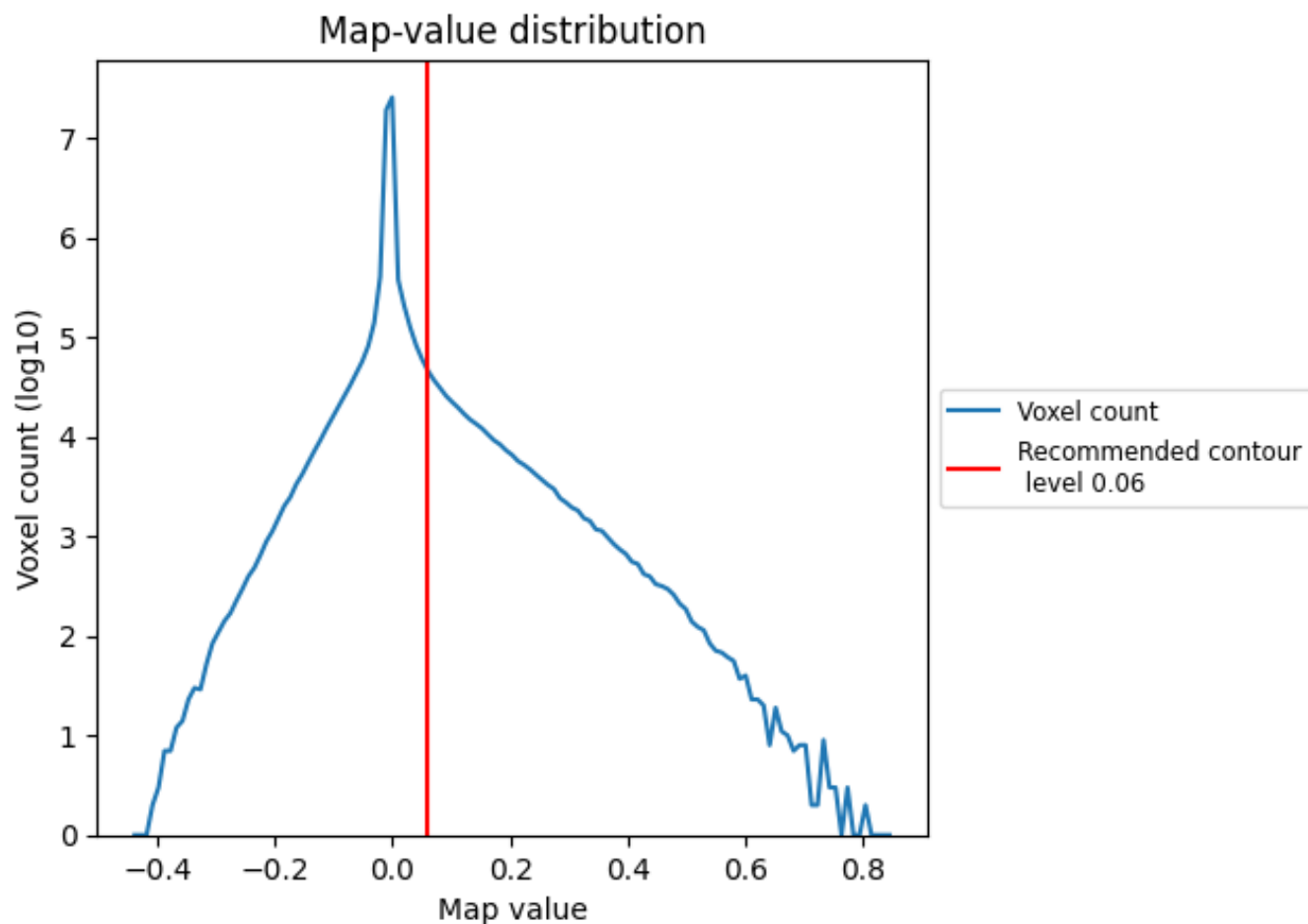


Z

7 Map analysis [i](#)

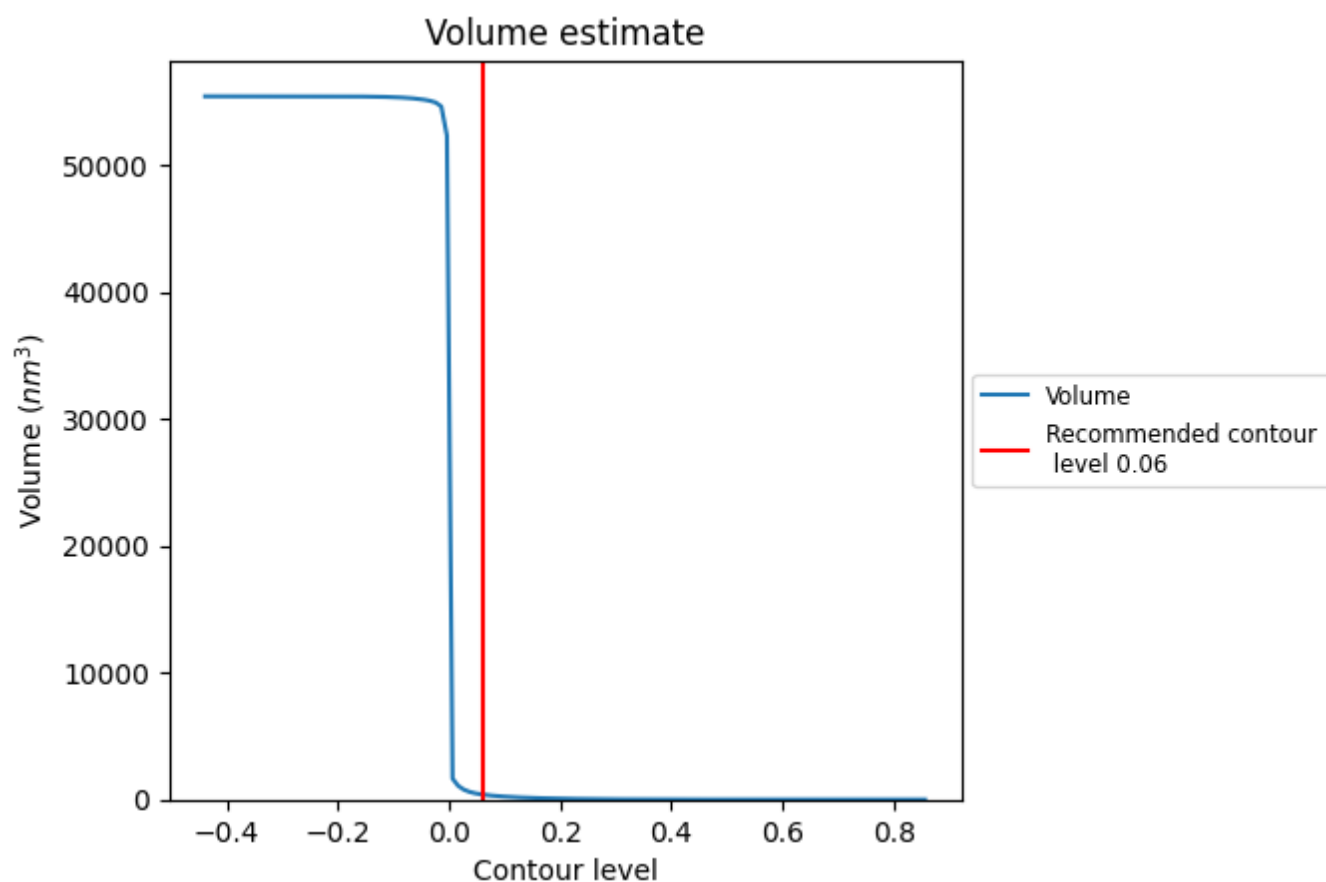
This section contains the results of statistical analysis of the map.

