



Full wwPDB EM Validation Report ⓘ

Mar 23, 2026 – 02:09 AM UTC

PDB ID : 8CSP / pdb_00008csp
EMDB ID : EMD-26966
Title : Human mitochondrial small subunit assembly intermediate (State A)
Authors : Harper, N.J.; Burnside, C.; Klinge, S.
Deposited on : 2022-05-13
Resolution : 2.66 Å(reported)
Based on initial models : 2C2N, 6RW4, 6AAX

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

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

A user guide is available at

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

The types of validation reports are described at

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

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

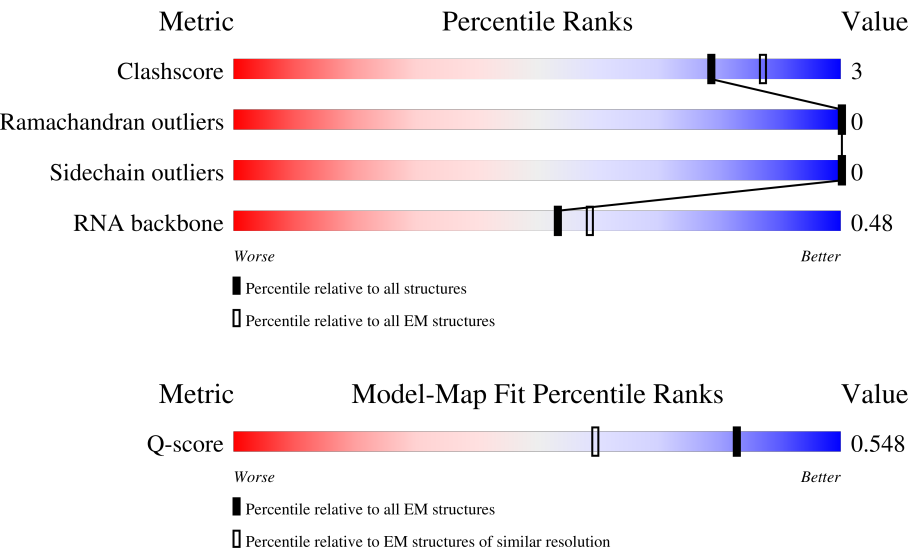
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

1 Overall quality at a glance i

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

The reported resolution of this entry is 2.66 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	9119 (2.16 - 3.16)

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

Mol	Chain	Length	Quality of chain
1	0	218	13% 90% 6% .
2	1	323	10% 77% 7% 16%
3	4	689	28% 72% 10% 18%

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Mol	Chain	Length	Quality of chain
4	5	346	
5	7	456	
6	8	390	
7	9	698	
8	A	955	
9	B	296	
10	C	167	
11	D	430	
12	E	125	
13	F	242	
14	G	396	
15	H	201	
16	J	138	
17	K	128	
18	L	257	
19	M	137	
20	N	130	
21	O	258	
22	P	142	
23	Q	87	
24	R	360	
25	S	190	
26	T	173	
27	U	205	
28	V	414	

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Mol	Chain	Length	Quality of chain
29	W	187	5%50%48%
30	X	398	82%6%12%
31	Y	395	5%28%69%
32	Z	106	14%45%8%47%
33	a	437	32%55%5%40%

2 Entry composition

There are 41 unique types of molecules in this entry. The entry contains 69302 atoms, of which 0 are hydrogens and 0 are deuteriums.

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

- Molecule 1 is a protein called 28S ribosomal protein S34, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	0	211	Total	C	N	O	S	0	0
			1754	1108	333	308	5		

- Molecule 2 is a protein called 28S ribosomal protein S35, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	1	271	Total	C	N	O	S	0	0
			2195	1393	369	422	11		

- Molecule 3 is a protein called Pentatricopeptide repeat domain-containing protein 3, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	4	566	Total	C	N	O	S	0	0
			4585	2940	774	843	28		

- Molecule 4 is a protein called Dimethyladenosine transferase 1, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	5	319	Total	C	N	O	S	0	0
			2568	1648	458	451	11		

- Molecule 5 is a protein called Methyltransferase-like protein 17, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	7	398	Total	C	N	O	S	0	0
			3145	2006	579	544	16		

- Molecule 6 is a protein called Malonyl-CoA-acyl carrier protein transacylase, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	8	326	Total	C	N	O	S	0	0
			2543	1617	463	446	17		

- Molecule 7 is a protein called Nitric oxide-associated protein 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	9	446	Total	C	N	O	S	0	0
			3558	2271	637	637	13		

- Molecule 8 is a RNA chain called 12S mitochondrial rRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	A	737	Total	C	N	O	P	0	0
			15655	7018	2818	5082	737		

- Molecule 9 is a protein called 28S ribosomal protein S2, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	B	220	Total	C	N	O	S	0	0
			1789	1142	324	313	10		

- Molecule 10 is a protein called 28S ribosomal protein S24, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	C	126	Total	C	N	O	S	0	0
			1042	679	181	177	5		

- Molecule 11 is a protein called 28S ribosomal protein S5, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
11	D	232	Total	C	N	O	S	0	0
			1838	1155	345	329	9		

- Molecule 12 is a protein called 28S ribosomal protein S6, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
12	E	105	Total	C	N	O	S	0	0
			839	526	154	155	4		

- Molecule 13 is a protein called 28S ribosomal protein S7, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
13	F	208	Total	C	N	O	S	0	0
			1724	1104	312	297	11		

- Molecule 14 is a protein called 28S ribosomal protein S9, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
14	G	292	Total	C	N	O	S	0	0
			2395	1522	418	441	14		

- Molecule 15 is a protein called 28S ribosomal protein S10, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
15	H	128	Total	C	N	O	S	0	0
			1045	675	172	195	3		

- Molecule 16 is a protein called 28S ribosomal protein S12, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
16	J	108	Total	C	N	O	S	0	0
			839	521	169	143	6		

- Molecule 17 is a protein called 28S ribosomal protein S14, mitochondrial.

Mol	Chain	Residues	Atoms				AltConf	Trace
17	K	33	Total	C	N	O	0	0
			283	173	66	44		

- Molecule 18 is a protein called 28S ribosomal protein S15, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
18	L	131	Total	C	N	O	S	0	0
			1091	698	190	197	6		

- Molecule 19 is a protein called 28S ribosomal protein S16, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
19	M	115	Total	C	N	O	S	0	0
			913	578	181	148	6		

- Molecule 20 is a protein called 28S ribosomal protein S17, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
20	N	109	Total	C	N	O	S	0	0
			859	557	155	144	3		

- Molecule 21 is a protein called 28S ribosomal protein S18b, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
21	O	190	Total	C	N	O	S	0	0
			1570	999	291	274	6		

- Molecule 22 is a protein called 28S ribosomal protein S18c, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
22	P	92	Total	C	N	O	S	0	0
			731	472	123	128	8		

- Molecule 23 is a protein called 28S ribosomal protein S21, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
23	Q	6	Total	C	N	O	S	0	0
			49	30	8	10	1		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
Q	50	ARG	CYS	conflict	UNP P82921

- Molecule 24 is a protein called 28S ribosomal protein S22, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
24	R	291	Total	C	N	O	S	0	0
			2382	1518	409	447	8		

- Molecule 25 is a protein called 28S ribosomal protein S23, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
25	S	133	Total	C	N	O	S	0	0
			1100	709	196	194	1		

- Molecule 26 is a protein called 28S ribosomal protein S25, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
26	T	164	Total	C	N	O	S	0	0
			1344	859	234	240	11		

- Molecule 27 is a protein called 28S ribosomal protein S26, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
27	U	174	Total	C	N	O	S	0	0
			1468	905	295	264	4		

- Molecule 28 is a protein called 28S ribosomal protein S27, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
28	V	359	Total	C	N	O	S	0	0
			2946	1891	491	552	12		

- Molecule 29 is a protein called 28S ribosomal protein S28, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
29	W	98	Total	C	N	O	S	0	0
			775	491	138	142	4		

- Molecule 30 is a protein called 28S ribosomal protein S29, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
30	X	352	Total	C	N	O	S	0	0
			2849	1822	499	517	11		

- Molecule 31 is a protein called 28S ribosomal protein S31, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
31	Y	123	Total	C	N	O	S	0	0
			1043	676	175	189	3		

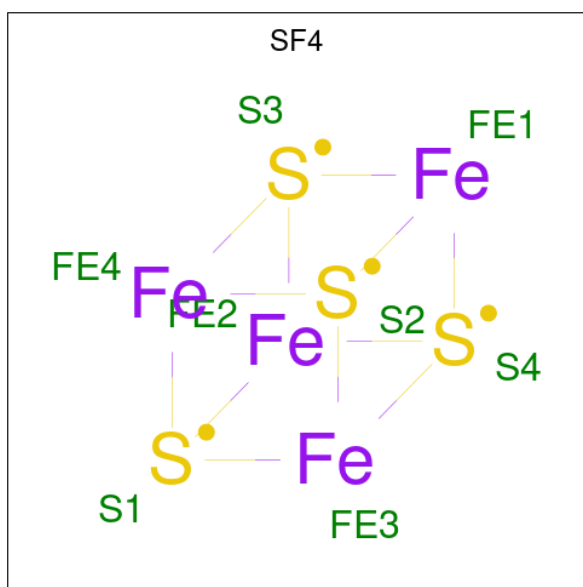
- Molecule 32 is a protein called 28S ribosomal protein S33, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
32	Z	56	Total	C	N	O	S	0	0
			465	298	84	80	3		

- Molecule 33 is a protein called GTPase Era, mitochondrial.

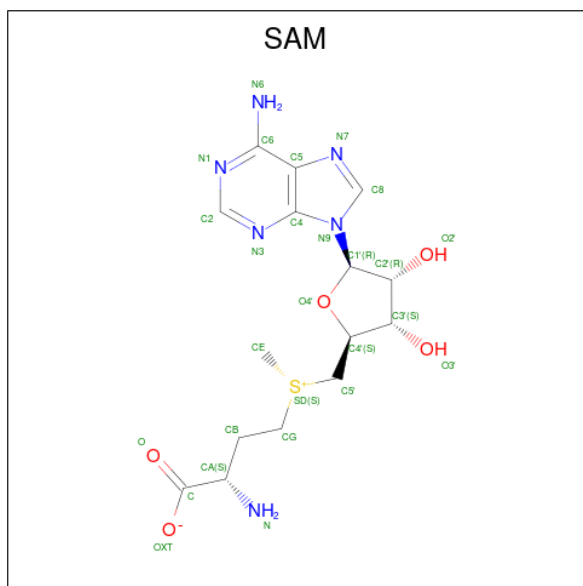
Mol	Chain	Residues	Atoms					AltConf	Trace
33	a	263	Total	C	N	O	S	0	0
			1785	1121	323	335	6		

- Molecule 34 is IRON/SULFUR CLUSTER (CCD ID: SF4) (formula: Fe₄S₄) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			AltConf
34	7	1	Total	Fe	S	0
			8	4	4	

- Molecule 35 is S-ADENOSYLMETHIONINE (CCD ID: SAM) (formula: $C_{15}H_{22}N_6O_5S$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					AltConf
35	7	1	Total	C	N	O	S	0
			27	15	6	5	1	

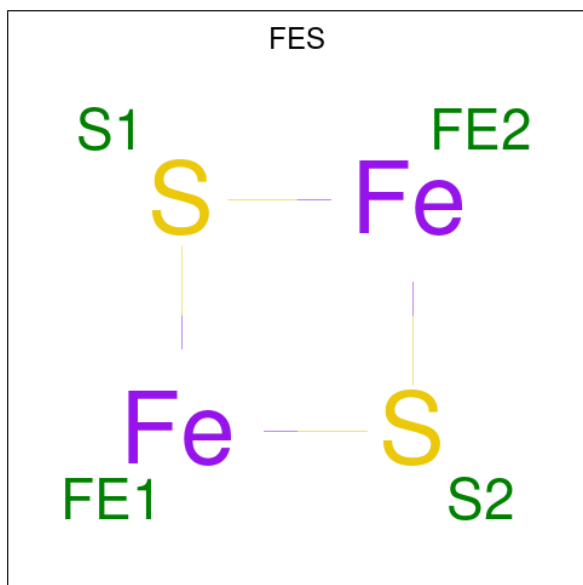
- Molecule 36 is POTASSIUM ION (CCD ID: K) (formula: K).

Mol	Chain	Residues	Atoms		AltConf
36	A	4	Total	K	0
			4	4	

- Molecule 37 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		AltConf
37	A	26	Total	Mg	0
			26	26	
37	B	1	Total	Mg	0
			1	1	
37	X	1	Total	Mg	0
			1	1	

- Molecule 38 is FE2/S2 (INORGANIC) CLUSTER (CCD ID: FES) (formula: Fe₂S₂).

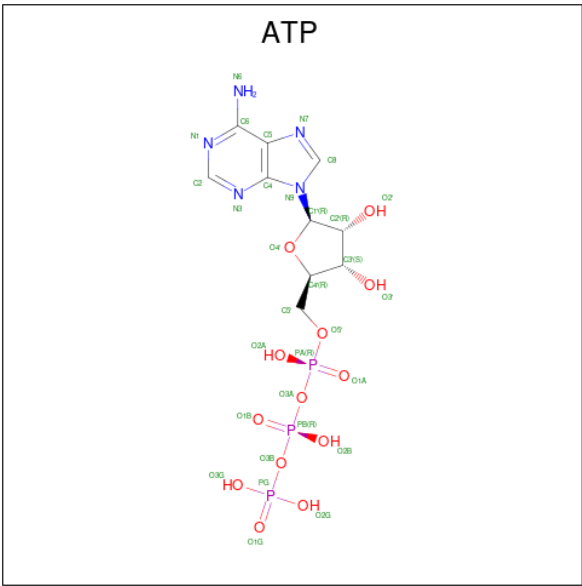


Mol	Chain	Residues	Atoms			AltConf
38	E	1	Total	Fe	S	0
			4	2	2	
38	T	1	Total	Fe	S	0
			4	2	2	

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

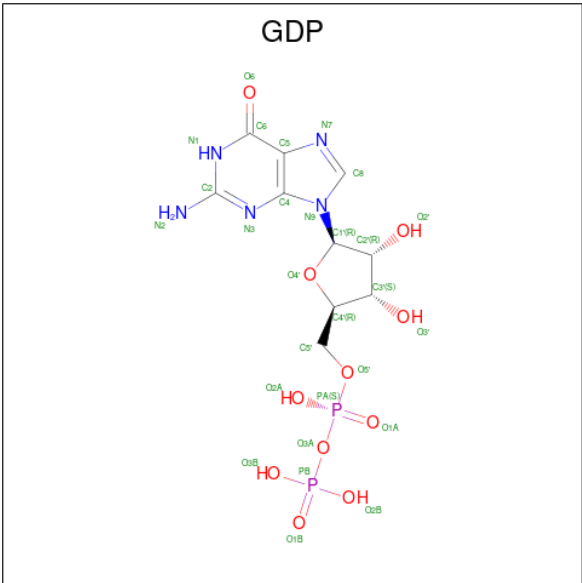
Mol	Chain	Residues	Atoms		AltConf
39	O	1	Total	Zn	0
			1	1	

- Molecule 40 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: ATP) (formula: C₁₀H₁₆N₅O₁₃P₃).



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

- Molecule 41 is GUANOSINE-5'-DIPHOSPHATE (CCD ID: GDP) (formula: C₁₀H₁₅N₅O₁₁P₂).

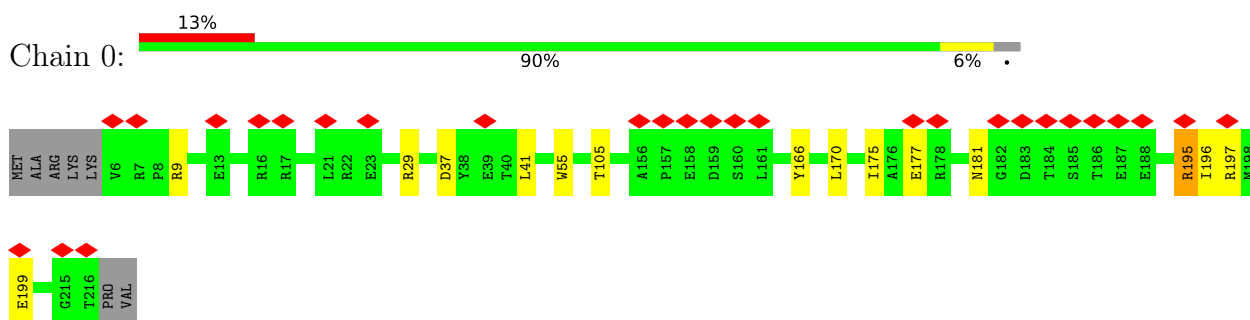


Mol	Chain	Residues	Atoms					AltConf
41	X	1	Total	C	N	O	P	0
			28	10	5	11	2	

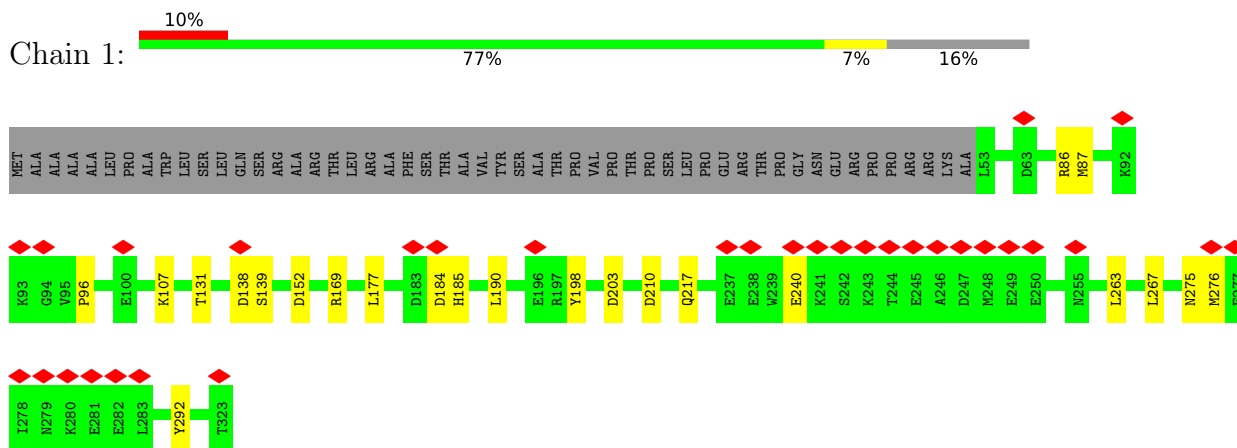
3 Residue-property plots

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

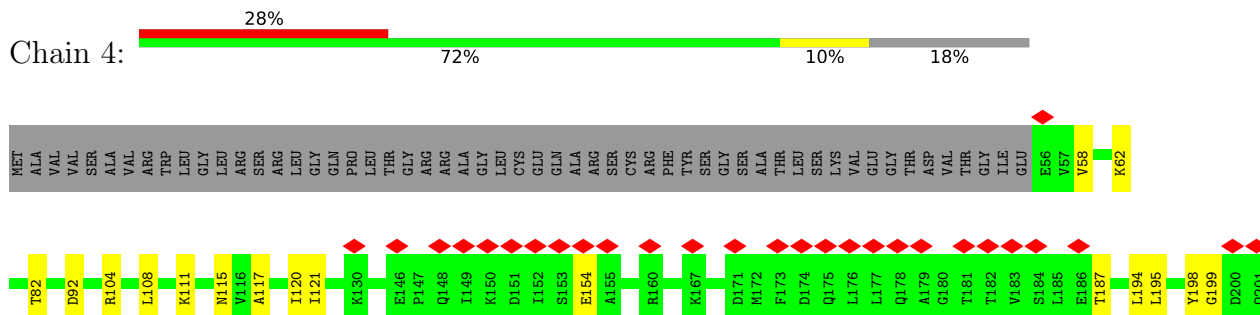
- Molecule 1: 28S ribosomal protein S34, mitochondrial

