



wwPDB EM Validation Summary Report ⓘ

Aug 4, 2026 – 11:17 PM EDT

PDB ID : 9CD6 / pdb_00009cd6
EMDB ID : EMD-45463
Title : Cryo-EM structure of respiratory supercomplex I in open state
Authors : Zhang, Z.; Maharjan, R.; Tringides, M.
Deposited on : 2024-06-24
Resolution : 3.20 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:

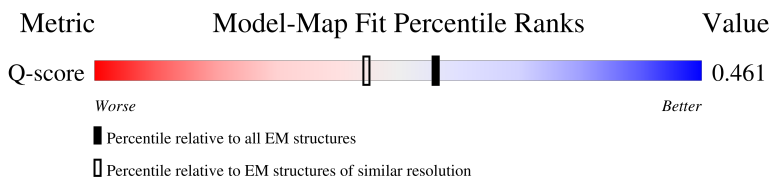
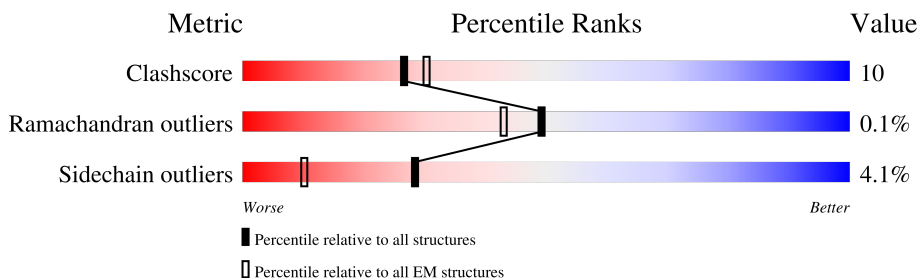
EMDB validation analysis : 0.0.1.dev133
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data 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) : 2.50

1 Overall quality at a glance

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

The reported resolution of this entry is 3.20 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.



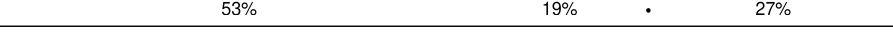
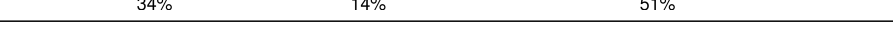
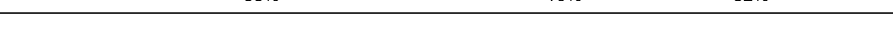
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	15020 (2.70 - 3.70)

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	B	464	
2	C	469	
3	D	264	
4	E	249	

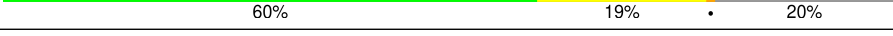
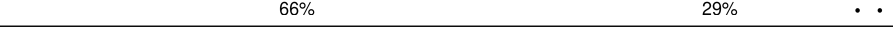
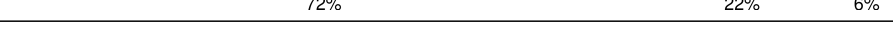
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Mol	Chain	Length	Quality of chain
5	F	123	
6	H	212	
7	I	196	
8	J	175	
9	K	145	
10	L	372	
11	N	116	
12	O	156	
12	X	156	
13	P	99	
14	Q	154	
15	R	110	
16	S	70	
17	T	169	
18	U	357	
19	V	141	
20	W	144	
21	Y	105	
22	Z	114	
23	a	189	
24	c	186	
25	d	176	
26	e	154	
27	f	76	
28	g	122	

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Mol	Chain	Length	Quality of chain
29	h	106	
30	i	347	
31	j	115	
32	k	98	
33	l	606	
34	m	175	
35	n	58	
36	o	129	
37	p	221	
38	q	459	
39	r	318	
40	s	249	
41	t	137	
42	G	727	
43	M	113	
44	b	128	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
46	SF4	B	502	-	-	X	-
46	SF4	I	301	-	-	X	-