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.

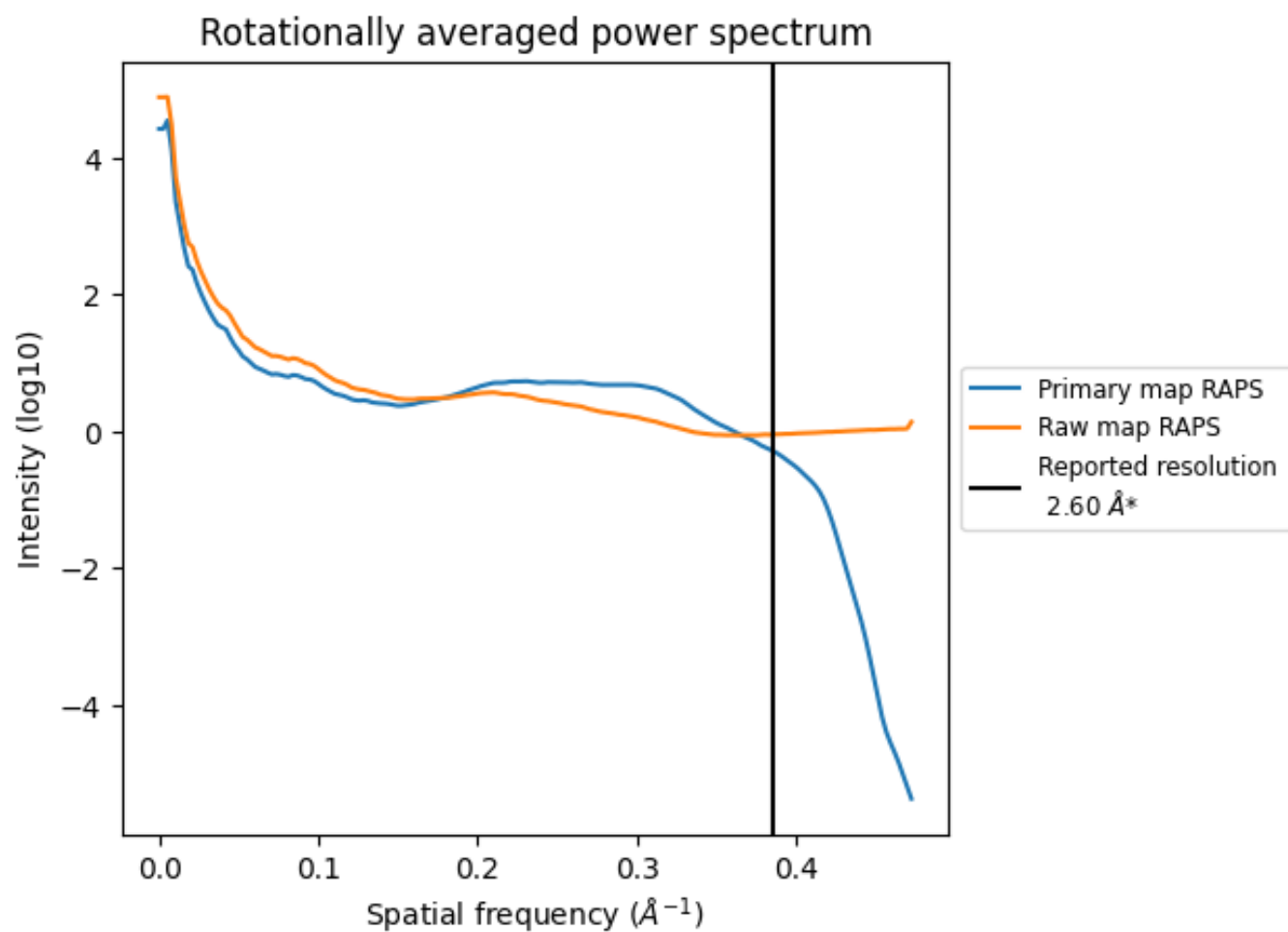
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 399 nm³; this corresponds to an approximate mass of 361 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 ⓘ