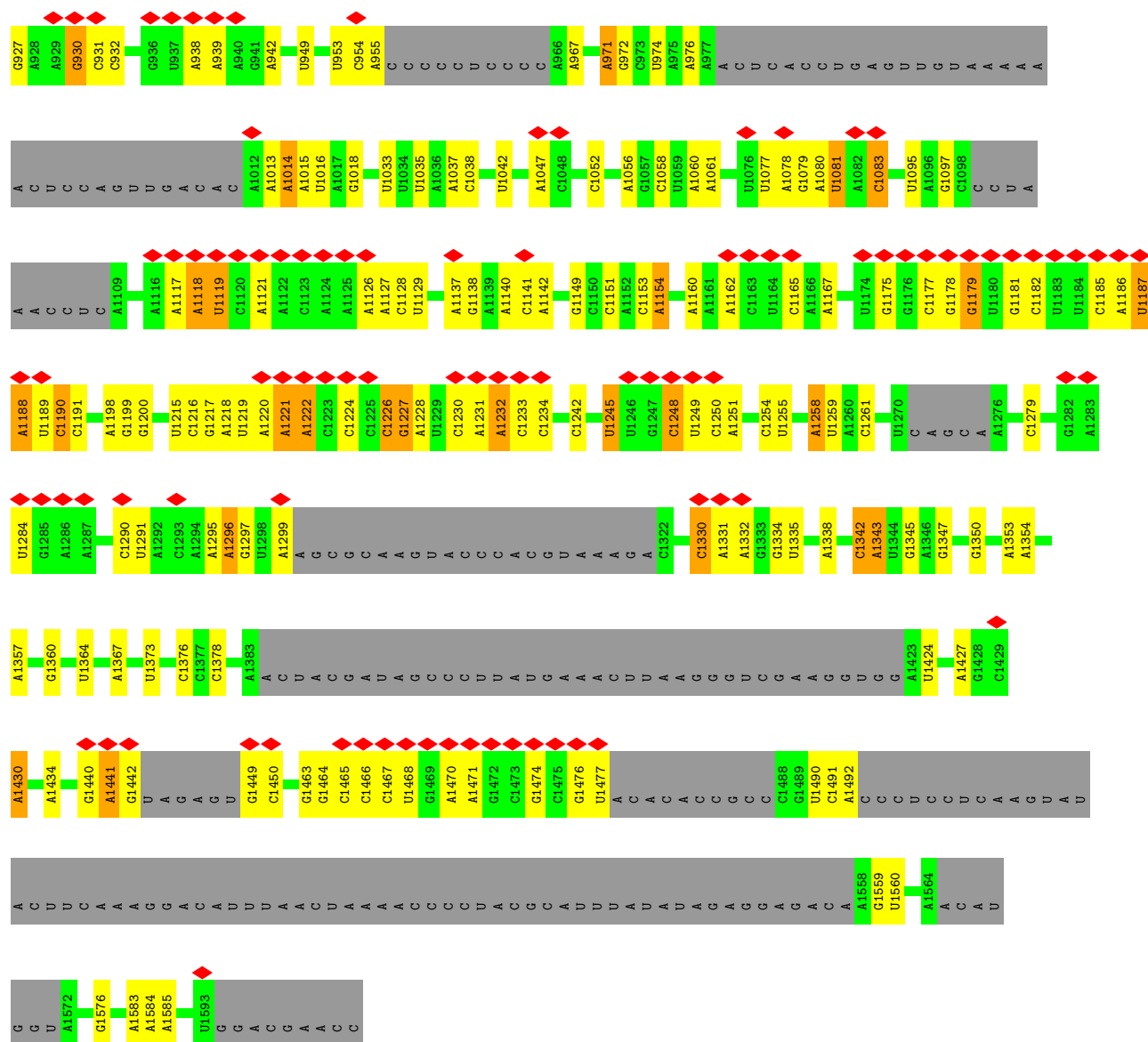


- Molecule 2: 28S ribosomal protein S35, mitochondrial



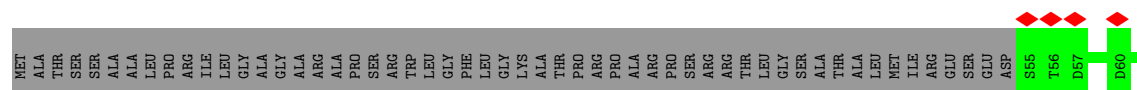
- Molecule 3: Pentatricopeptide repeat domain-containing protein 3, mitochondrial





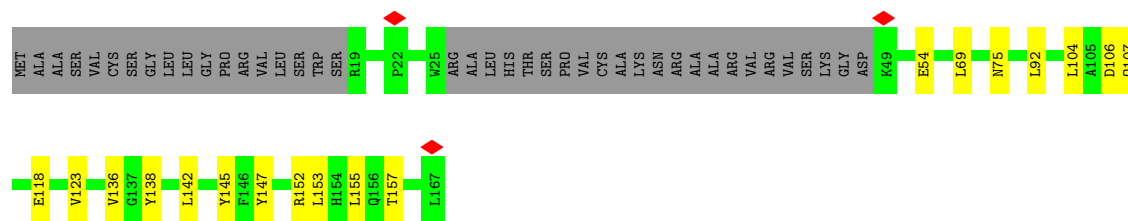
- Molecule 9: 28S ribosomal protein S2, mitochondrial

Chain B: 70% 26%

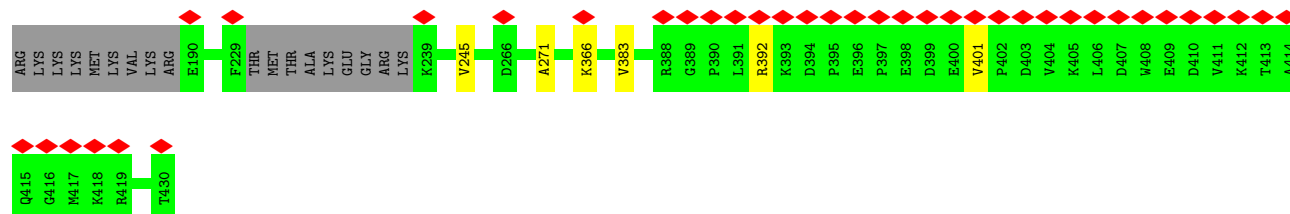
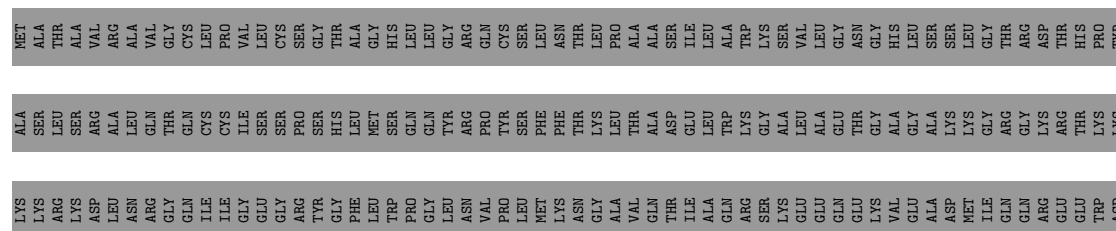


- Molecule 10: 28S ribosomal protein S24, mitochondrial

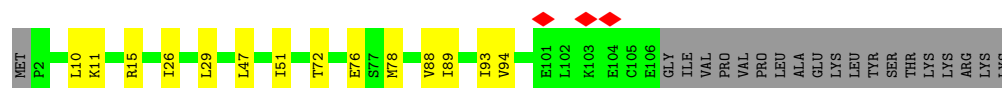
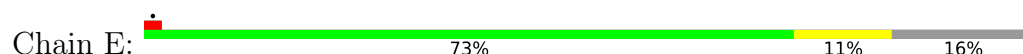
Chain C: 65% 11% 25%



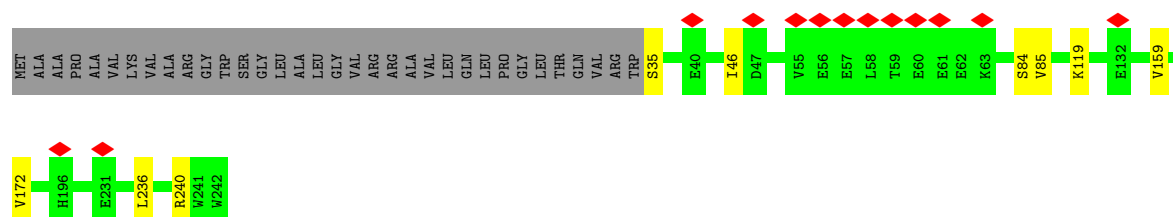
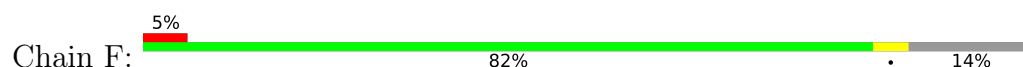
- Molecule 11: 28S ribosomal protein S5, mitochondrial



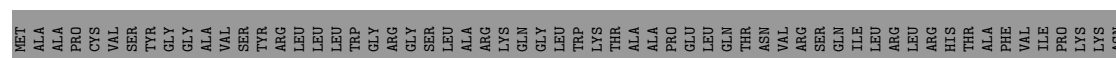
- Molecule 12: 28S ribosomal protein S6, mitochondrial

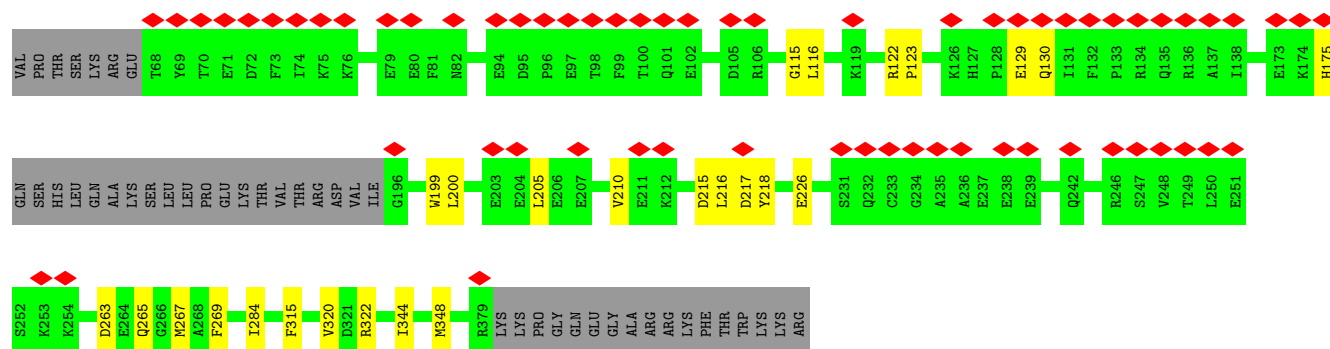


- Molecule 13: 28S ribosomal protein S7, mitochondrial



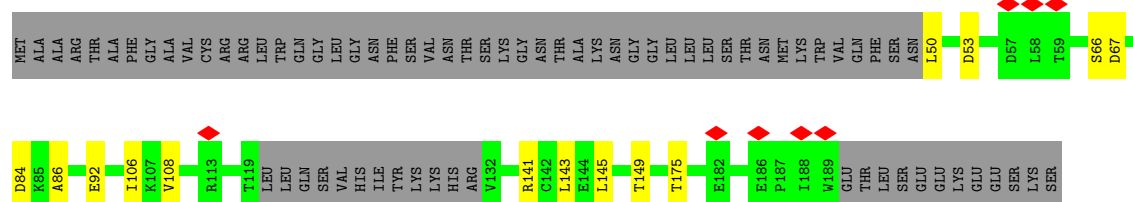
- Molecule 14: 28S ribosomal protein S9, mitochondrial





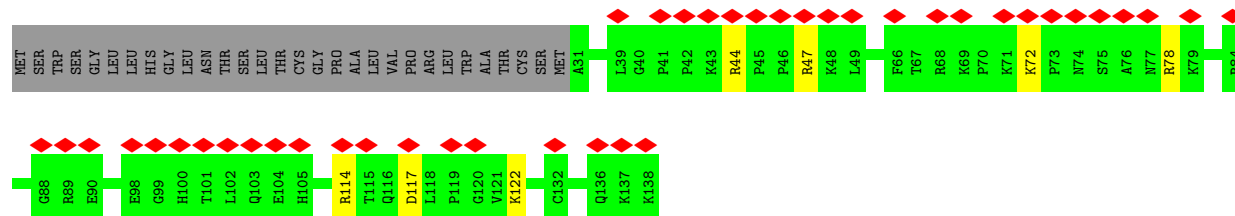
- Molecule 15: 28S ribosomal protein S10, mitochondrial

Chain H: 57% 7% 36%



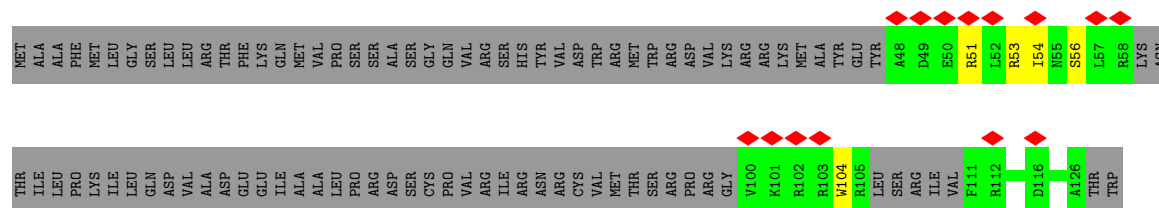
- Molecule 16: 28S ribosomal protein S12, mitochondrial

Chain J: 30% 73% 5% 22%



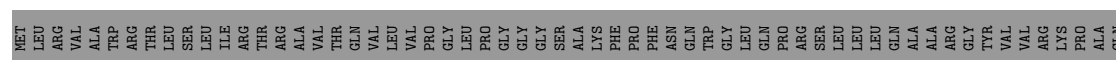
- Molecule 17: 28S ribosomal protein S14, mitochondrial

Chain K: 11% 22% 74%

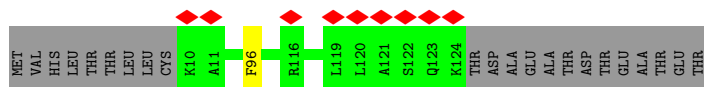
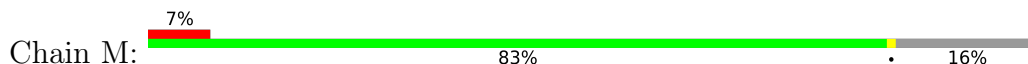


- Molecule 18: 28S ribosomal protein S15, mitochondrial

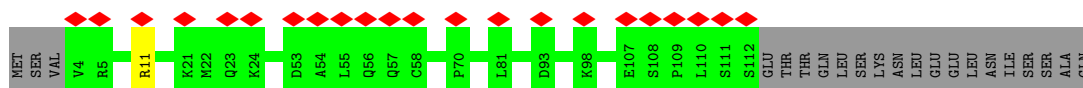
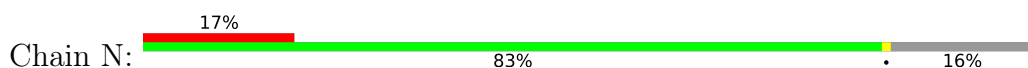
Chain L: 45% 6% 49%



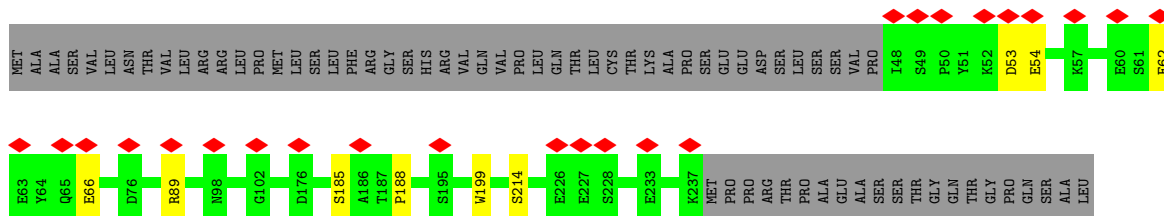
- Molecule 19: 28S ribosomal protein S16, mitochondrial



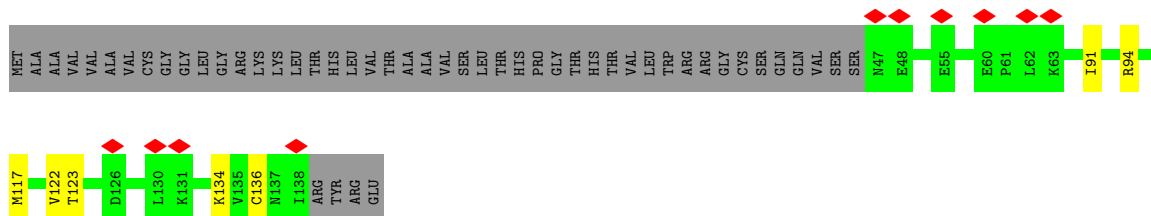
- Molecule 20: 28S ribosomal protein S17, mitochondrial



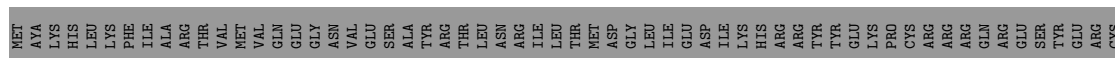
- Molecule 21: 28S ribosomal protein S18b, mitochondrial

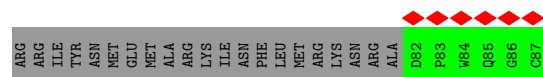


- Molecule 22: 28S ribosomal protein S18c, mitochondrial

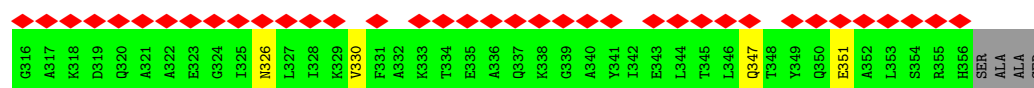
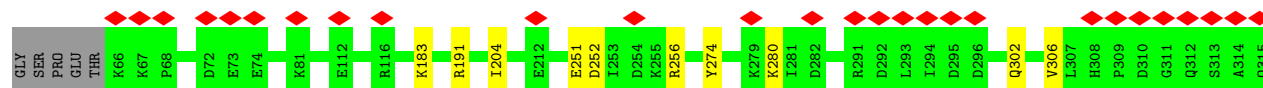
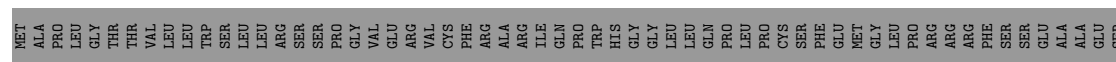
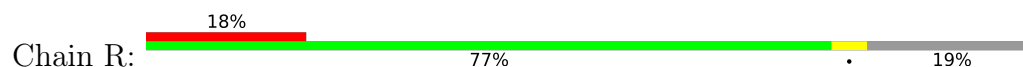