2 Entry composition [i](#)

There are 49 unique types of molecules in this entry. The entry contains 65688 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 NADH dehydrogenase [ubiquinone] flavoprotein 1, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	B	429	Total	C	N	O	S	0	0
			3300	2084	587	609	20		

- Molecule 2 is a protein called NADH dehydrogenase [ubiquinone] iron-sulfur protein 2, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	C	430	Total	C	N	O	S	0	0
			3453	2207	592	630	24		

- Molecule 3 is a protein called NADH dehydrogenase [ubiquinone] iron-sulfur protein 3, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	D	208	Total	C	N	O	S	0	0
			1737	1123	298	314	2		

- Molecule 4 is a protein called NADH dehydrogenase [ubiquinone] flavoprotein 2, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	E	214	Total	C	N	O	S	0	0
			1658	1058	278	312	10		

- Molecule 5 is a protein called NADH dehydrogenase [ubiquinone] iron-sulfur protein 6, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	F	93	Total	C	N	O	S	0	0
			720	439	135	143	3		

- Molecule 6 is a protein called NADH dehydrogenase [ubiquinone] iron-sulfur protein 8, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	H	176	Total	C	N	O	S	0	0
			1412	887	243	269	13		

- Molecule 7 is a protein called NADH dehydrogenase [ubiquinone] iron-sulfur protein 7, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	I	156	Total	C	N	O	S	0	0
			1249	794	227	214	14		

- Molecule 8 is a protein called NADH dehydrogenase [ubiquinone] iron-sulfur protein 4, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	J	118	Total	C	N	O	S	0	0
			963	608	173	179	3		

- Molecule 9 is a protein called NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 12.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	K	144	Total	C	N	O	S	0	0
			1203	769	217	212	5		

- Molecule 10 is a protein called NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 9, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	L	328	Total	C	N	O	S	0	0
			2625	1696	458	463	8		

- Molecule 11 is a protein called NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 5 isoform X1.

Mol	Chain	Residues	Atoms					AltConf	Trace
11	N	112	Total	C	N	O	S	0	0
			911	588	154	166	3		

- Molecule 12 is a protein called Acyl carrier protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
12	O	83	Total	C	N	O	S	0	0
			668	431	99	133	5		

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Mol	Chain	Residues	Atoms					AltConf	Trace
12	X	85	Total	C	N	O	S	0	0
			689	445	101	138	5		

- Molecule 13 is a protein called NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
13	P	83	Total	C	N	O	S	0	0
			669	419	125	123	2		

- Molecule 14 is a protein called NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 6.

Mol	Chain	Residues	Atoms					AltConf	Trace
14	Q	112	Total	C	N	O	S	0	0
			955	610	176	164	5		

- Molecule 15 is a protein called NADH dehydrogenase [ubiquinone] flavoprotein 3, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
15	R	35	Total	C	N	O	S	0	0
			295	185	55	54	1		

- Molecule 16 is a protein called NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
16	S	70	Total	C	N	O	S	0	0
			562	361	101	94	6		

- Molecule 17 is a protein called NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 3.

Mol	Chain	Residues	Atoms					AltConf	Trace
17	T	82	Total	C	N	O	S	0	0
			638	414	109	114	1		

- Molecule 18 is a protein called NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 10, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
18	U	318	Total	C	N	O	S	0	0
			2573	1638	437	488	10		

- Molecule 19 is a protein called NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 11.

Mol	Chain	Residues	Atoms					AltConf	Trace
19	V	140	Total	C	N	O	S	0	0
			1021	651	174	190	6		

- Molecule 20 is a protein called NADH:ubiquinone oxidoreductase subunit A13.

Mol	Chain	Residues	Atoms					AltConf	Trace
20	W	140	Total	C	N	O	S	0	0
			1162	749	201	203	9		

- Molecule 21 is a protein called NADH:ubiquinone oxidoreductase subunit B2.

Mol	Chain	Residues	Atoms					AltConf	Trace
21	Y	66	Total	C	N	O	S	0	0
			571	378	94	98	1		

- Molecule 22 is a protein called NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 3.