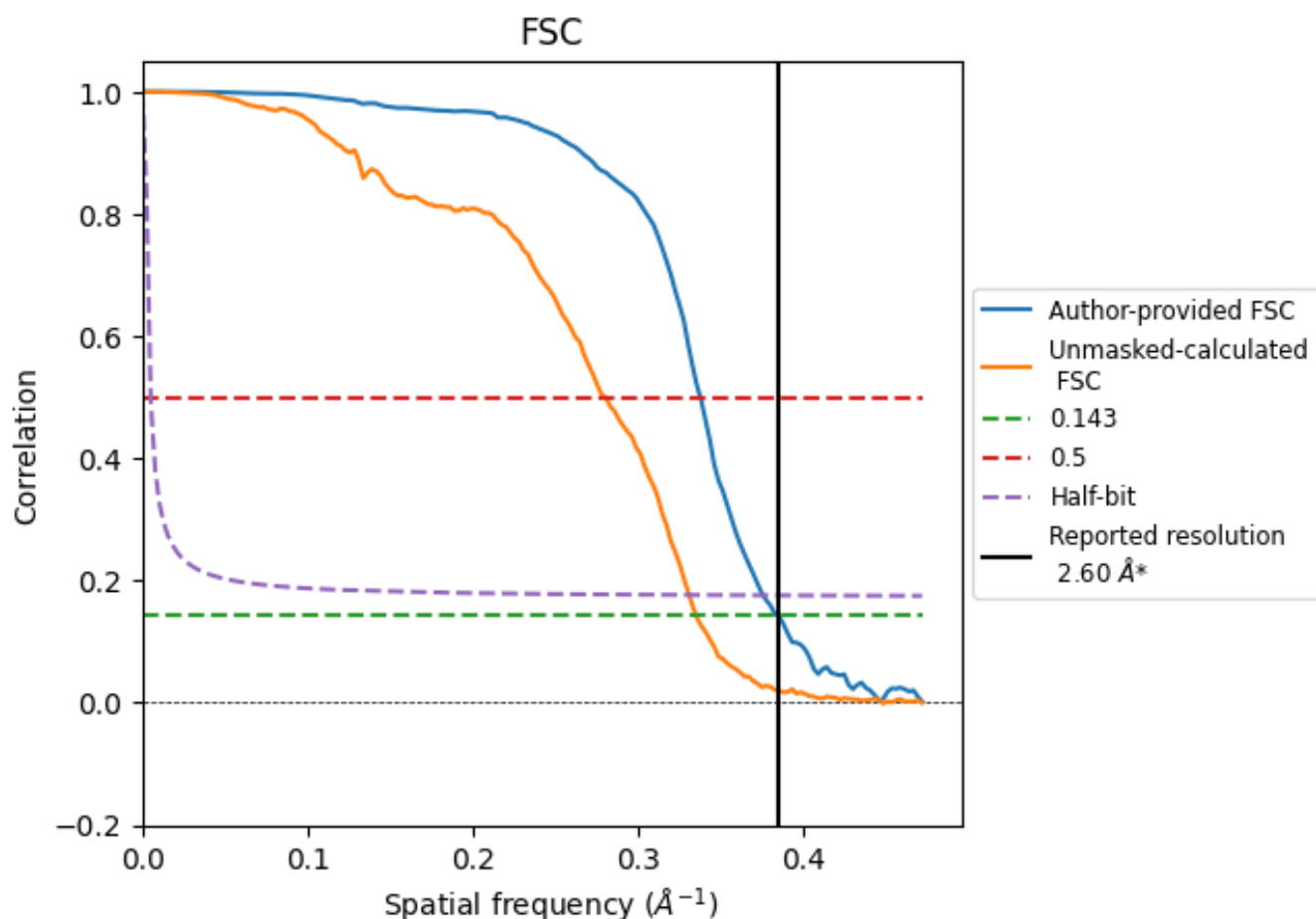


*Reported resolution corresponds to spatial frequency of 0.385 Å⁻¹

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 0.385 Å⁻¹

8.2 Resolution estimates [i](#)

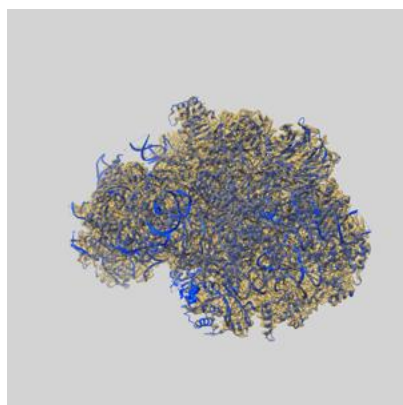
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.60	-	-
Author-provided FSC curve	2.60	2.96	2.66
Unmasked-calculated*	2.98	3.57	3.02