- Molecule 23: 28S ribosomal protein S21, mitochondrial

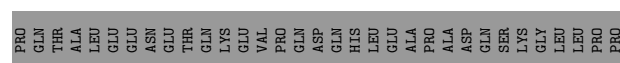
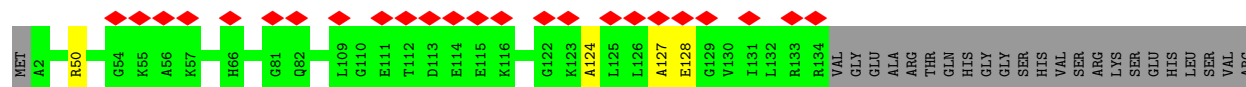




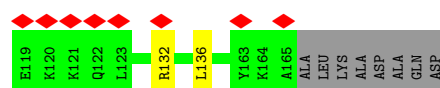
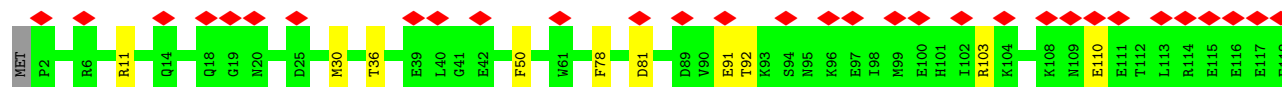
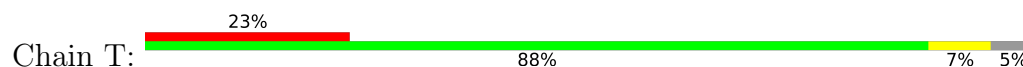
- Molecule 24: 28S ribosomal protein S22, mitochondrial



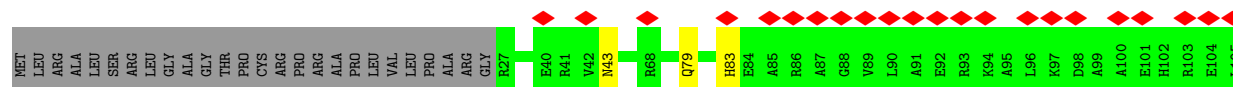
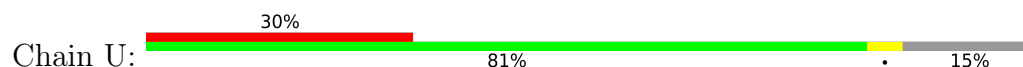
- Molecule 25: 28S ribosomal protein S23, mitochondrial



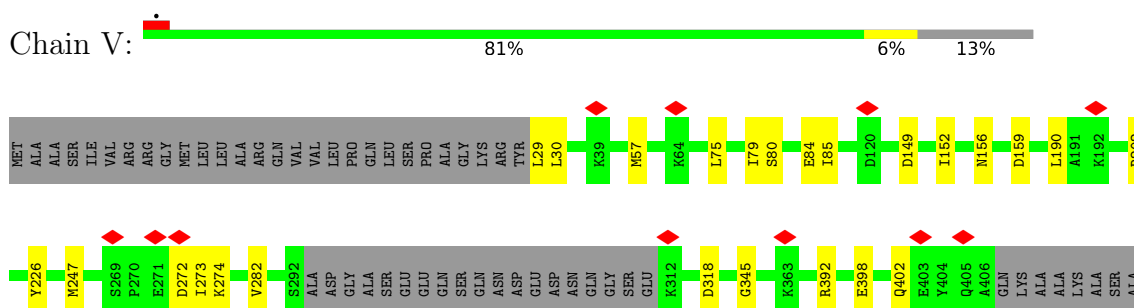
- Molecule 26: 28S ribosomal protein S25, mitochondrial



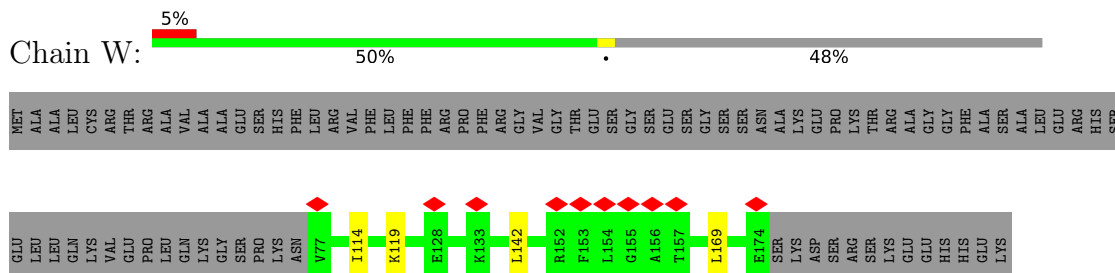
- Molecule 27: 28S ribosomal protein S26, mitochondrial



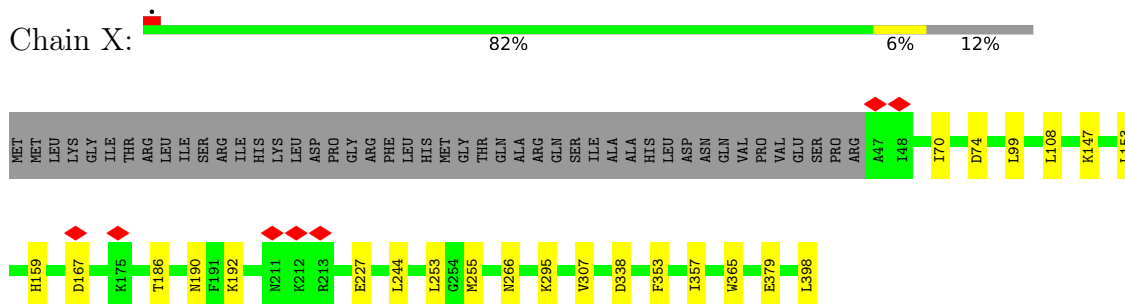
- Molecule 28: 28S ribosomal protein S27, mitochondrial



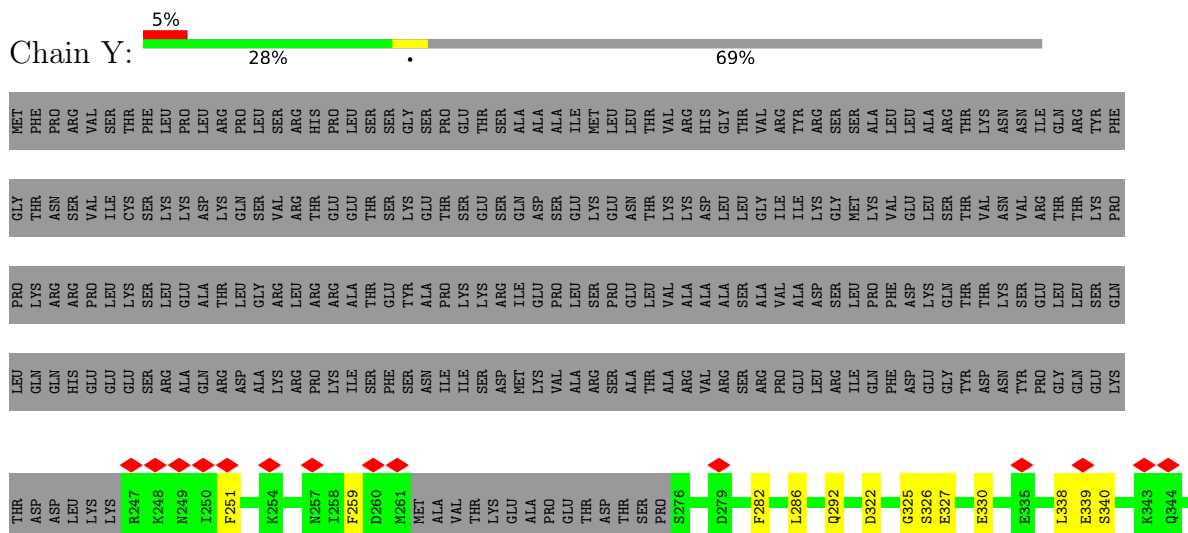
- Molecule 29: 28S ribosomal protein S28, mitochondrial

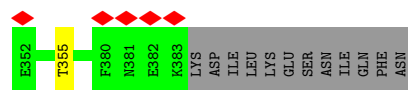


- Molecule 30: 28S ribosomal protein S29, mitochondrial

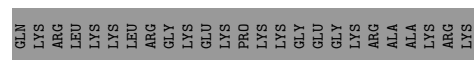
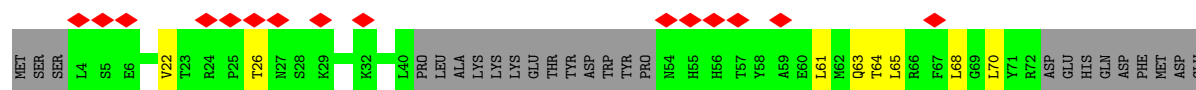
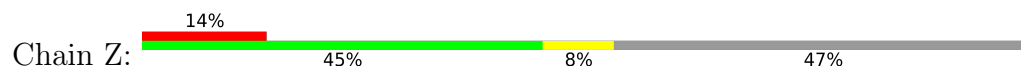


- Molecule 31: 28S ribosomal protein S31, mitochondrial

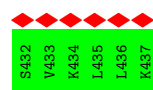
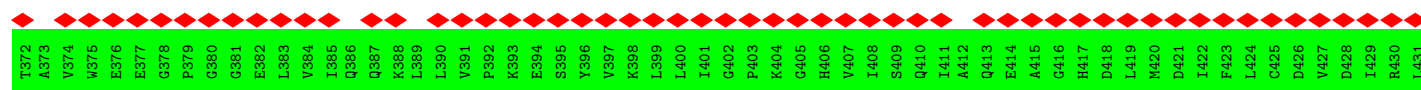
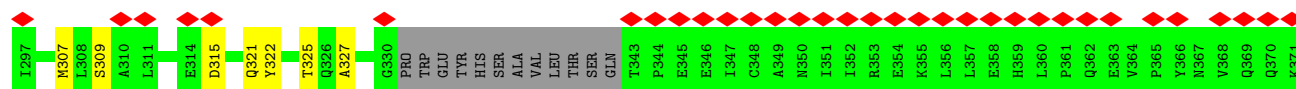
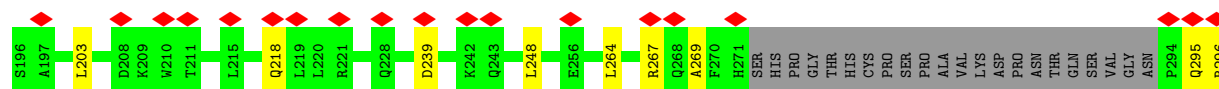
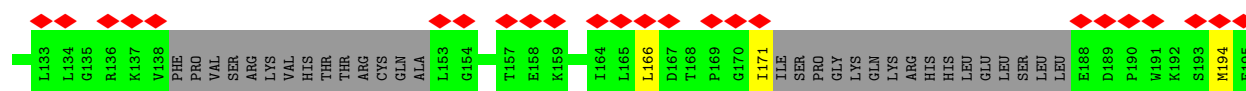
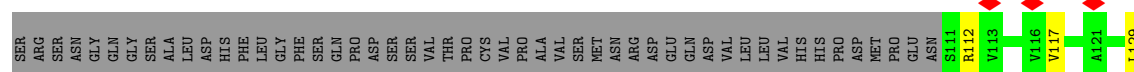




- Molecule 32: 28S ribosomal protein S33, mitochondrial



- Molecule 33: GTPase Era, mitochondrial



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	263184	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 ($e^-/\text{\AA}^2$)	51	Depositor
Minimum defocus (nm)	700	Depositor
Maximum defocus (nm)	1800	Depositor
Magnification	64000	Depositor
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	66.216	Depositor
Minimum map value	-26.873	Depositor
Average map value	0.045	Depositor
Map value standard deviation	1.467	Depositor
Recommended contour level	8.5	Depositor
Map size (Å)	424.80002, 424.80002, 424.80002	wwPDB
Map dimensions	400, 400, 400	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.062, 1.062, 1.062	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SAM, SF4, GDP, ATP, FES, K, ZN, MG

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	0	0.17	0/1800	0.36	0/2440
2	1	0.18	0/2241	0.36	0/3033
3	4	0.18	0/4686	0.38	0/6335
4	5	0.19	0/2624	0.40	0/3552
5	7	0.19	0/3234	0.37	0/4390
6	8	0.19	0/2603	0.40	0/3520
7	9	0.19	0/3637	0.41	0/4935
8	A	0.15	0/17503	0.30	0/27229
9	B	0.18	0/1832	0.35	0/2480
10	C	0.19	0/1074	0.40	0/1456
11	D	0.15	0/1877	0.34	0/2523
12	E	0.19	0/853	0.38	0/1151
13	F	0.16	0/1766	0.34	0/2373
14	G	0.18	0/2446	0.35	0/3283
15	H	0.17	0/1067	0.35	0/1448
16	J	0.13	0/855	0.33	0/1148
17	K	0.14	0/285	0.31	0/375
18	L	0.17	0/1111	0.37	0/1490
19	M	0.15	0/934	0.33	0/1255
20	N	0.14	0/877	0.31	0/1187
21	O	0.17	0/1624	0.34	0/2210
22	P	0.16	0/747	0.33	0/1004
23	Q	0.12	0/51	0.29	0/68
24	R	0.14	0/2429	0.30	0/3280
25	S	0.16	0/1127	0.34	0/1518
26	T	0.15	0/1375	0.34	0/1847
27	U	0.18	0/1490	0.36	0/1999
28	V	0.18	0/3007	0.35	0/4062
29	W	0.15	0/787	0.28	0/1060
30	X	0.19	0/2921	0.37	0/3954
31	Y	0.17	0/1073	0.36	0/1444
32	Z	0.17	0/473	0.45	0/631

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
33	a	0.15	0/1801	0.36	0/2447
All	All	0.17	0/72210	0.35	0/101127

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
1	0	0	1
16	J	0	1
All	All	0	2

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	0	195	ARG	Sidechain
16	J	72	LYS	Peptide

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	0	1754	0	1754	11	0
2	1	2195	0	2217	16	0
3	4	4585	0	4595	49	0
4	5	2568	0	2655	20	0
5	7	3145	0	3159	16	0
6	8	2543	0	2561	30	0
7	9	3558	0	3636	29	0
8	A	15655	0	7953	50	0
9	B	1789	0	1781	8	0
10	C	1042	0	1033	16	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
11	D	1838	0	1858	4	0
12	E	839	0	848	12	0
13	F	1724	0	1769	13	0
14	G	2395	0	2364	16	0
15	H	1045	0	1065	14	0
16	J	839	0	887	5	0
17	K	283	0	289	3	0
18	L	1091	0	1137	11	0
19	M	913	0	943	1	0
20	N	859	0	922	1	0
21	O	1570	0	1533	7	0
22	P	731	0	761	6	0
23	Q	49	0	36	0	0
24	R	2382	0	2405	8	0
25	S	1100	0	1103	6	0
26	T	1344	0	1359	9	0
27	U	1468	0	1478	6	0
28	V	2946	0	2942	17	0
29	W	775	0	791	3	0
30	X	2849	0	2843	17	0
31	Y	1043	0	996	10	0
32	Z	465	0	475	6	0
33	a	1785	0	1599	16	0
34	7	8	0	0	0	0
35	7	27	0	22	1	0
36	A	4	0	0	0	0
37	A	26	0	0	0	0
37	B	1	0	0	0	0
37	X	1	0	0	0	0
38	E	4	0	0	0	0
38	T	4	0	0	0	0
39	O	1	0	0	0	0
40	X	31	0	12	0	0
41	X	28	0	12	1	0
All	All	69302	0	61793	364	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 3.