Mol	Chain	Residues	Atoms					AltConf	Trace
22	Z	77	Total	C	N	O	S	0	0
			620	407	104	108	1		

- Molecule 23 is a protein called NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 5, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
23	a	138	Total	C	N	O	S	0	0
			1151	754	195	199	3		

- Molecule 24 is a protein called NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 8, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
24	c	153	Total	C	N	O	S	0	0
			1291	838	208	237	8		

- Molecule 25 is a protein called NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 10.

Mol	Chain	Residues	Atoms					AltConf	Trace
25	d	169	Total	C	N	O	S	0	0
			1426	895	259	264	8		

- Molecule 26 is a protein called NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 11, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
26	e	99	Total	C	N	O	S	0	0
			826	530	137	155	4		

- Molecule 27 is a protein called NADH dehydrogenase [ubiquinone] 1 subunit C1, mitochondrial.

Mol	Chain	Residues	Atoms				AltConf	Trace
27	f	46	Total	C	N	O	0	0
			391	259	67	65		

- Molecule 28 is a protein called NADH dehydrogenase [ubiquinone] 1 subunit C2.

Mol	Chain	Residues	Atoms					AltConf	Trace
28	g	121	Total	C	N	O	S	0	0
			1000	650	173	171	6		

- Molecule 29 is a protein called NADH dehydrogenase [ubiquinone] iron-sulfur protein 5.

Mol	Chain	Residues	Atoms					AltConf	Trace
29	h	105	Total	C	N	O	S	0	0
			867	550	161	150	6		

- Molecule 30 is a protein called NADH-ubiquinone oxidoreductase chain 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
30	i	347	Total	C	N	O	S	0	0
			2710	1782	420	462	46		

- Molecule 31 is a protein called NADH-ubiquinone oxidoreductase chain 3.

Mol	Chain	Residues	Atoms					AltConf	Trace
31	j	106	Total	C	N	O	S	0	0
			845	573	122	144	6		

- Molecule 32 is a protein called NADH-ubiquinone oxidoreductase chain 4L.

Mol	Chain	Residues	Atoms					AltConf	Trace
32	k	98	Total	C	N	O	S	0	0
			748	493	113	128	14		

- Molecule 33 is a protein called NADH-ubiquinone oxidoreductase chain 5.

Mol	Chain	Residues	Atoms					AltConf	Trace
33	l	603	Total	C	N	O	S	0	0
			4785	3174	741	819	51		

- Molecule 34 is a protein called NADH-ubiquinone oxidoreductase chain 6.

Mol	Chain	Residues	Atoms					AltConf	Trace
34	m	173	Total	C	N	O	S	0	0
			1321	888	187	234	12		

- Molecule 35 is a protein called NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
35	n	56	Total	C	N	O	S	0	0
			479	311	88	79	1		

- Molecule 36 is a protein called NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 4.

Mol	Chain	Residues	Atoms					AltConf	Trace
36	o	126	Total	C	N	O		0	0
			1041	677	179	185			

- Molecule 37 is a protein called NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 9.

Mol	Chain	Residues	Atoms					AltConf	Trace
37	p	177	Total	C	N	O	S	0	0
			1529	979	278	264	8		

- Molecule 38 is a protein called NADH-ubiquinone oxidoreductase chain 4.

Mol	Chain	Residues	Atoms					AltConf	Trace
38	q	459	Total	C	N	O	S	0	0
			3630	2410	572	609	39		

- Molecule 39 is a protein called NADH-ubiquinone oxidoreductase chain 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
39	r	309	Total	C	N	O	S	0	0
			2437	1632	376	408	21		

- Molecule 40 is a protein called NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 8.

Mol	Chain	Residues	Atoms					AltConf	Trace
40	s	171	Total	C	N	O	S	0	0
			1398	887	250	251	10		

- Molecule 41 is a protein called NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 7.