*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 2.98 differs from the reported value 2.6 by more than 10 %

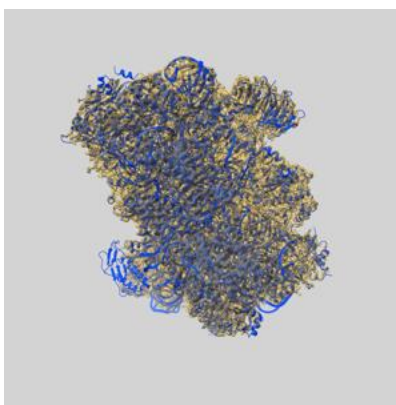
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-11520 and PDB model 6ZXG. Per-residue inclusion information can be found in section 3 on page 13.

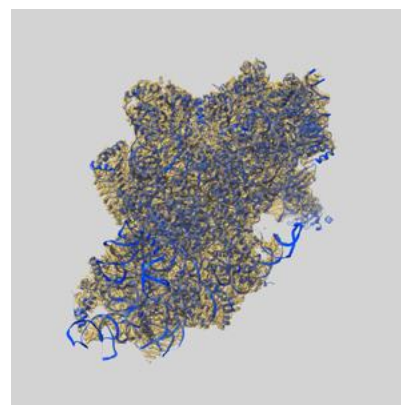
9.1 Map-model overlay [i](#)