All (364) 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)
2:1:198:TYR:OH	2:1:203:ASP:OD1	1.96	0.84
4:5:17:THR:OG1	4:5:20:GLU:OE1	1.96	0.82
2:1:185:HIS:NE2	2:1:240:GLU:OE2	2.13	0.82
8:A:1081:U:O2'	8:A:1083:C:OP2	1.99	0.79
3:4:243:ASN:OD1	3:4:246:ARG:NH2	2.16	0.78
5:7:300:LYS:NZ	5:7:344:ASN:O	2.16	0.78
1:0:177:GLU:OE2	1:0:181:ASN:ND2	2.17	0.77
13:F:240:ARG:NH2	33:a:307:MET:O	2.17	0.76
26:T:103:ARG:NH2	26:T:110:GLU:OE2	2.18	0.76
5:7:193:GLY:O	35:7:502:SAM:N	2.19	0.76
8:A:1179:G:O6	8:A:1476:G:N2	2.17	0.76
7:9:87:ARG:NH1	7:9:90:GLN:OE1	2.21	0.74
10:C:118:GLU:OE2	10:C:152:ARG:NH2	2.22	0.72
13:F:240:ARG:NH1	33:a:315:ASP:OD2	2.23	0.71
18:L:115:ILE:HG21	18:L:181:ILE:HD13	1.73	0.71
8:A:1248:C:O2'	8:A:1251:A:O2'	2.10	0.69
7:9:256:ASP:OD2	7:9:257:ALA:N	2.24	0.69
6:8:152:PHE:O	6:8:300:ASN:ND2	2.26	0.69
3:4:612:ASN:OD1	3:4:648:ARG:NH1	2.25	0.69
27:U:79:GLN:O	27:U:83:HIS:ND1	2.23	0.68
6:8:95:ARG:NH2	6:8:101:ASP:OD1	2.26	0.68
8:A:1016:U:O2'	22:P:91:ILE:O	2.11	0.67
8:A:1248:C:HO2'	8:A:1251:A:HO2'	1.42	0.67
18:L:86:ASP:OD2	18:L:87:ASP:N	2.28	0.67
30:X:153:LEU:HD11	30:X:244:LEU:CD2	2.25	0.67
28:V:226:TYR:HE1	28:V:282:VAL:HG21	1.59	0.66
20:N:11:ARG:NH1	26:T:81:ASP:OD1	2.29	0.66
3:4:290:ASP:OD1	3:4:293:THR:OG1	2.07	0.65
6:8:290:ILE:HD12	6:8:313:HIS:HD2	1.61	0.65
4:5:36:ASN:ND2	4:5:219:PRO:O	2.29	0.65
5:7:41:SER:O	5:7:46:LYS:NZ	2.25	0.65
14:G:265:GLN:N	14:G:265:GLN:OE1	2.30	0.65
1:0:197:ARG:NH1	1:0:199:GLU:OE1	2.30	0.64
4:5:280:LEU:HB2	4:5:310:MET:HE1	1.78	0.64
18:L:76:TYR:CD2	18:L:93:LEU:HD22	2.32	0.64
9:B:65:GLU:OE2	9:B:69:HIS:NE2	2.28	0.64
33:a:322:TYR:O	33:a:325:THR:OG1	2.10	0.64
6:8:56:ARG:NE	8:A:1227:G:OP2	2.30	0.64
3:4:619:LYS:NZ	3:4:652:ASP:O	2.29	0.64
6:8:194:LEU:HD11	6:8:260:THR:CG2	2.28	0.64
6:8:210:GLU:OE1	6:8:210:GLU:N	2.30	0.64
14:G:116:LEU:O	14:G:122:ARG:NE	2.31	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
24:R:252:ASP:OD2	24:R:256:ARG:NH2	2.31	0.64
11:D:366:LYS:O	11:D:392:ARG:NH1	2.31	0.63
13:F:119:LYS:NZ	30:X:398:LEU:OXT	2.25	0.63
13:F:85:VAL:HG22	30:X:379:GLU:OE1	1.98	0.63
8:A:702:C:OP1	8:A:848:U:O2'	2.15	0.63
8:A:1037:A:O2'	18:L:152:HIS:NE2	2.30	0.63
28:V:274:LYS:NZ	28:V:345:GLY:O	2.32	0.63
8:A:1118:A:OP2	25:S:50:ARG:NH2	2.31	0.63
6:8:278:VAL:O	6:8:282:THR:OG1	2.13	0.63
7:9:569:THR:HG21	7:9:577:LEU:HD22	1.81	0.62
2:1:169:ARG:NH1	2:1:210:ASP:O	2.32	0.62
14:G:129:GLU:OE2	14:G:130:GLN:NE2	2.32	0.62
14:G:344:ILE:HG22	14:G:348:MET:HE2	1.81	0.62
24:R:251:GLU:OE2	24:R:274:TYR:OH	2.07	0.62
2:1:131:THR:OG1	15:H:149:THR:OG1	2.15	0.62
8:A:1226:C:O2'	8:A:1227:G:OP1	2.11	0.62
3:4:302:VAL:HG21	3:4:341:CYS:HB3	1.81	0.61
8:A:932:C:N3	26:T:11:ARG:NH2	2.48	0.61
28:V:226:TYR:CE1	28:V:282:VAL:HG21	2.35	0.61
8:A:927:G:OP2	16:J:47:ARG:NH1	2.32	0.61
33:a:267:ARG:NH1	33:a:269:ALA:O	2.34	0.61
1:0:195:ARG:NH1	1:0:196:ILE:HD11	2.16	0.61
2:1:131:THR:HG1	15:H:149:THR:HG1	1.42	0.61
4:5:64:PRO:CG	4:5:83:VAL:HG13	2.31	0.60
8:A:1035:U:H5''	12:E:72:THR:HG23	1.83	0.60
24:R:302:GLN:O	24:R:306:VAL:HG23	2.02	0.60
3:4:530:GLU:O	3:4:534:LEU:HD13	2.02	0.59
8:A:691:A:N7	8:A:716:U:O2'	2.35	0.59
31:Y:339:GLU:OE2	31:Y:340:SER:OG	2.21	0.59
4:5:22:ILE:HD13	4:5:33:LEU:HD21	1.85	0.58
3:4:656:ASN:OD1	3:4:657:GLN:N	2.36	0.58
7:9:548:LEU:HD11	7:9:642:HIS:HD2	1.68	0.58
3:4:448:ILE:HG21	3:4:454:ARG:HG3	1.85	0.58
9:B:239:ASN:OD1	29:W:119:LYS:NZ	2.35	0.58
14:G:199:TRP:NE1	14:G:226:GLU:OE2	2.36	0.57
3:4:244:ALA:HB1	3:4:262:MET:HE1	1.85	0.57
18:L:68:PRO:O	18:L:71:THR:HG22	2.04	0.57
2:1:292:TYR:OH	30:X:338:ASP:OD1	2.23	0.57
6:8:290:ILE:HD12	6:8:313:HIS:CD2	2.40	0.57
26:T:132:ARG:NH1	26:T:136:LEU:O	2.38	0.56
4:5:21:ILE:HD12	4:5:39:LEU:HD11	1.86	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:C:106:ASP:OD1	10:C:107:GLN:N	2.34	0.56
8:A:1221:A:O2'	8:A:1222:A:OP2	2.21	0.56
10:C:136:VAL:HG22	10:C:153:LEU:HD22	1.87	0.56
28:V:57:MET:CE	28:V:75:LEU:HD13	2.35	0.56
28:V:30:LEU:HD12	28:V:149:ASP:HB2	1.88	0.56
10:C:123:VAL:HG23	10:C:157:THR:HG22	1.87	0.56
13:F:119:LYS:NZ	30:X:365:TRP:O	2.28	0.56
25:S:124:ALA:O	25:S:127:ALA:N	2.39	0.55
33:a:117:VAL:HG22	33:a:166:LEU:HD12	1.87	0.55
3:4:195:LEU:O	3:4:199:GLY:N	2.39	0.55
28:V:57:MET:HE1	28:V:75:LEU:HD13	1.88	0.55
3:4:508:VAL:HG12	3:4:512:ILE:CD1	2.36	0.55
5:7:268:LEU:HB2	5:7:305:LEU:HD21	1.88	0.55
10:C:123:VAL:CG2	10:C:157:THR:HG22	2.36	0.55
12:E:15:ARG:NH1	27:U:182:ASP:OD2	2.39	0.55
7:9:351:LYS:NZ	7:9:497:THR:O	2.39	0.55
14:G:115:GLY:N	15:H:84:ASP:OD2	2.38	0.55
33:a:239:ASP:OD1	33:a:309:SER:OG	2.14	0.55
3:4:62:LYS:O	15:H:66:SER:OG	2.21	0.55
24:R:251:GLU:OE2	24:R:280:LYS:NZ	2.39	0.55
7:9:326:GLU:OE2	7:9:326:GLU:N	2.41	0.54
7:9:572:ASP:OD1	7:9:573:ARG:N	2.40	0.54
3:4:154:GLU:OE2	3:4:187:THR:OG1	2.23	0.54
5:7:109:ARG:NH1	5:7:142:THR:HG22	2.22	0.54
9:B:186:THR:HG22	9:B:186:THR:O	2.08	0.54
4:5:306:VAL:HG12	4:5:310:MET:HE2	1.90	0.53
6:8:194:LEU:HD11	6:8:260:THR:HG22	1.88	0.53
2:1:138:ASP:OD2	2:1:139:SER:N	2.42	0.53
9:B:202:ILE:HG22	9:B:202:ILE:O	2.08	0.53
7:9:184:LEU:HD23	7:9:372:ILE:CG2	2.38	0.53
5:7:178:ARG:HD3	5:7:206:ILE:HG23	1.90	0.53
5:7:423:ASP:OD1	5:7:424:LEU:N	2.41	0.53
7:9:209:LEU:O	7:9:241:LYS:NZ	2.31	0.52
3:4:120:ILE:HD12	32:Z:68:LEU:HD13	1.91	0.52
8:A:1013:A:O2'	8:A:1014:A:O5'	2.24	0.52
3:4:324:VAL:HG22	31:Y:259:PHE:CE2	2.45	0.52
17:K:53:ARG:O	17:K:56:SER:OG	2.18	0.52
1:0:170:LEU:HD21	19:M:96:PHE:CE1	2.44	0.52
3:4:372:TYR:CE2	3:4:400:LEU:HD21	2.45	0.52
8:A:918:A:OP1	21:O:89:ARG:NH1	2.42	0.51
15:H:92:GLU:OE1	15:H:141:ARG:NH1	2.43	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
30:X:153:LEU:HD11	30:X:244:LEU:HD23	1.93	0.51
6:8:282:THR:HG22	6:8:286:LYS:NZ	2.24	0.51
6:8:328:GLN:OE1	6:8:328:GLN:N	2.44	0.51
3:4:527:LEU:HD23	3:4:531:ILE:HD12	1.93	0.51
7:9:265:ARG:NH2	8:A:1149:G:OP2	2.43	0.51
15:H:106:ILE:HD12	15:H:145:LEU:HD21	1.92	0.51
3:4:357:LEU:HD21	31:Y:251:PHE:HE1	1.76	0.51
24:R:326:ASN:O	24:R:330:VAL:HG23	2.10	0.51
7:9:634:THR:HG21	7:9:664:VAL:HG21	1.92	0.51
5:7:415:LEU:HD22	5:7:428:ALA:HB2	1.92	0.51
6:8:194:LEU:HD13	6:8:262:MET:SD	2.50	0.50
3:4:343:ARG:HA	3:4:378:LEU:HD13	1.93	0.50
10:C:138:TYR:O	10:C:142:LEU:HD23	2.11	0.50
14:G:217:ASP:OD1	14:G:218:TYR:N	2.44	0.50
21:O:214:SER:OG	28:V:318:ASP:OD1	2.16	0.50
24:R:191:ARG:HG3	24:R:204:ILE:HG23	1.93	0.50
3:4:92:ASP:OD1	10:C:157:THR:OG1	2.22	0.50
6:8:348:VAL:HG23	6:8:348:VAL:O	2.12	0.50
10:C:104:LEU:HD23	10:C:123:VAL:HG12	1.94	0.50
2:1:184:ASP:OD1	2:1:184:ASP:N	2.45	0.50
4:5:189:GLY:O	4:5:293:ARG:NH1	2.44	0.50
6:8:111:GLN:NE2	6:8:115:ASP:OD1	2.44	0.50
30:X:70:ILE:CD1	30:X:99:LEU:HD13	2.42	0.50
8:A:843:G:N2	8:A:846:A:OP2	2.38	0.50
30:X:74:ASP:OD1	30:X:147:LYS:NZ	2.40	0.49
3:4:508:VAL:HG12	3:4:512:ILE:HD11	1.94	0.49
5:7:259:PHE:O	5:7:289:GLY:N	2.44	0.49
33:a:171:ILE:HD12	33:a:218:GLN:HB2	1.95	0.49
12:E:94:VAL:CG1	22:P:117:MET:HE1	2.43	0.49
3:4:282:LEU:HD12	3:4:287:LEU:HB2	1.95	0.49
13:F:159:VAL:HG23	13:F:172:VAL:HG11	1.95	0.49
28:V:29:LEU:HD13	28:V:152:ILE:HG21	1.95	0.49
3:4:526:ASP:O	3:4:529:GLU:N	2.46	0.49
6:8:194:LEU:HD11	6:8:260:THR:HG23	1.93	0.49
7:9:231:LEU:HD12	7:9:234:LEU:HD12	1.95	0.49
6:8:231:LEU:HD11	6:8:235:CYS:SG	2.53	0.49
21:O:53:ASP:OD2	21:O:54:GLU:N	2.46	0.49
3:4:526:ASP:O	3:4:530:GLU:OE1	2.31	0.48
5:7:82:THR:HG23	8:A:1190:C:H5'	1.95	0.48
6:8:206:PHE:O	6:8:210:GLU:OE1	2.30	0.48
8:A:899:G:O2'	8:A:907:A:N1	2.43	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:E:11:LYS:HA	12:E:89:ILE:HD11	1.95	0.48
3:4:82:THR:HG22	15:H:50:LEU:HD23	1.94	0.48
6:8:178:ARG:HB2	6:8:281:LEU:HD13	1.96	0.48
8:A:1033:U:O2'	12:E:93:ILE:O	2.31	0.48
18:L:112:MET:HE1	18:L:127:ALA:CB	2.43	0.48
33:a:112:ARG:NH2	33:a:327:ALA:O	2.44	0.48
3:4:357:LEU:HD21	31:Y:251:PHE:CE1	2.48	0.48
15:H:84:ASP:OD1	15:H:86:ALA:N	2.46	0.48
1:0:41:LEU:HD13	1:0:55:TRP:CG	2.49	0.48
1:0:166:TYR:O	21:O:199:TRP:NE1	2.44	0.48
2:1:275:ASN:OD1	2:1:276:MET:N	2.47	0.48
7:9:634:THR:CG2	7:9:664:VAL:HG21	2.43	0.48
14:G:175:HIS:O	14:G:175:HIS:ND1	2.46	0.48
2:1:87:MET:HE3	14:G:123:PRO:HD3	1.95	0.47
10:C:138:TYR:CE1	32:Z:70:LEU:HD13	2.50	0.47
3:4:120:ILE:CD1	32:Z:68:LEU:HD13	2.44	0.47
9:B:67:LEU:HD11	9:B:138:ARG:NH2	2.28	0.47
25:S:128:GLU:OE1	25:S:128:GLU:N	2.47	0.47
26:T:92:THR:O	26:T:92:THR:HG22	2.14	0.47
8:A:1188:A:H62	8:A:1466:C:H42	1.62	0.47
3:4:429:LEU:HA	3:4:464:LEU:HD21	1.95	0.47
8:A:738:A:O2'	8:A:739:C:O2	2.26	0.47
6:8:335:GLU:O	8:A:1227:G:N2	2.39	0.47
3:4:117:ALA:HB1	10:C:142:LEU:HD22	1.96	0.47
6:8:307:ARG:NH2	8:A:1230:C:N3	2.63	0.47
5:7:109:ARG:CZ	5:7:142:THR:HG22	2.45	0.47
8:A:1430:A:OP1	13:F:35:SER:N	2.48	0.47
8:A:1559:G:OP2	8:A:1559:G:N2	2.41	0.47
12:E:94:VAL:HG11	22:P:117:MET:HE1	1.95	0.47
29:W:114:ILE:HG21	29:W:142:LEU:HD11	1.96	0.47
32:Z:22:VAL:HG21	32:Z:26:THR:HG21	1.97	0.47
2:1:217:GLN:NE2	31:Y:326:SER:O	2.48	0.47
28:V:190:LEU:HD22	28:V:226:TYR:OH	2.14	0.47
3:4:310:GLU:N	3:4:310:GLU:OE2	2.48	0.46
10:C:69:LEU:HD11	10:C:107:GLN:OE1	2.15	0.46
14:G:215:ASP:OD1	14:G:216:LEU:N	2.48	0.46
3:4:111:LYS:NZ	3:4:115:ASN:OD1	2.48	0.46
6:8:77:MET:HE2	6:8:350:PRO:HB3	1.97	0.46
11:D:383:VAL:HG13	11:D:383:VAL:O	2.16	0.46
6:8:77:MET:HE2	6:8:350:PRO:CB	2.46	0.46
7:9:96:GLN:O	7:9:99:GLU:N	2.48	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:4:58:VAL:HG23	3:4:58:VAL:O	2.15	0.46
3:4:336:ASN:ND2	3:4:409:ASP:OD2	2.42	0.46
6:8:82:LEU:HD21	6:8:103:LEU:HD11	1.98	0.46
7:9:70:GLU:O	28:V:392:ARG:NH1	2.45	0.46
2:1:107:LYS:NZ	10:C:54:GLU:OE2	2.49	0.46
3:4:538:ASP:OD1	3:4:539:LYS:N	2.47	0.46
13:F:46:ILE:O	13:F:46:ILE:HG22	2.15	0.46
3:4:645:LEU:O	3:4:649:VAL:HG23	2.15	0.46
33:a:321:GLN:O	33:a:325:THR:HG23	2.16	0.46
3:4:194:LEU:HD12	3:4:198:TYR:HD2	1.81	0.46
4:5:64:PRO:HG3	4:5:83:VAL:HG13	1.98	0.46
6:8:307:ARG:NH1	8:A:1231:A:N1	2.64	0.46
8:A:953:U:O4	26:T:36:THR:HG21	2.16	0.46
26:T:30:MET:HE2	26:T:78:PHE:CE2	2.51	0.46
32:Z:63:GLN:NE2	32:Z:64:THR:HG23	2.30	0.46
2:1:263:LEU:O	2:1:267:LEU:HD23	2.16	0.46
4:5:64:PRO:HG2	4:5:83:VAL:HG13	1.97	0.46
26:T:91:GLU:OE1	27:U:120:ARG:NH1	2.50	0.45
3:4:392:ILE:HD12	3:4:392:ILE:H	1.81	0.45
5:7:29:ALA:C	5:7:30:LEU:HD12	2.41	0.45
18:L:112:MET:HE1	18:L:127:ALA:HB2	1.98	0.45
2:1:152:ASP:OD1	15:H:175:THR:OG1	2.28	0.45
8:A:703:A:OP2	27:U:43:ASN:ND2	2.49	0.45
22:P:122:VAL:HG13	22:P:123:THR:HG23	1.98	0.45
1:0:9:ARG:NE	8:A:805:C:O2	2.49	0.45
8:A:971:A:O2'	12:E:76:GLU:OE2	2.31	0.45
8:A:1231:A:H2'	8:A:1232:A:H4'	1.97	0.45
8:A:1343:A:O2'	8:A:1345:G:N7	2.32	0.45
7:9:617:SER:OG	7:9:665:ASN:OD1	2.32	0.45
27:U:126:GLN:O	27:U:130:GLU:OE1	2.35	0.45
2:1:86:ARG:NH1	2:1:96:PRO:O	2.47	0.45
7:9:220:MET:CE	7:9:354:LEU:HD22	2.47	0.45
7:9:642:HIS:C	7:9:643:LEU:HD12	2.42	0.45
30:X:159:HIS:NE2	30:X:266:ASN:OD1	2.49	0.45
21:O:185:SER:O	24:R:183:LYS:NZ	2.49	0.44
1:0:37:ASP:O	1:0:41:LEU:N	2.50	0.44
3:4:104:ARG:O	3:4:108:LEU:HD13	2.17	0.44
8:A:897:C:OP1	16:J:114:ARG:NH2	2.51	0.44
8:A:1119:U:C4'	25:S:50:ARG:HH12	2.31	0.44
7:9:184:LEU:HD23	7:9:372:ILE:HG22	2.00	0.44
3:4:472:ASP:OD1	3:4:473:VAL:N	2.49	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:V:272:ASP:OD1	28:V:273:ILE:N	2.51	0.44
30:X:253:LEU:HD12	30:X:255:MET:HE3	2.00	0.44
6:8:252:SER:O	6:8:256:HIS:N	2.51	0.44
7:9:642:HIS:O	7:9:643:LEU:HD12	2.18	0.44
8:A:1245:U:O2'	8:A:1342:C:N4	2.51	0.44
10:C:75:ASN:OD1	17:K:104:TRP:NE1	2.48	0.44
24:R:347:GLN:NE2	24:R:351:GLU:OE2	2.50	0.44
30:X:192:LYS:NZ	30:X:227:GLU:OE1	2.42	0.44
31:Y:322:ASP:OD2	31:Y:325:GLY:N	2.51	0.44
4:5:288:PRO:HB2	7:9:400:MET:HE1	2.00	0.43
3:4:454:ARG:NH1	31:Y:292:GLN:OE1	2.51	0.43
18:L:112:MET:O	18:L:116:VAL:HG22	2.18	0.43
30:X:186:THR:O	30:X:190:ASN:ND2	2.51	0.43
30:X:108:LEU:HD21	30:X:307:VAL:CG1	2.48	0.43
8:A:1035:U:O3'	12:E:72:THR:HG21	2.18	0.43
13:F:84:SER:OG	30:X:379:GLU:OE2	2.33	0.43
14:G:200:LEU:HD23	14:G:205:LEU:HD13	1.99	0.43
3:4:327:LYS:HG3	3:4:327:LYS:O	2.19	0.43
13:F:46:ILE:HG23	14:G:315:PHE:CZ	2.53	0.43
1:0:29:ARG:CD	1:0:105:THR:HG21	2.49	0.43
4:5:83:VAL:HG21	4:5:105:LEU:HD11	2.00	0.43
6:8:282:THR:HG22	6:8:286:LYS:HZ3	1.83	0.43
8:A:1018:G:OP2	22:P:94:ARG:NE	2.37	0.43
30:X:167:ASP:OD1	30:X:167:ASP:O	2.37	0.43
3:4:259:TYR:CB	3:4:282:LEU:HD13	2.49	0.43
8:A:1441:A:H2	8:A:1449:G:H22	1.67	0.43
14:G:320:VAL:HG23	14:G:322:ARG:HG2	2.01	0.43
31:Y:327:GLU:N	31:Y:330:GLU:OE1	2.50	0.43
5:7:344:ASN:OD1	5:7:345:LEU:N	2.52	0.42
6:8:73:GLN:HA	6:8:77:MET:HE1	2.00	0.42
6:8:246:ARG:HG3	6:8:250:LYS:NZ	2.34	0.42
7:9:666:ILE:HG23	7:9:666:ILE:O	2.18	0.42
7:9:344:VAL:HG12	7:9:345:GLY:N	2.34	0.42
7:9:619:ALA:HA	7:9:634:THR:HG22	2.01	0.42
11:D:401:VAL:HG21	26:T:50:PHE:CD1	2.55	0.42
4:5:39:LEU:HD12	4:5:218:VAL:HG21	2.02	0.42
12:E:47:LEU:HD23	12:E:51:ILE:HD12	2.01	0.42
14:G:210:VAL:HG12	14:G:210:VAL:O	2.19	0.42
3:4:513:TRP:HD1	3:4:531:ILE:HD13	1.84	0.42
13:F:236:LEU:HD23	33:a:248:LEU:HD22	2.01	0.42
22:P:134:LYS:NZ	22:P:136:CYS:O	2.51	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:0:195:ARG:HH11	1:0:196:ILE:HD11	1.84	0.42
15:H:92:GLU:OE2	15:H:108:VAL:HG11	2.19	0.42
25:S:124:ALA:O	25:S:128:GLU:OE1	2.38	0.42
1:0:175:ILE:HG21	21:O:188:PRO:HD2	2.01	0.42
8:A:875:U:O2'	8:A:882:A:N1	2.49	0.42
11:D:245:VAL:HG22	11:D:271:ALA:HB1	2.01	0.42
16:J:78:ARG:NE	16:J:117:ASP:OD2	2.33	0.42
10:C:123:VAL:HG11	10:C:155:LEU:HD22	2.02	0.42
3:4:397:MET:HE3	3:4:397:MET:HA	2.01	0.42
4:5:148:THR:HB	4:5:149:PRO:HD3	2.02	0.42
6:8:166:MET:HE3	6:8:288:VAL:HG11	2.01	0.42
33:a:112:ARG:NE	33:a:327:ALA:O	2.51	0.42
8:A:700:A:N1	8:A:709:G:O2'	2.53	0.42
12:E:29:LEU:HD23	12:E:78:MET:SD	2.59	0.42
14:G:269:PHE:CD2	14:G:284:ILE:HG23	2.55	0.42
3:4:397:MET:HE1	3:4:438:LEU:HD22	2.02	0.42
3:4:512:ILE:H	3:4:512:ILE:HD12	1.85	0.42
6:8:231:LEU:HD12	6:8:232:PHE:CD1	2.55	0.42
30:X:353:PHE:CZ	30:X:357:ILE:HD11	2.54	0.42
14:G:263:ASP:OD2	14:G:267:MET:N	2.52	0.41
28:V:398:GLU:O	28:V:402:GLN:OE1	2.38	0.41
3:4:302:VAL:O	3:4:344:ARG:NH2	2.50	0.41
9:B:144:PHE:CE1	9:B:158:MET:HE1	2.54	0.41
15:H:108:VAL:HG22	15:H:143:LEU:CD2	2.49	0.41
17:K:51:ARG:HA	17:K:54:ILE:HG12	2.03	0.41
30:X:295:LYS:NZ	41:X:503:GDP:O3A	2.53	0.41
3:4:121:ILE:HD13	10:C:145:TYR:CG	2.55	0.41
28:V:80:SER:N	28:V:84:GLU:OE1	2.42	0.41
2:1:177:LEU:HD22	2:1:190:LEU:HD23	2.00	0.41
5:7:400:HIS:CD2	5:7:414:VAL:HG22	2.55	0.41
25:S:124:ALA:O	25:S:127:ALA:HB3	2.20	0.41
32:Z:61:LEU:O	32:Z:65:LEU:HD13	2.21	0.41
4:5:209:ILE:HD11	4:5:229:HIS:HB2	2.03	0.41
28:V:30:LEU:HD12	28:V:149:ASP:CB	2.49	0.41
28:V:79:ILE:HG21	28:V:85:ILE:HD13	2.02	0.41
8:A:1279:C:O2'	8:A:1296:A:N1	2.50	0.41
15:H:50:LEU:HD13	15:H:53:ASP:O	2.19	0.41
15:H:67:ASP:OD1	15:H:67:ASP:N	2.48	0.41
7:9:266:GLU:O	7:9:270:GLU:OE1	2.38	0.41
7:9:622:ASP:OD1	7:9:632:SER:OG	2.37	0.41
8:A:1258:A:H61	8:A:1330:C:H5'	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:B:229:PRO:HA	9:B:232:ILE:HD13	2.03	0.41
4:5:15:LEU:HD12	4:5:16:PRO:HD2	2.02	0.41
13:F:236:LEU:CD2	33:a:248:LEU:HD22	2.51	0.41
5:7:170:VAL:HG21	5:7:199:VAL:HG22	2.03	0.41
5:7:268:LEU:CB	5:7:305:LEU:HD21	2.51	0.41
15:H:108:VAL:HG13	15:H:143:LEU:HD23	2.02	0.41
18:L:122:THR:O	18:L:122:THR:HG22	2.20	0.41
21:O:62:GLU:O	21:O:66:GLU:OE1	2.37	0.41
29:W:142:LEU:HD23	29:W:169:LEU:HA	2.02	0.41
3:4:286:ARG:O	3:4:286:ARG:HG2	2.21	0.41
8:A:930:G:H22	16:J:44:ARG:HB3	1.85	0.41
33:a:117:VAL:HG11	33:a:194:MET:CE	2.51	0.41
7:9:234:LEU:HD13	7:9:243:LEU:HD11	2.01	0.40
8:A:1187:U:O2'	8:A:1465:C:N4	2.54	0.40
18:L:158:MET:O	18:L:162:GLN:OE1	2.38	0.40
28:V:156:ASN:ND2	28:V:159:ASP:OD2	2.52	0.40
28:V:202:ARG:HG2	28:V:247:MET:HE2	2.03	0.40
31:Y:282:PHE:CE2	31:Y:286:LEU:HD11	2.56	0.40
12:E:26:ILE:HG21	27:U:173:LEU:CD1	2.51	0.40
33:a:129:LEU:HD13	33:a:203:LEU:HD21	2.02	0.40
4:5:21:ILE:CD1	4:5:39:LEU:HD11	2.50	0.40
8:A:1165:C:OP2	16:J:122:LYS:NZ	2.47	0.40
10:C:92:LEU:HD21	10:C:147:TYR:HE2	1.86	0.40
12:E:10:LEU:CD2	12:E:88:VAL:HG22	2.51	0.40
31:Y:338:LEU:HD11	31:Y:355:THR:HG21	2.03	0.40
33:a:295:GLN:NE2	33:a:296:ARG:O	2.54	0.40
4:5:215:GLN:NE2	8:A:1490:U:OP1	2.53	0.40
7:9:548:LEU:C	7:9:548:LEU:HD12	2.46	0.40
8:A:1151:C:H6	8:A:1154:A:H61	1.68	0.40
13:F:236:LEU:CD1	33:a:264:LEU:HD21	2.51	0.40
4:5:22:ILE:CD1	4:5:33:LEU:HD21	2.49	0.40
4:5:257:ARG:O	4:5:291:ARG:NH1	2.54	0.40
7:9:517:ILE:HG21	7:9:589:ILE:HD11	2.02	0.40
9:B:269:ALA:O	9:B:273:LEU:HD23	2.21	0.40
18:L:116:VAL:HG12	18:L:126:GLU:HG3	2.03	0.40