Mol	Chain	Residues	Atoms					AltConf	Trace
41	t	117	Total	C	N	O	S	0	0
			1014	632	193	180	9		

- Molecule 42 is a protein called NADH-ubiquinone oxidoreductase 75 kDa subunit, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
42	G	684	Total	C	N	O	S	0	0
			5260	3298	917	1006	39		

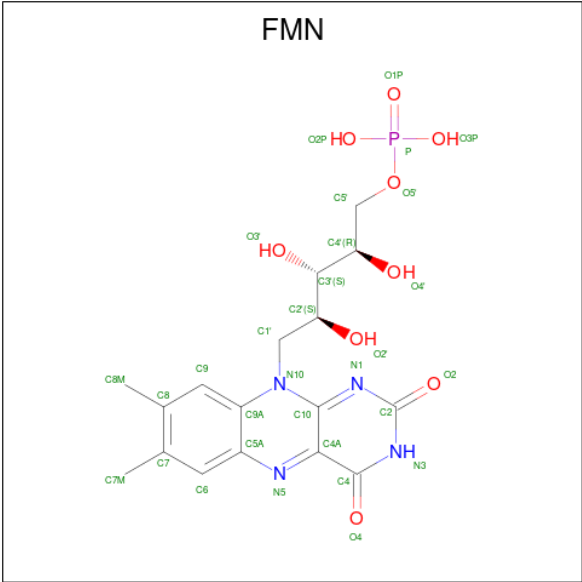
- Molecule 43 is a protein called NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 7.

Mol	Chain	Residues	Atoms					AltConf	Trace
43	M	96	Total	C	N	O	S	0	0
			774	487	146	138	3		

- Molecule 44 is a protein called NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 6.

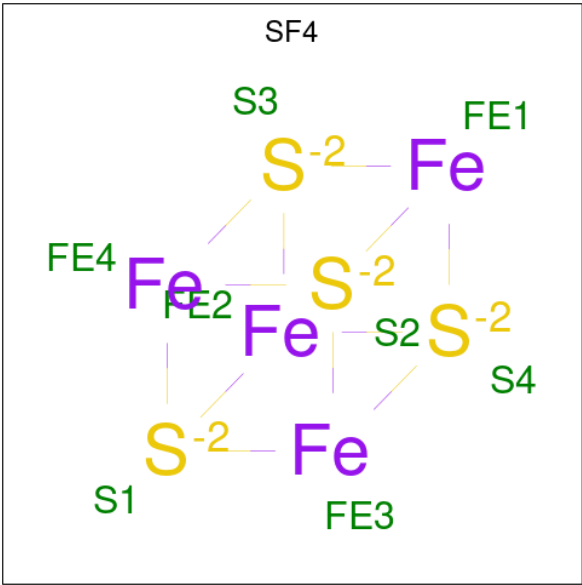
Mol	Chain	Residues	Atoms					AltConf	Trace
44	b	111	Total	C	N	O	S	0	0
			946	623	163	159	1		

- Molecule 45 is FLAVIN MONONUCLEOTIDE (CCD ID: FMN) (formula: C₁₇H₂₁N₄O₉P).



Mol	Chain	Residues	Atoms					AltConf
45	B	1	Total	C	N	O	P	0
			31	17	4	9	1	

- Molecule 46 is IRON/SULFUR CLUSTER (CCD ID: SF4) (formula: Fe₄S₄).



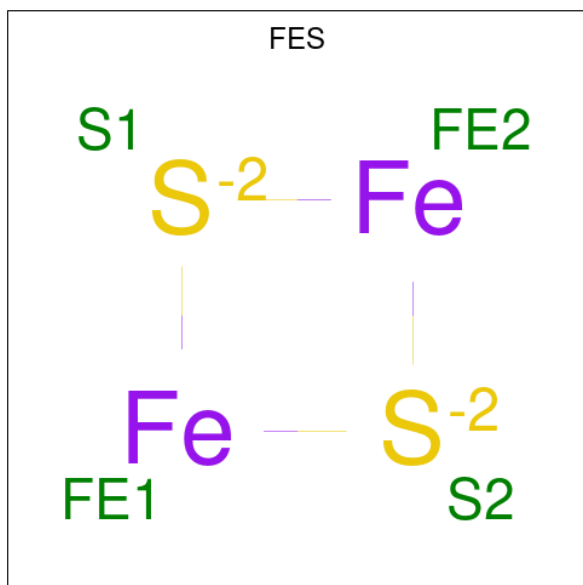
Mol	Chain	Residues	Atoms			AltConf
46	B	1	Total	Fe	S	0
			8	4	4	
46	H	1	Total	Fe	S	0
			8	4	4	
46	H	1	Total	Fe	S	0
			8	4	4	