X



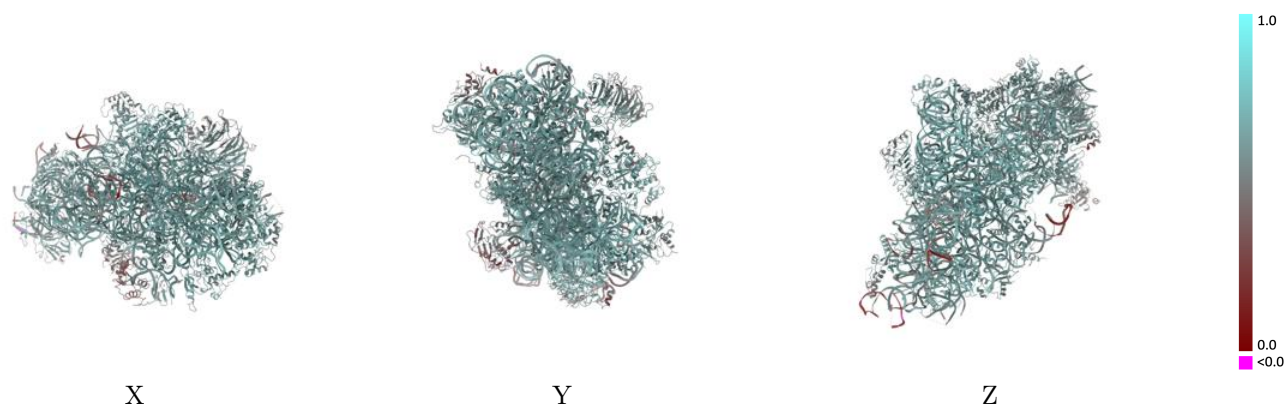
Y



Z

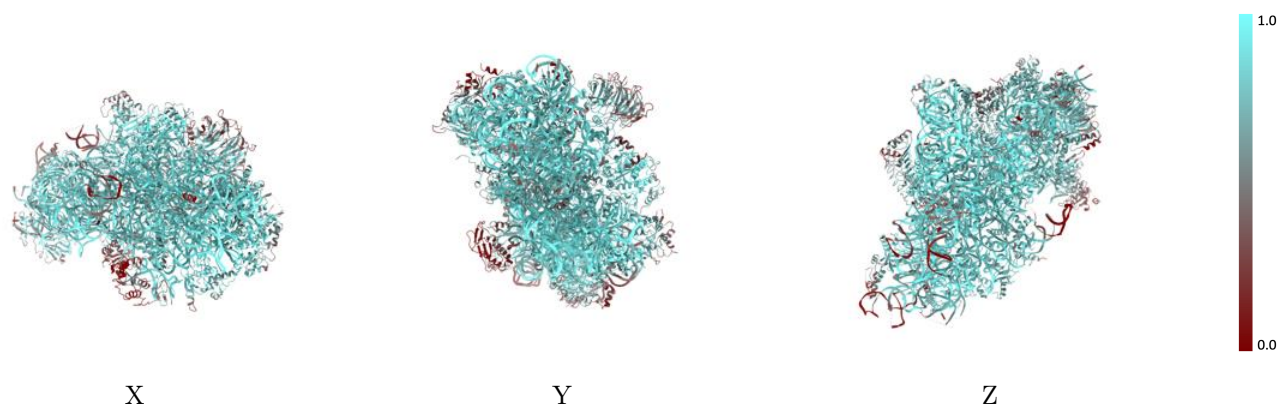
The images above show the 3D surface view of the map at the recommended contour level 0.06 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