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	
1	0	209/218 (96%)	206 (99%)	3 (1%)	0	100	100
2	1	269/323 (83%)	263 (98%)	6 (2%)	0	100	100
3	4	558/689 (81%)	550 (99%)	8 (1%)	0	100	100
4	5	317/346 (92%)	313 (99%)	4 (1%)	0	100	100
5	7	394/456 (86%)	383 (97%)	11 (3%)	0	100	100
6	8	324/390 (83%)	315 (97%)	9 (3%)	0	100	100
7	9	438/698 (63%)	425 (97%)	13 (3%)	0	100	100
9	B	218/296 (74%)	216 (99%)	2 (1%)	0	100	100
10	C	122/167 (73%)	119 (98%)	3 (2%)	0	100	100
11	D	228/430 (53%)	224 (98%)	4 (2%)	0	100	100
12	E	103/125 (82%)	100 (97%)	3 (3%)	0	100	100
13	F	206/242 (85%)	204 (99%)	2 (1%)	0	100	100
14	G	288/396 (73%)	282 (98%)	6 (2%)	0	100	100
15	H	124/201 (62%)	121 (98%)	3 (2%)	0	100	100
16	J	106/138 (77%)	101 (95%)	5 (5%)	0	100	100
17	K	27/128 (21%)	27 (100%)	0	0	100	100
18	L	129/257 (50%)	129 (100%)	0	0	100	100
19	M	113/137 (82%)	113 (100%)	0	0	100	100
20	N	107/130 (82%)	107 (100%)	0	0	100	100
21	O	188/258 (73%)	187 (100%)	1 (0%)	0	100	100
22	P	90/142 (63%)	88 (98%)	2 (2%)	0	100	100
23	Q	4/87 (5%)	4 (100%)	0	0	100	100
24	R	289/360 (80%)	283 (98%)	6 (2%)	0	100	100
25	S	131/190 (69%)	130 (99%)	1 (1%)	0	100	100
26	T	162/173 (94%)	160 (99%)	2 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
27	U	172/205 (84%)	170 (99%)	2 (1%)	0	100	100
28	V	355/414 (86%)	351 (99%)	4 (1%)	0	100	100
29	W	96/187 (51%)	95 (99%)	1 (1%)	0	100	100
30	X	350/398 (88%)	346 (99%)	4 (1%)	0	100	100
31	Y	119/395 (30%)	118 (99%)	1 (1%)	0	100	100
32	Z	52/106 (49%)	52 (100%)	0	0	100	100
33	a	253/437 (58%)	247 (98%)	6 (2%)	0	100	100
All	All	6541/9119 (72%)	6429 (98%)	112 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	0	184/190 (97%)	184 (100%)	0	100	100
2	1	250/291 (86%)	250 (100%)	0	100	100
3	4	506/609 (83%)	506 (100%)	0	100	100
4	5	286/309 (93%)	286 (100%)	0	100	100
5	7	339/385 (88%)	339 (100%)	0	100	100
6	8	270/317 (85%)	270 (100%)	0	100	100
7	9	388/600 (65%)	388 (100%)	0	100	100
9	B	194/249 (78%)	194 (100%)	0	100	100
10	C	111/143 (78%)	111 (100%)	0	100	100
11	D	196/357 (55%)	196 (100%)	0	100	100
12	E	89/107 (83%)	89 (100%)	0	100	100
13	F	185/209 (88%)	185 (100%)	0	100	100
14	G	254/342 (74%)	254 (100%)	0	100	100
15	H	118/180 (66%)	118 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
16	J	93/118 (79%)	93 (100%)	0	100	100
17	K	28/113 (25%)	28 (100%)	0	100	100
18	L	123/226 (54%)	123 (100%)	0	100	100
19	M	94/113 (83%)	94 (100%)	0	100	100
20	N	95/115 (83%)	95 (100%)	0	100	100
21	O	171/230 (74%)	171 (100%)	0	100	100
22	P	83/123 (68%)	83 (100%)	0	100	100
23	Q	5/79 (6%)	5 (100%)	0	100	100
24	R	261/318 (82%)	261 (100%)	0	100	100
25	S	115/164 (70%)	115 (100%)	0	100	100
26	T	151/157 (96%)	151 (100%)	0	100	100
27	U	150/174 (86%)	150 (100%)	0	100	100
28	V	323/364 (89%)	323 (100%)	0	100	100
29	W	85/158 (54%)	85 (100%)	0	100	100
30	X	311/351 (89%)	311 (100%)	0	100	100
31	Y	113/357 (32%)	113 (100%)	0	100	100
32	Z	51/95 (54%)	51 (100%)	0	100	100
33	a	151/386 (39%)	151 (100%)	0	100	100
All	All	5773/7929 (73%)	5773 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (63) such sidechains are listed below:

Mol	Chain	Res	Type
1	0	111	HIS
2	1	231	HIS
2	1	268	GLN
3	4	269	HIS
4	5	53	ASN
4	5	109	HIS
4	5	137	HIS
4	5	229	HIS
4	5	254	GLN
5	7	36	GLN

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Mol	Chain	Res	Type
5	7	69	ASN
5	7	86	GLN
5	7	144	HIS
5	7	398	HIS
5	7	400	HIS
5	7	412	HIS
6	8	108	HIS
6	8	136	HIS
6	8	311	HIS
6	8	313	HIS
7	9	106	GLN
7	9	187	HIS
7	9	242	GLN
7	9	551	ASN
7	9	584	HIS
7	9	642	HIS
7	9	669	GLN
9	B	167	HIS
10	C	60	HIS
10	C	156	GLN
11	D	227	ASN
11	D	361	GLN
13	F	151	ASN
14	G	156	GLN
15	H	147	HIS
15	H	183	HIS
16	J	37	HIS
16	J	77	ASN
19	M	75	HIS
20	N	9	HIS
21	O	65	GLN
21	O	169	GLN
22	P	137	ASN
24	R	76	GLN
24	R	214	ASN
24	R	278	ASN
25	S	66	HIS
25	S	82	GLN
26	T	63	GLN
27	U	31	HIS
27	U	61	GLN
27	U	62	HIS

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Mol	Chain	Res	Type
27	U	110	GLN
28	V	87	HIS
28	V	188	HIS
30	X	69	ASN
30	X	110	HIS
30	X	291	HIS
30	X	367	GLN
31	Y	337	HIS
33	a	123	ASN
33	a	243	GLN
33	a	271	HIS

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
8	A	726/955 (76%)	184 (25%)	3 (0%)

All (184) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
8	A	650	U
8	A	651	A
8	A	680	U
8	A	688	A
8	A	689	U
8	A	704	U
8	A	721	U
8	A	722	C
8	A	730	A
8	A	736	C
8	A	737	C
8	A	739	C
8	A	741	A
8	A	753	A
8	A	759	C
8	A	761	A
8	A	766	G
8	A	769	G
8	A	772	A
8	A	773	U
8	A	777	G

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Mol	Chain	Res	Type
8	A	782	A
8	A	791	G
8	A	794	U
8	A	796	G
8	A	807	A
8	A	830	U
8	A	831	U
8	A	832	U
8	A	835	C
8	A	851	A
8	A	860	A
8	A	861	U
8	A	866	A
8	A	867	C
8	A	868	C
8	A	870	C
8	A	871	A
8	A	883	U
8	A	886	C
8	A	889	G
8	A	893	G
8	A	902	G
8	A	903	U
8	A	919	A
8	A	930	G
8	A	931	C
8	A	938	A
8	A	939	A
8	A	942	A
8	A	949	U
8	A	954	C
8	A	955	A
8	A	967	A
8	A	971	A
8	A	972	G
8	A	974	U
8	A	976	A
8	A	1014	A
8	A	1015	A
8	A	1038	C
8	A	1042	U
8	A	1047	A

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Mol	Chain	Res	Type
8	A	1052	C
8	A	1056	A
8	A	1058	C
8	A	1060	A
8	A	1061	A
8	A	1077	U
8	A	1078	A
8	A	1079	G
8	A	1080	A
8	A	1081	U
8	A	1083	C
8	A	1095	U
8	A	1097	G
8	A	1117	A
8	A	1118	A
8	A	1119	U
8	A	1121	A
8	A	1126	A
8	A	1127	A
8	A	1128	C
8	A	1129	U
8	A	1137	A
8	A	1138	G
8	A	1140	A
8	A	1141	C
8	A	1142	A
8	A	1153	C
8	A	1154	A
8	A	1160	A
8	A	1162	A
8	A	1167	A
8	A	1175	G
8	A	1177	C
8	A	1178	G
8	A	1179	G
8	A	1181	G
8	A	1182	C
8	A	1185	C
8	A	1186	A
8	A	1187	U
8	A	1188	A
8	A	1189	U

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Mol	Chain	Res	Type
8	A	1190	C
8	A	1191	C
8	A	1198	A
8	A	1199	G
8	A	1200	G
8	A	1215	U
8	A	1216	C
8	A	1217	G
8	A	1218	A
8	A	1219	U
8	A	1220	A
8	A	1221	A
8	A	1222	A
8	A	1224	C
8	A	1226	C
8	A	1227	G
8	A	1228	A
8	A	1232	A
8	A	1233	C
8	A	1234	C
8	A	1242	C
8	A	1245	U
8	A	1248	C
8	A	1249	U
8	A	1250	C
8	A	1254	C
8	A	1255	U
8	A	1258	A
8	A	1259	U
8	A	1261	C
8	A	1284	U
8	A	1290	C
8	A	1291	U
8	A	1295	A
8	A	1296	A
8	A	1297	G
8	A	1299	A
8	A	1330	C
8	A	1331	A
8	A	1332	A
8	A	1334	G
8	A	1335	U

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Mol	Chain	Res	Type
8	A	1338	A
8	A	1342	C
8	A	1343	A
8	A	1347	G
8	A	1350	G
8	A	1353	A
8	A	1354	A
8	A	1357	A
8	A	1360	G
8	A	1364	U
8	A	1367	A
8	A	1373	U
8	A	1376	C
8	A	1378	C
8	A	1424	U
8	A	1427	A
8	A	1430	A
8	A	1434	A
8	A	1440	G
8	A	1441	A
8	A	1442	G
8	A	1450	C
8	A	1463	G
8	A	1464	G
8	A	1467	C
8	A	1468	U
8	A	1470	A
8	A	1471	A
8	A	1474	G
8	A	1477	U
8	A	1491	C
8	A	1492	A
8	A	1560	U
8	A	1576	G
8	A	1583	A
8	A	1584	A
8	A	1585	A

All (3) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
8	A	1077	U

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Mol	Chain	Res	Type
8	A	1226	C
8	A	1330	C

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 39 ligands modelled in this entry, 33 are monoatomic - leaving 6 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	$\# Z > 2$
38	FES	T	201	26,19	0,4,4	-	-	-		
34	SF4	7	501	5	0,12,12	-	-	-		
38	FES	E	201	12,22	0,4,4	-	-	-		
35	SAM	7	502	-	27,29,29	1.10	5 (18%)	34,42,42	2.00	8 (23%)
40	ATP	X	501	37	32,33,33	0.53	0	48,52,52	0.68	0
41	GDP	X	503	-	29,30,30	3.28	10 (34%)	45,47,47	2.32	13 (28%)

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
38	FES	T	201	26,19	-	-	0/1/1/1
34	SF4	7	501	5	-	-	0/6/5/5

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
38	FES	E	201	12,22	-	-	0/1/1/1
35	SAM	7	502	-	-	6/17/33/33	0/3/3/3
40	ATP	X	501	37	-	2/22/38/38	0/3/3/3
41	GDP	X	503	-	-	4/16/32/32	0/3/3/3

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
41	X	503	GDP	O6-C6	9.78	1.42	1.23
41	X	503	GDP	C2-N1	7.58	1.56	1.37
41	X	503	GDP	PA-O3A	7.38	1.67	1.59
41	X	503	GDP	C2-N3	5.23	1.45	1.33
41	X	503	GDP	C2-N2	5.17	1.46	1.34
41	X	503	GDP	C2'-C3'	-2.91	1.45	1.53
41	X	503	GDP	C6-N1	2.48	1.43	1.38
35	7	502	SAM	C2-N1	2.47	1.38	1.33
35	7	502	SAM	C2-N3	2.46	1.38	1.33
41	X	503	GDP	PB-O3B	-2.39	1.45	1.54
41	X	503	GDP	PA-O5'	2.35	1.68	1.59
35	7	502	SAM	OXT-C	-2.29	1.23	1.30
41	X	503	GDP	PB-O2B	-2.17	1.46	1.54
35	7	502	SAM	C8-N7	2.09	1.35	1.31
35	7	502	SAM	C5-N7	-2.02	1.35	1.39

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
41	X	503	GDP	C8-N7-C5	6.82	116.41	104.26
41	X	503	GDP	N9-C8-N7	-6.81	100.78	113.40
35	7	502	SAM	N3-C2-N1	-5.55	120.18	128.58
41	X	503	GDP	C2-N1-C6	-4.80	116.40	125.11
35	7	502	SAM	C5-C4-N3	-4.59	120.40	126.72
41	X	503	GDP	C8-N9-C4	4.04	113.59	106.03
35	7	502	SAM	N9-C8-N7	-3.82	108.52	113.94
41	X	503	GDP	C6-C5-C4	3.58	124.21	118.83
35	7	502	SAM	C2-N3-C4	3.50	120.37	111.83
35	7	502	SAM	C5-N7-C8	3.45	108.87	103.45
41	X	503	GDP	C4-C5-N7	-3.22	105.56	110.67
41	X	503	GDP	O3B-PB-O3A	2.98	114.62	104.64
35	7	502	SAM	N3-C4-N9	2.80	131.92	127.17
41	X	503	GDP	C2-N3-C4	2.79	117.10	112.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
41	X	503	GDP	C3'-C2'-C1'	2.71	106.60	101.46
35	7	502	SAM	OXT-C-O	-2.68	117.99	124.08
41	X	503	GDP	N9-C4-N3	2.67	131.29	125.95
41	X	503	GDP	O2B-PB-O3A	2.64	113.47	104.64
35	7	502	SAM	C4-C5-N7	-2.57	107.64	110.58
41	X	503	GDP	O2A-PA-O1A	-2.31	101.72	112.44
41	X	503	GDP	C5-C4-N3	-2.10	125.04	128.39

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
41	X	503	GDP	C5'-O5'-PA-O3A
41	X	503	GDP	C5'-O5'-PA-O1A
40	X	501	ATP	PA-O3A-PB-O1B
35	7	502	SAM	O4'-C4'-C5'-SD
35	7	502	SAM	CB-CG-SD-C5'
35	7	502	SAM	C3'-C4'-C5'-SD
35	7	502	SAM	OXT-C-CA-CB
35	7	502	SAM	O-C-CA-CB
41	X	503	GDP	PB-O3A-PA-O1A
35	7	502	SAM	CB-CG-SD-CE
41	X	503	GDP	PA-O3A-PB-O1B
40	X	501	ATP	PA-O3A-PB-O2B

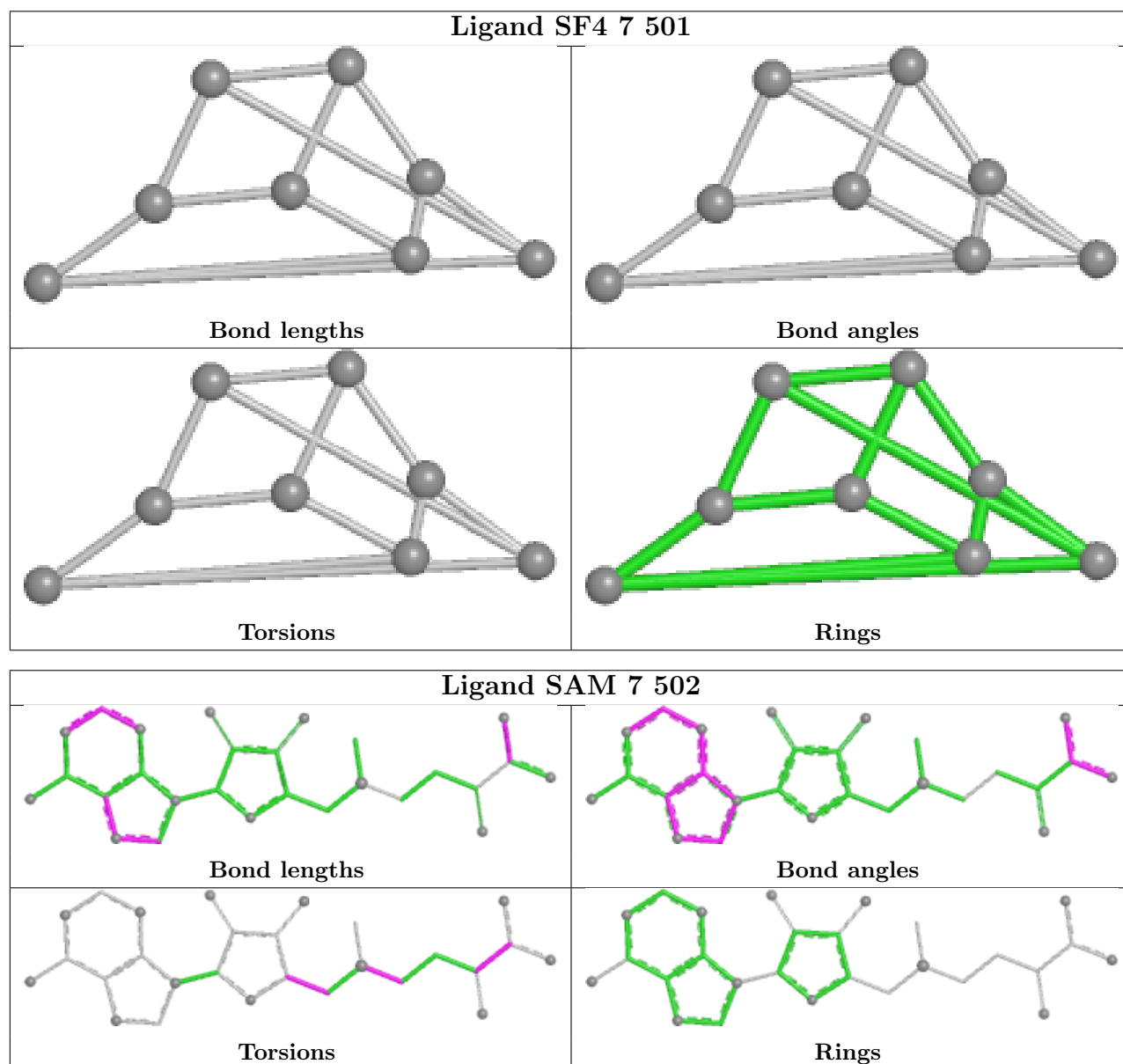
There are no ring outliers.

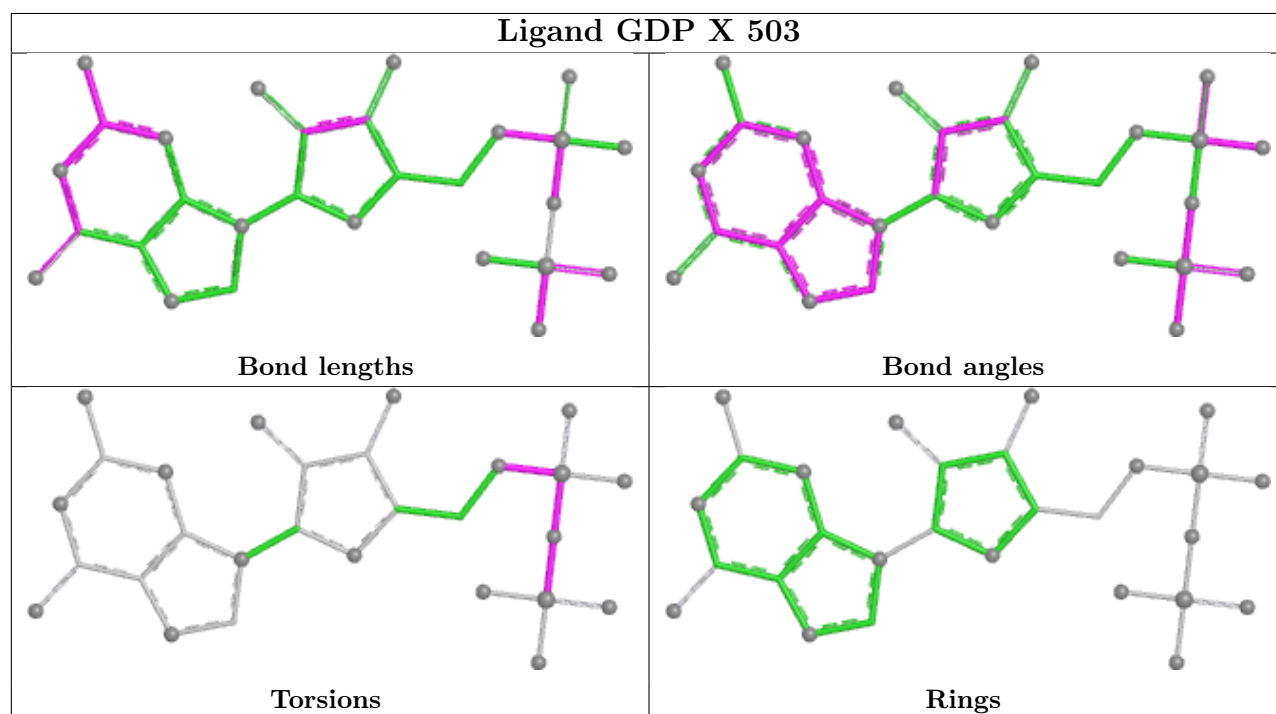
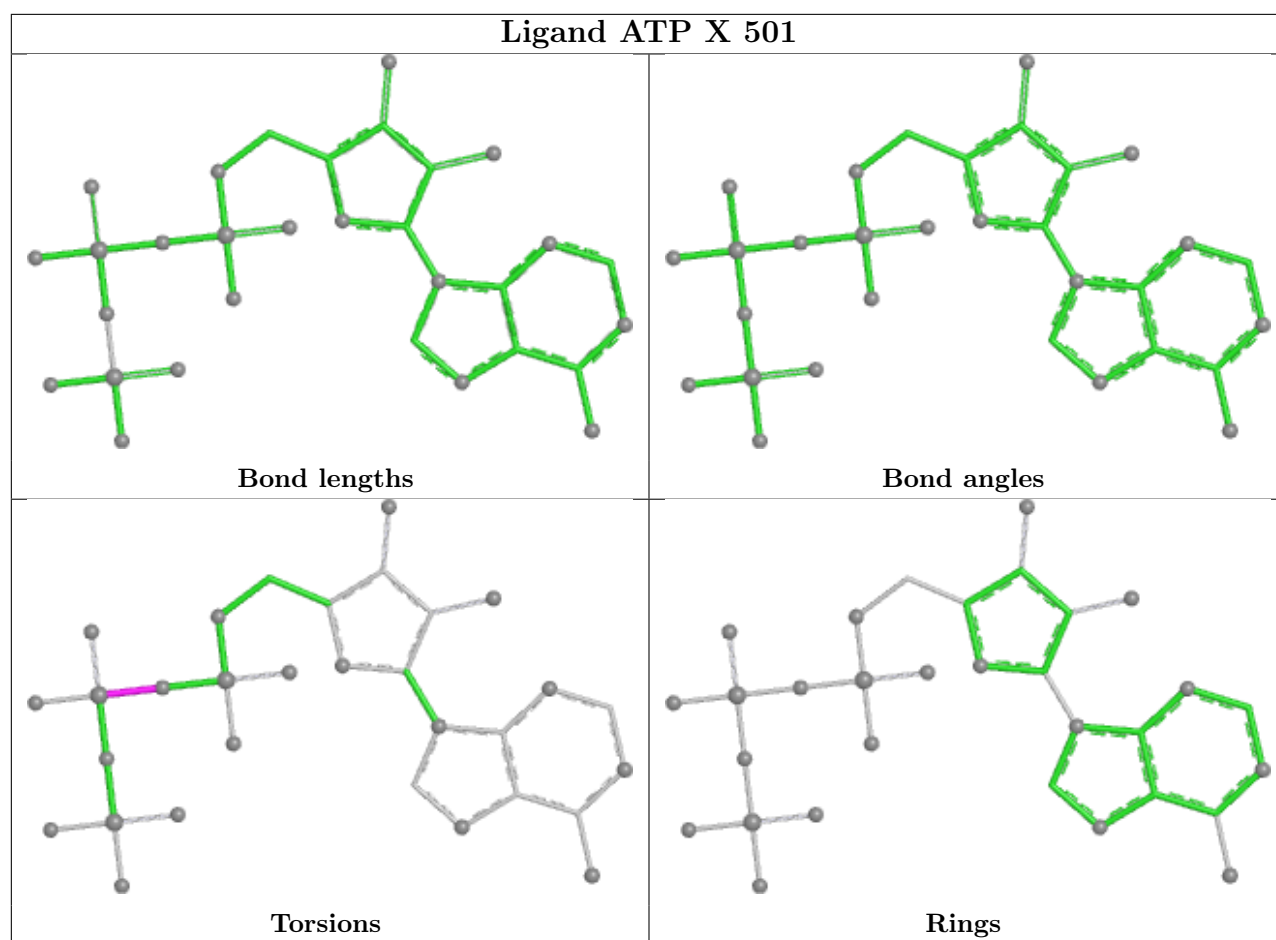
2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
35	7	502	SAM	1	0
41	X	503	GDP	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](#)

There are no chain breaks in this entry.

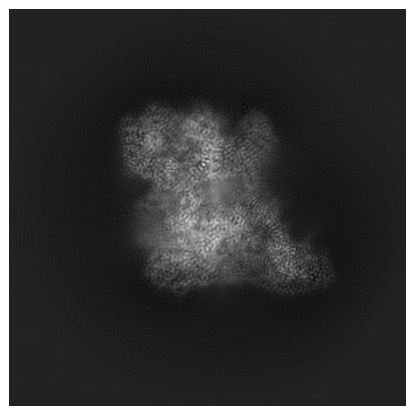
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-26966. These allow visual inspection of the internal detail of the map and identification of artifacts.

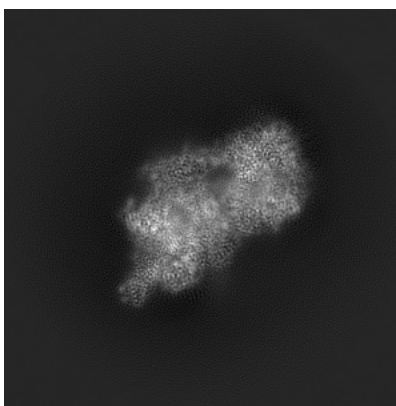
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

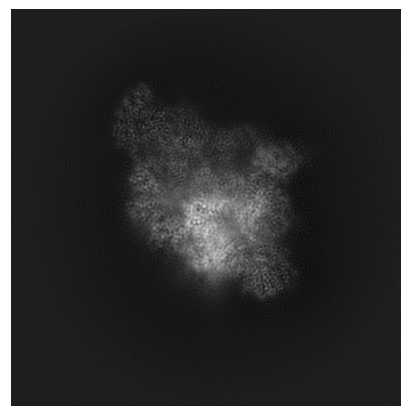
6.1.1 Primary map



X

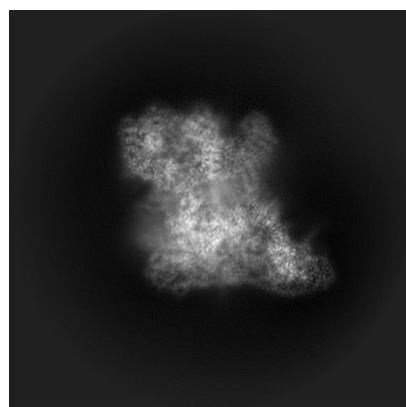


Y

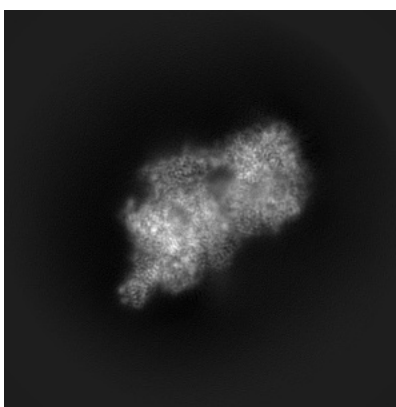


Z

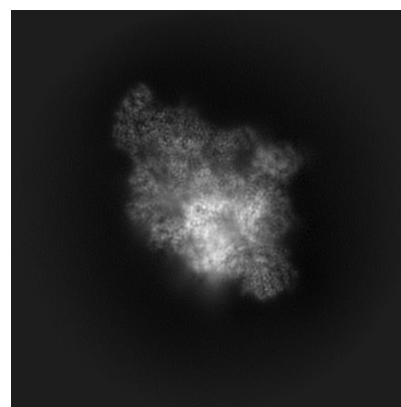
6.1.2 Raw map



X



Y

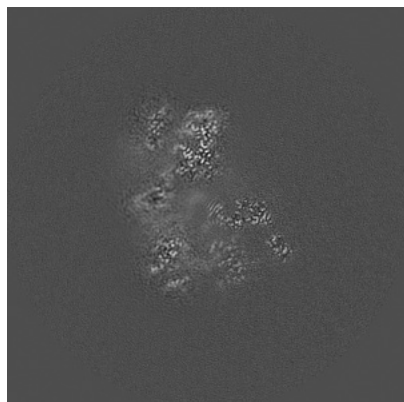


Z

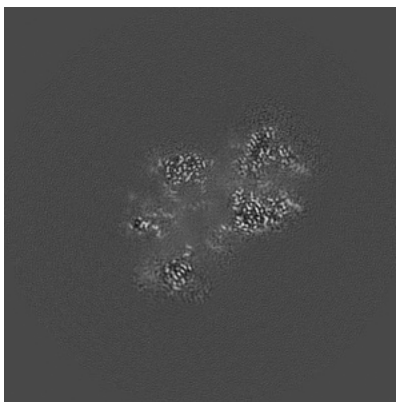
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