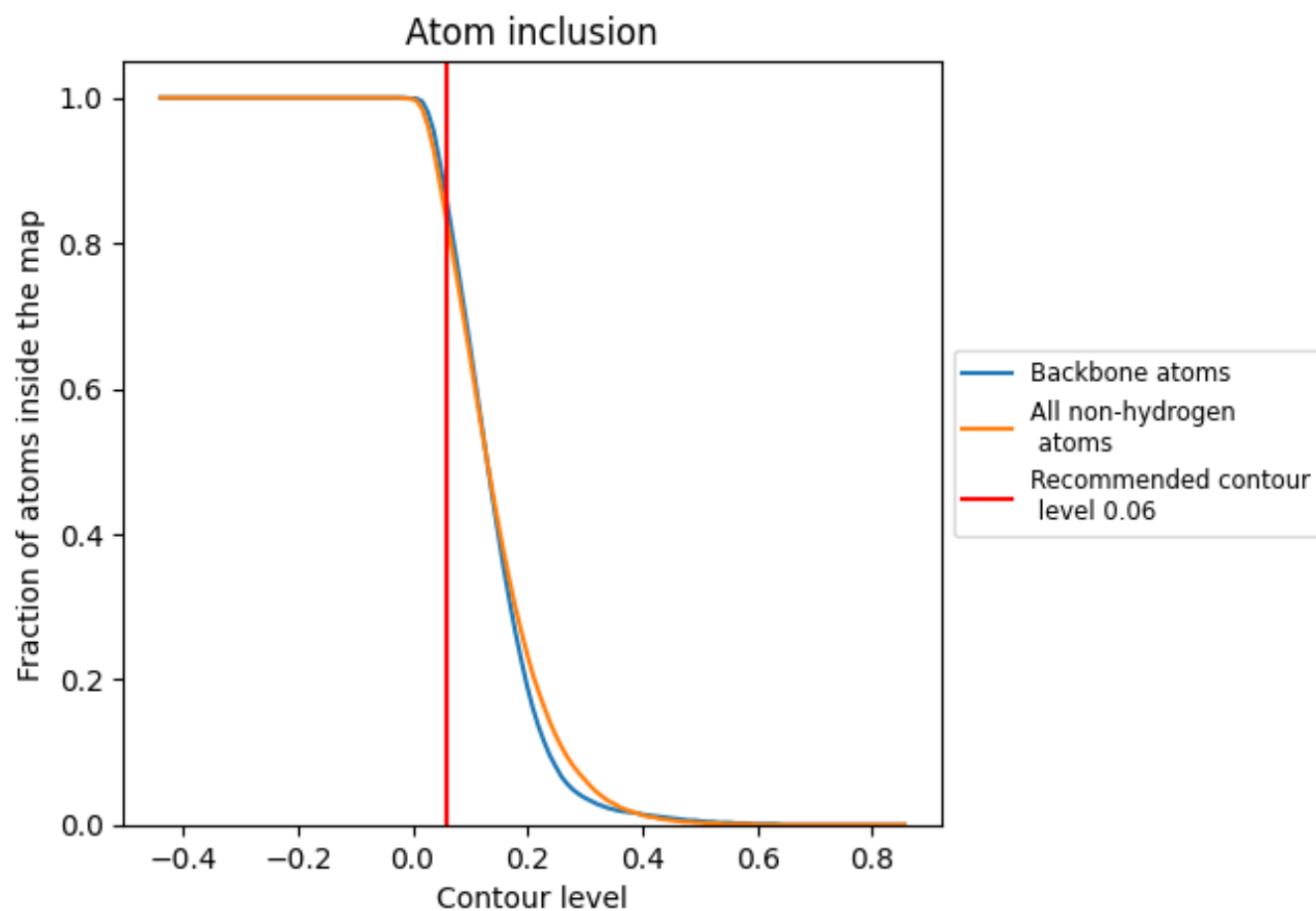
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

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



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.06).

9.4 Atom inclusion ⓘ



At the recommended contour level, 86% of all backbone atoms, 83% 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 (0.06) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8280	 0.6210
2	 0.9030	 0.6360
A	 0.8850	 0.6390
B	 0.8150	 0.6260
C	 0.9340	 0.6710
D	 0.7790	 0.6150
E	 0.8910	 0.6490
F	 0.7920	 0.6050
G	 0.6600	 0.5700
H	 0.6630	 0.5670
I	 0.7740	 0.6070
J	 0.8170	 0.6060
K	 0.7930	 0.6090
L	 0.8990	 0.6570
M	 0.1720	 0.4090
N	 0.8940	 0.6490
O	 0.8750	 0.6330
P	 0.7990	 0.6180
Q	 0.8630	 0.6330
R	 0.7390	 0.5990
S	 0.7900	 0.6170
T	 0.8550	 0.6420
U	 0.7010	 0.5890
V	 0.8860	 0.6530
W	 0.9560	 0.6800
X	 0.9330	 0.6760
Y	 0.8370	 0.6190
Z	 0.6460	 0.5780
b	 0.8080	 0.6330
c	 0.5940	 0.5520
d	 0.9470	 0.6770
e	 0.8120	 0.6300
f	 0.3920	 0.5110
g	 0.5860	 0.5510
h	 0.8910	 0.6580



Continued on next page...

Continued from previous page...

Chain	Atom inclusion	Q-score
j	 0.7610	 0.6040
k	 0.5120	 0.5310
z	 0.7770	 0.6120