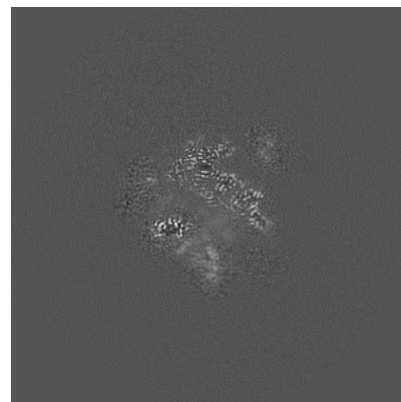
6.2.1 Primary map



X Index: 200

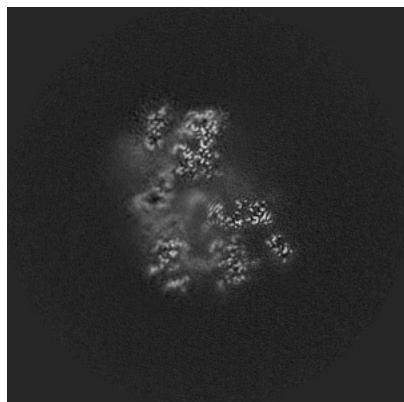


Y Index: 200

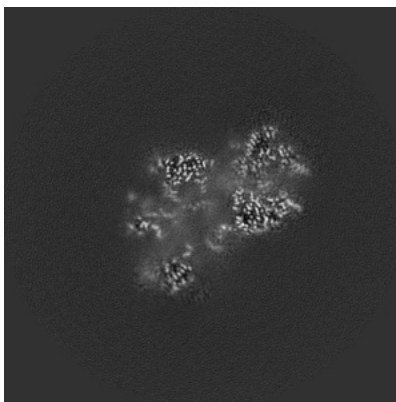


Z Index: 200

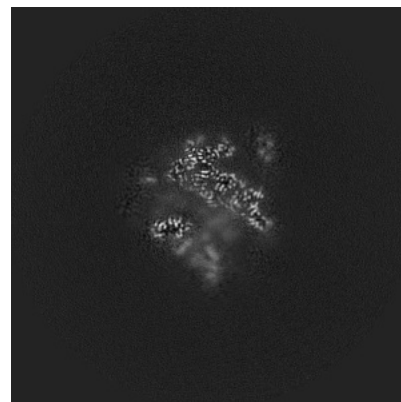
6.2.2 Raw map



X Index: 200



Y Index: 200

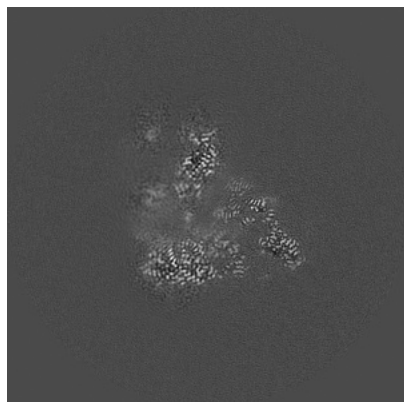


Z Index: 200

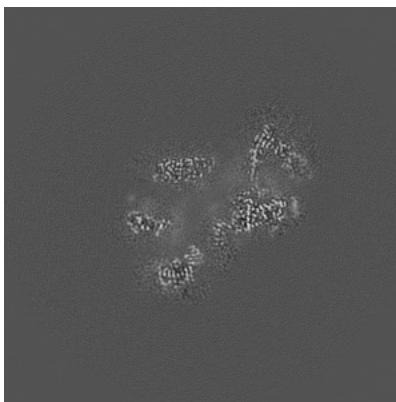
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

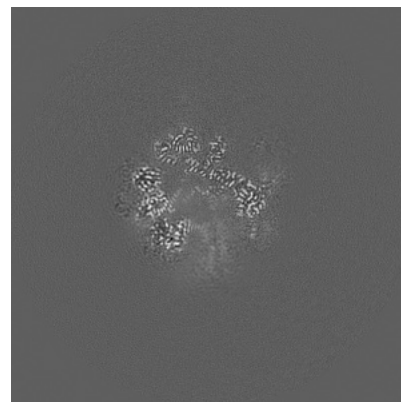
6.3.1 Primary map



X Index: 185

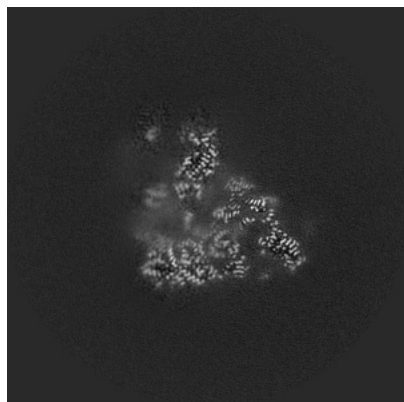


Y Index: 194

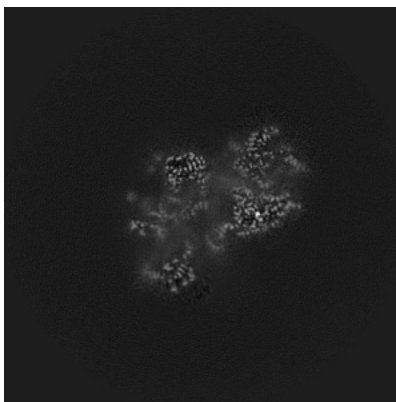


Z Index: 185

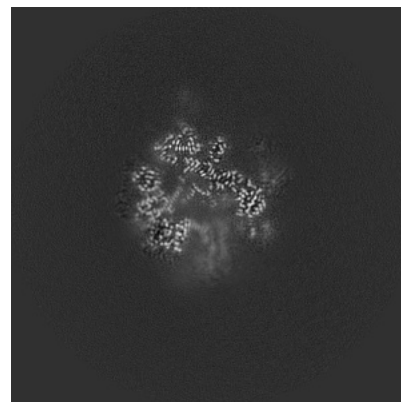
6.3.2 Raw map



X Index: 185



Y Index: 202

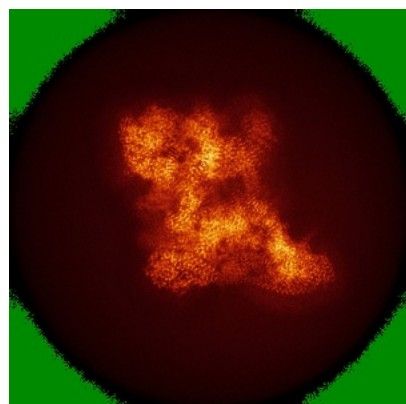


Z Index: 186

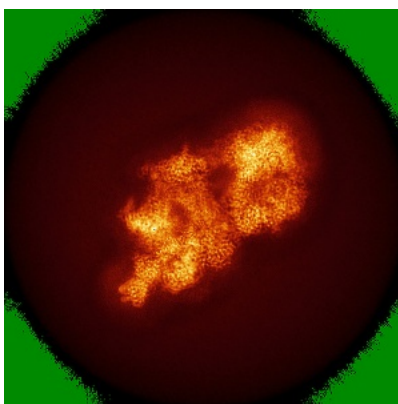
The images above show the largest variance slices of the map in three orthogonal directions.

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

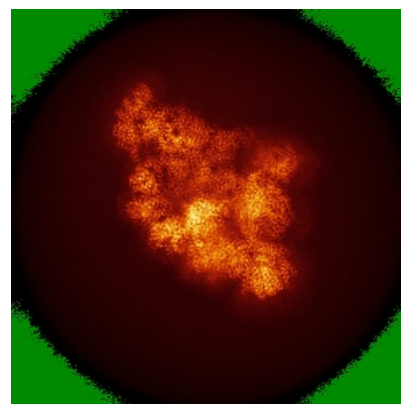
6.4.1 Primary map



X



Y

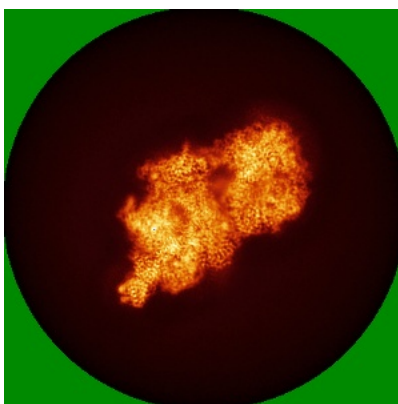


Z

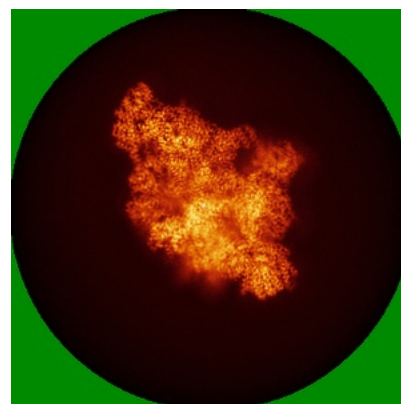
6.4.2 Raw map



X



Y

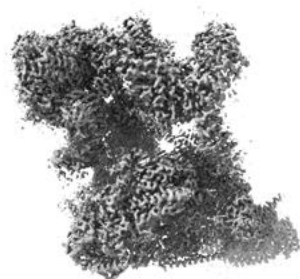


Z

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](#)

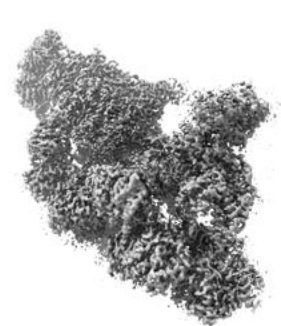
6.5.1 Primary map



X



Y



Z

The images above show the 3D surface view of the map at the recommended contour level 8.5. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

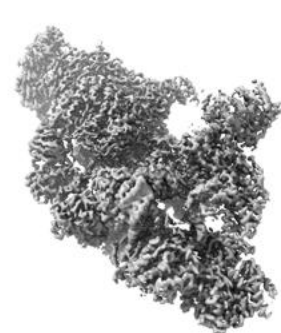
6.5.2 Raw map



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.

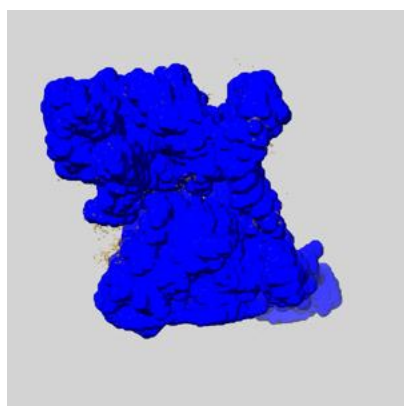
6.6 Mask visualisation [i](#)

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

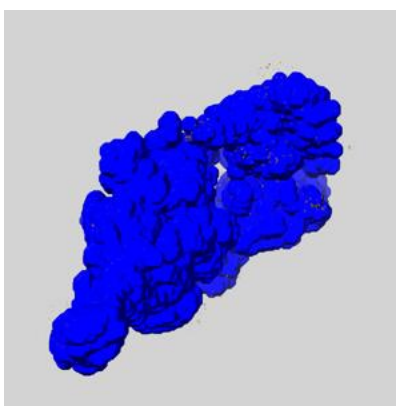
A mask typically either:

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

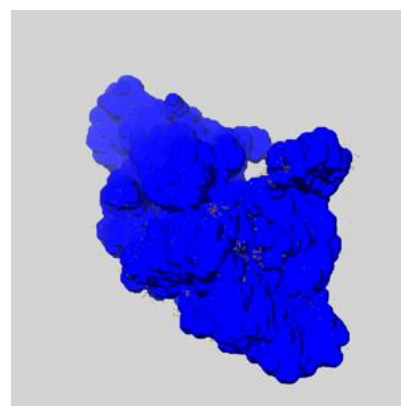
6.6.1 emd_26966_msk_1.map [i](#)



X



Y

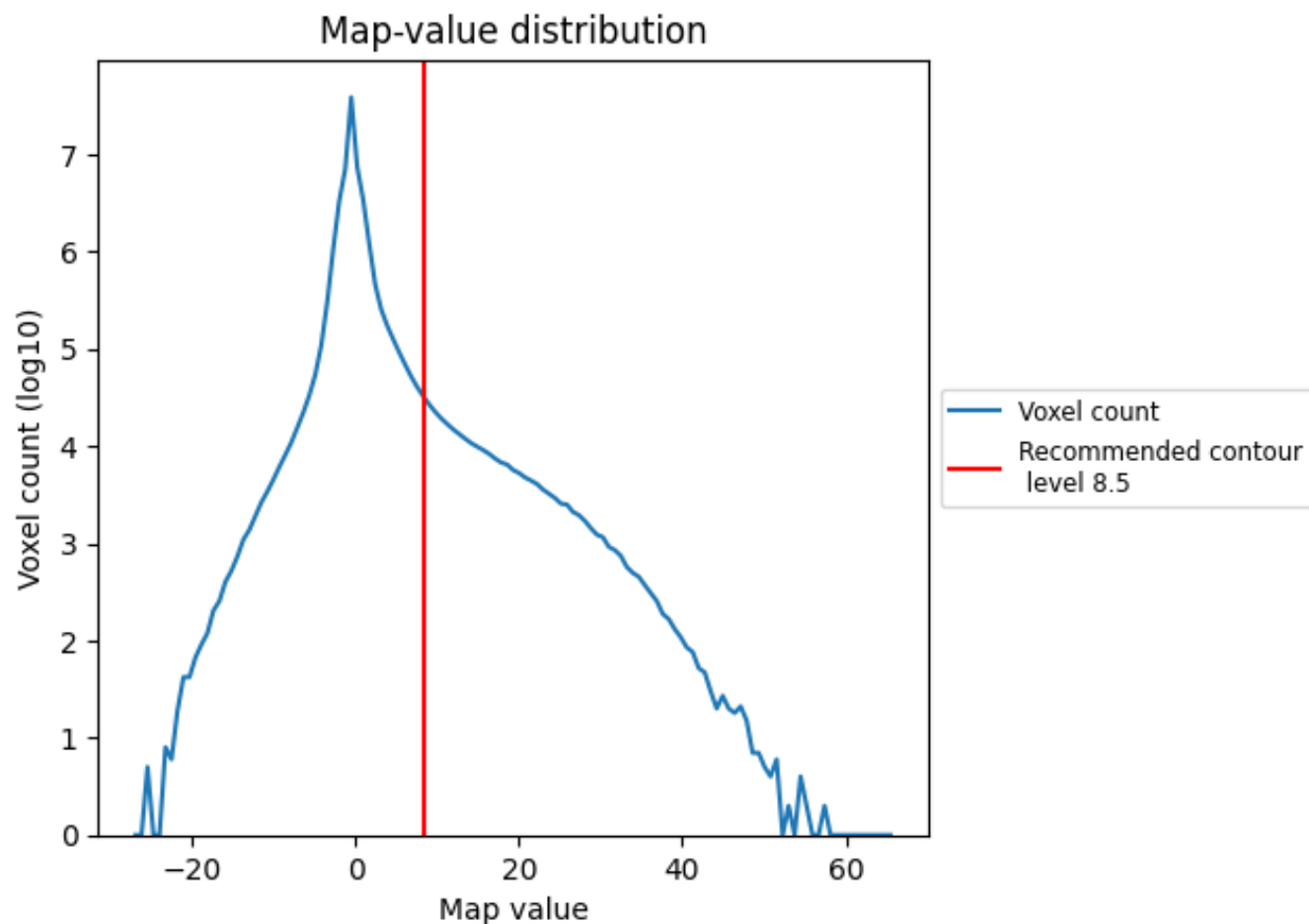


Z

7 Map analysis [i](#)

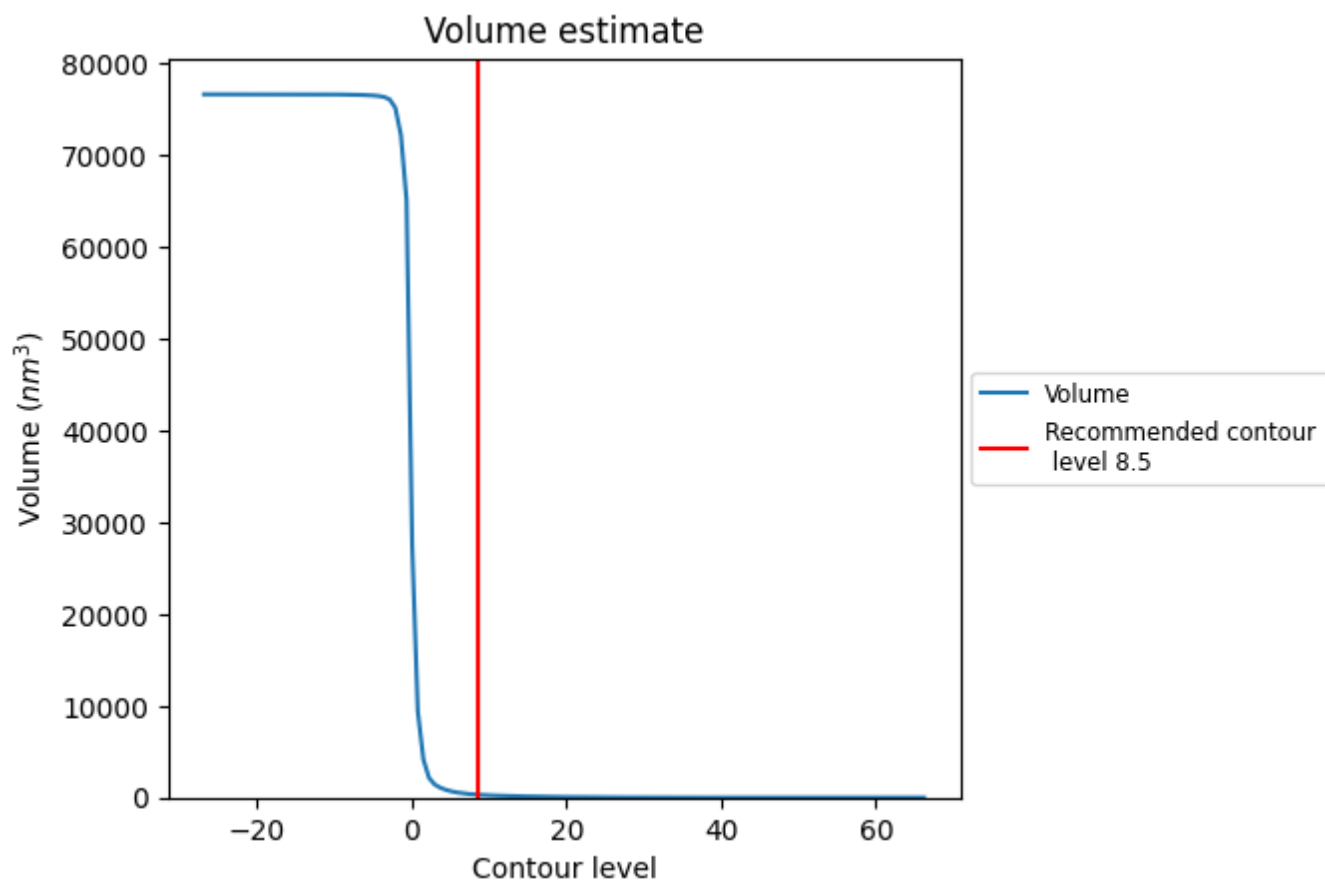
This section contains the results of statistical analysis of the map.

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.

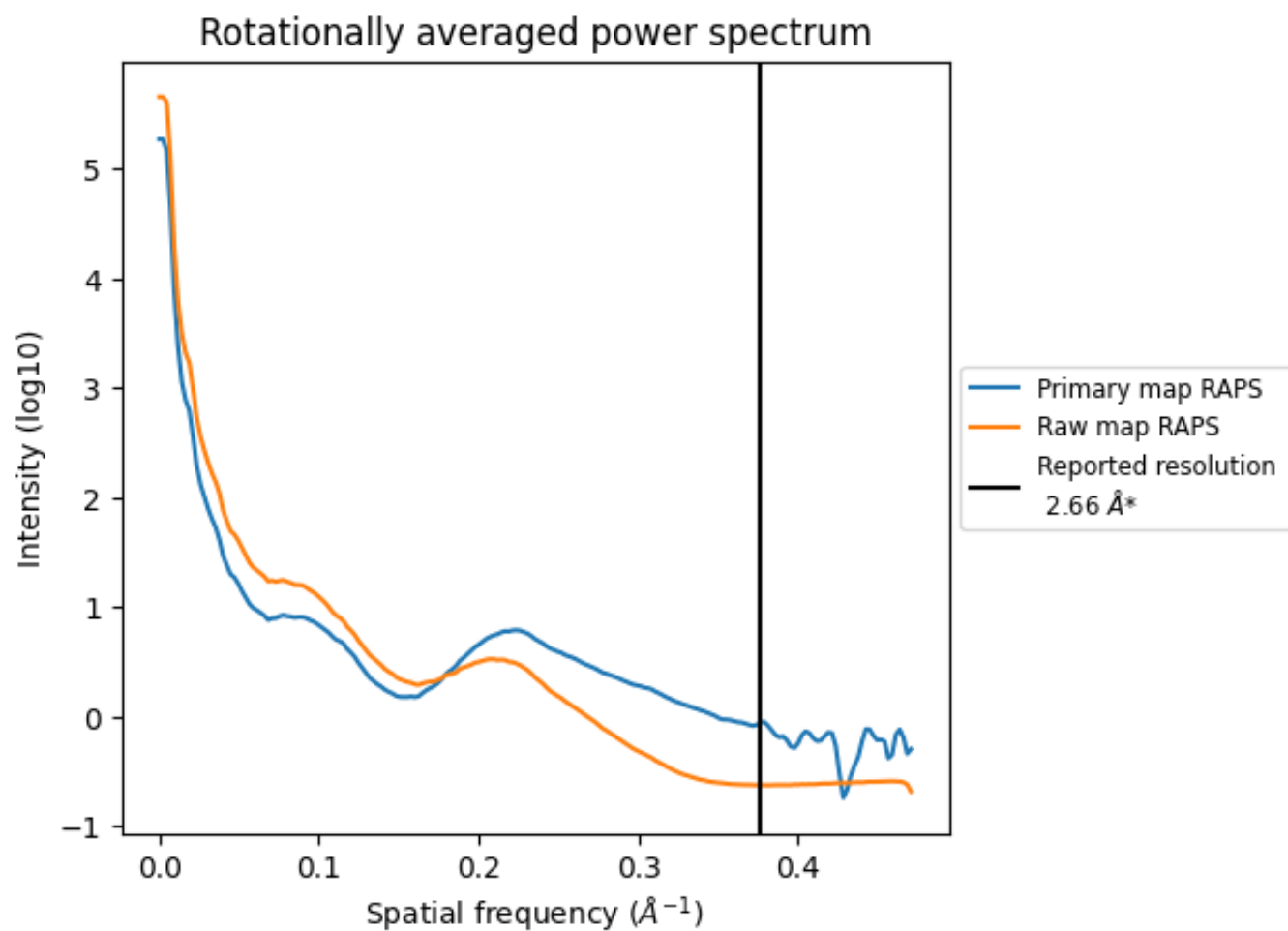
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 315 nm³; this corresponds to an approximate mass of 285 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 ⓘ

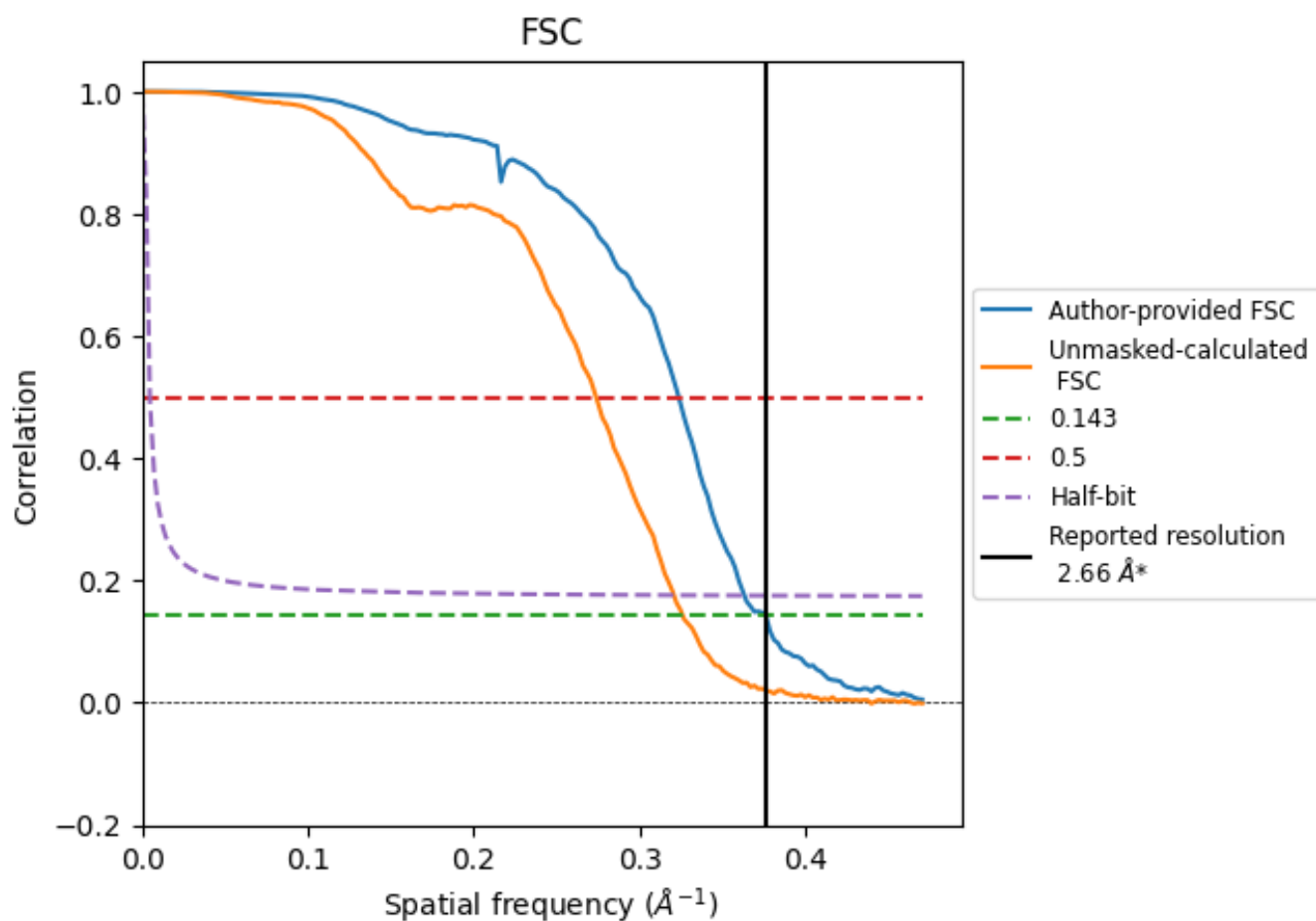


*Reported resolution corresponds to spatial frequency of 0.376 Å⁻¹

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 0.376 Å⁻¹

8.2 Resolution estimates [i](#)

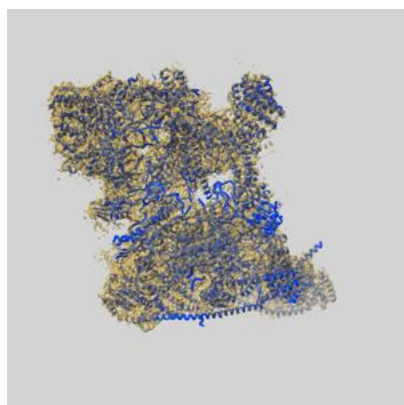
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.66	-	-
Author-provided FSC curve	2.66	3.09	2.75
Unmasked-calculated*	3.07	3.65	3.12

*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 3.07 differs from the reported value 2.66 by more than 10 %

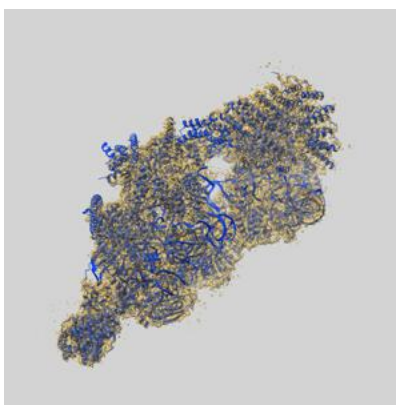
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-26966 and PDB model 8CSP. Per-residue inclusion information can be found in section [3](#) on page [13](#).

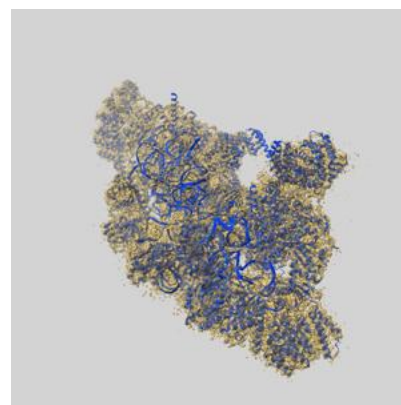
9.1 Map-model overlay [i](#)



X



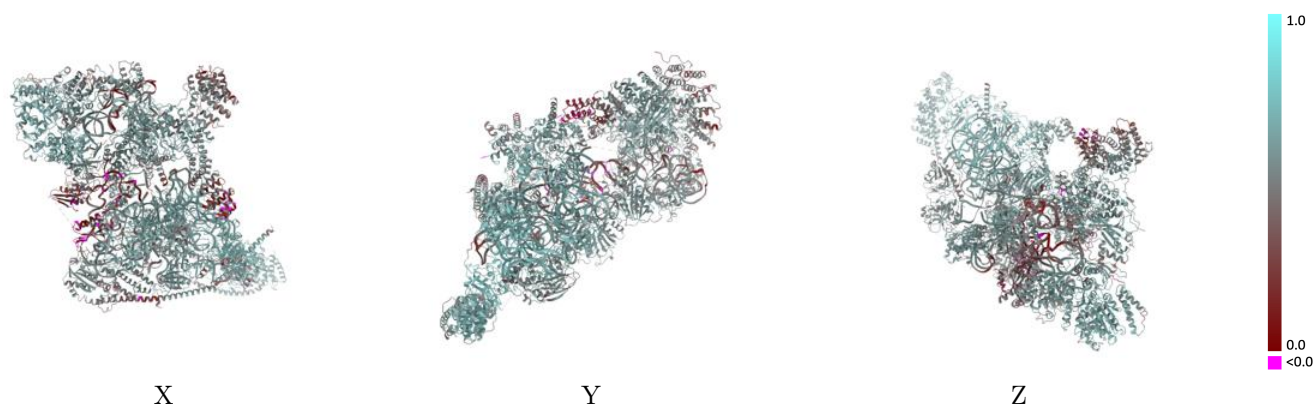
Y



Z

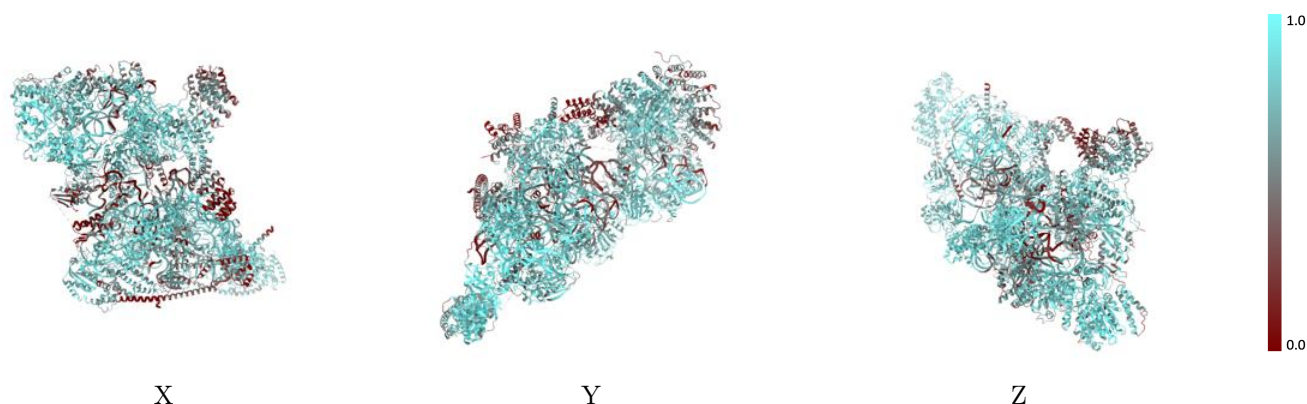
The images above show the 3D surface view of the map at the recommended contour level 8.5 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



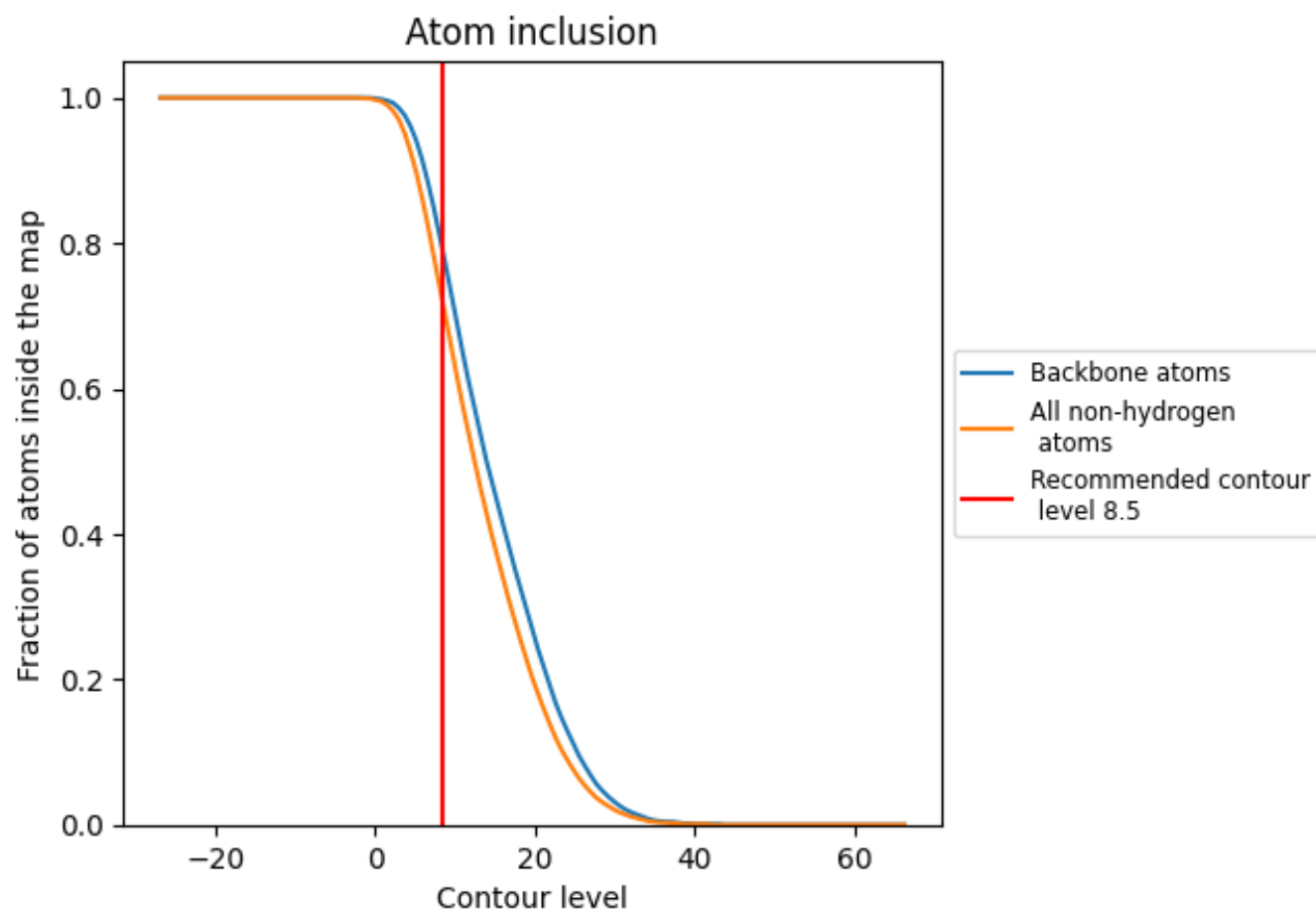
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

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



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (8.5).

9.4 Atom inclusion ⓘ



At the recommended contour level, 79% of all backbone atoms, 71% 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 (8.5) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.7140	 0.5480
0	 0.7930	 0.6370
1	 0.7310	 0.5570
4	 0.5240	 0.4470
5	 0.8610	 0.5750
7	 0.8520	 0.5670
8	 0.7250	 0.5320
9	 0.7320	 0.5100
A	 0.7140	 0.5260
B	 0.8760	 0.6230
C	 0.8070	 0.5810
D	 0.6980	 0.6110
E	 0.7980	 0.5240
F	 0.7780	 0.5580
G	 0.6860	 0.5470
H	 0.7530	 0.5760
J	 0.4840	 0.5870
K	 0.4960	 0.4670
L	 0.7660	 0.5230
M	 0.8680	 0.6740
N	 0.6130	 0.6040
O	 0.7640	 0.6500
P	 0.7470	 0.5070
Q	 0.1250	 0.4210
R	 0.6380	 0.6180
S	 0.6580	 0.5420
T	 0.6300	 0.6060
U	 0.5520	 0.5220
V	 0.8010	 0.6160
W	 0.7320	 0.5690
X	 0.8690	 0.6220
Y	 0.6530	 0.5240
Z	 0.4890	 0.4550
a	 0.3870	 0.2740