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Mol	Chain	Residues	Atoms			AltConf
46	I	1	Total	Fe	S	0
			8	4	4	
46	G	1	Total	Fe	S	0
			8	4	4	
46	G	1	Total	Fe	S	0
			8	4	4	

- Molecule 47 is FE2/S2 (INORGANIC) CLUSTER (CCD ID: FES) (formula: Fe_2S_2).



Mol	Chain	Residues	Atoms			AltConf
47	E	1	Total	Fe	S	0
			4	2	2	
47	G	1	Total	Fe	S	0
			4	2	2	

- Molecule 48 is NADPH DIHYDRO-NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (CCD ID: NDP) (formula: $\text{C}_{21}\text{H}_{30}\text{N}_7\text{O}_{17}\text{P}_3$).



Mol	Chain	Residues	Atoms					AltCon
48	L	1	Total	C	N	O	P	0
			48	21	7	17	3	

- Molecule 49 is S-[2-($\{N-[(2S)-2\text{-hydroxy-3,3-dimethyl-4-(phosphonooxy)butanoyl]}\text{-beta-alanine}\}$)amino)ethyl] tetradecanethioate (CCD ID: ZMP) (formula: $C_{25}H_{49}N_2O_8PS$).

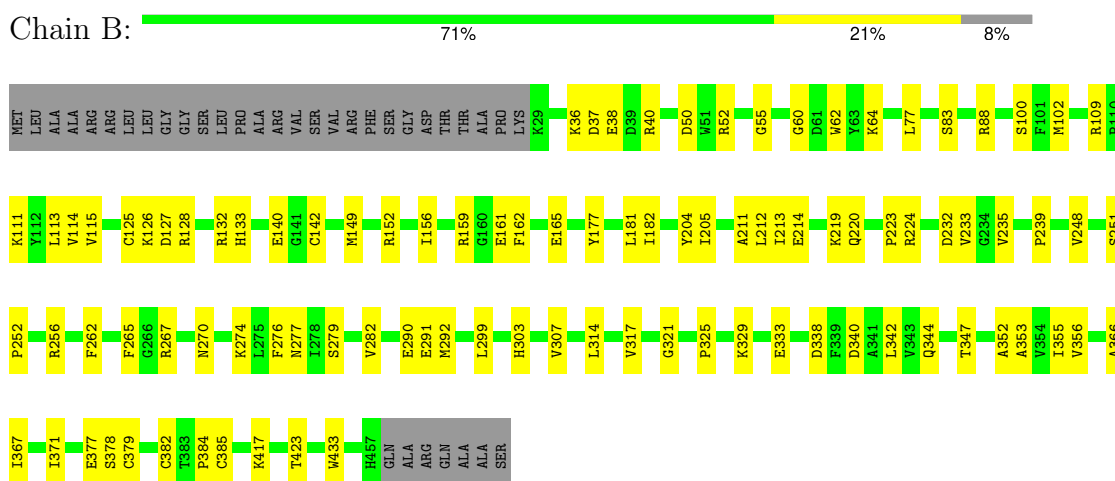


Mol	Chain	Residues	Atoms						AltCon
49	Q	1	Total	C	N	O	P	S	0
			30	18	2	8	1	1	

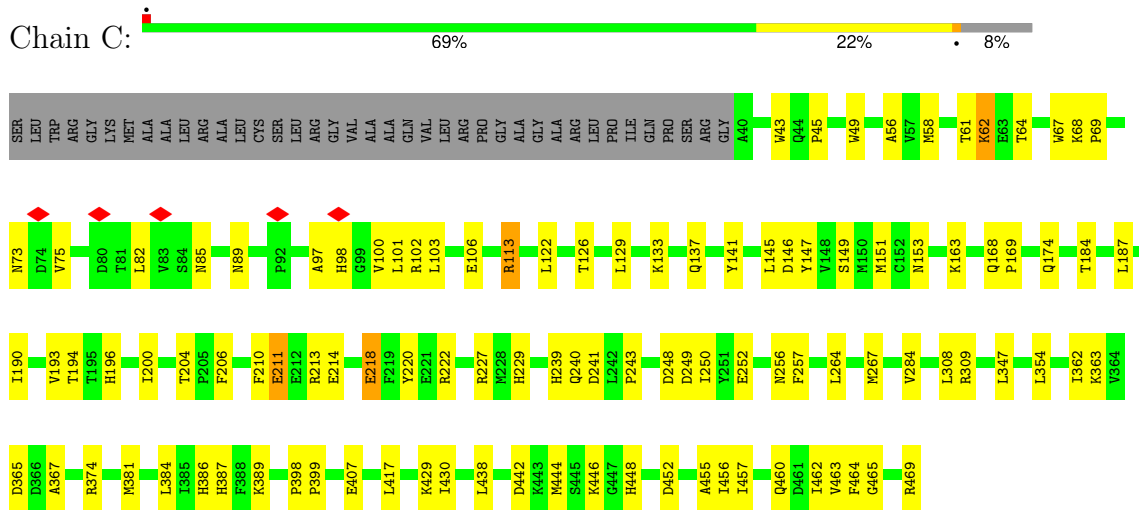
3 Residue-property plots

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: NADH dehydrogenase [ubiquinone] flavoprotein 1, mitochondrial

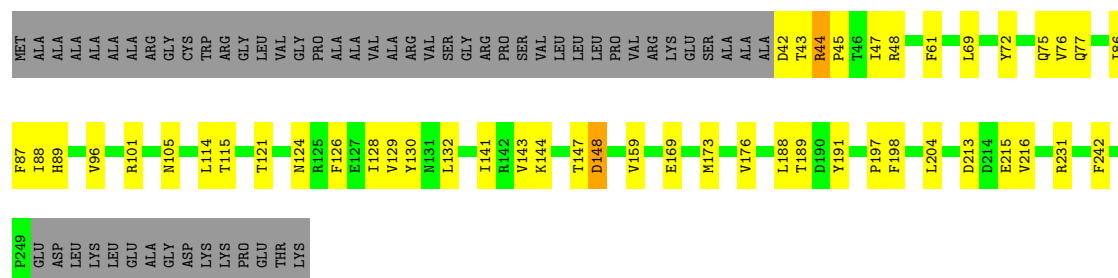


- Molecule 2: NADH dehydrogenase [ubiquinone] iron-sulfur protein 2, mitochondrial



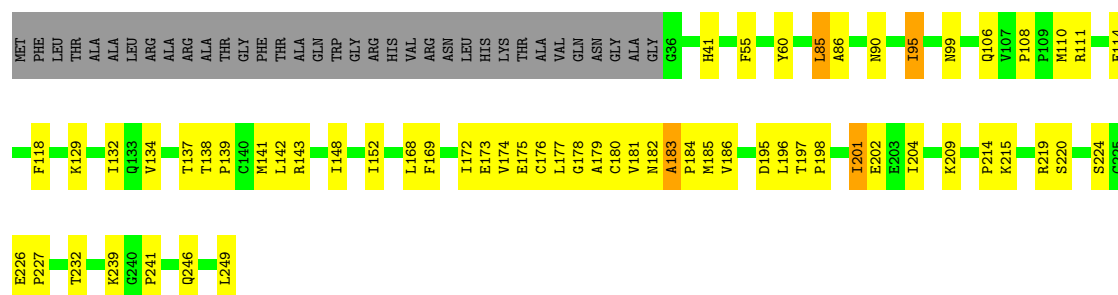
- Molecule 3: NADH dehydrogenase [ubiquinone] iron-sulfur protein 3, mitochondrial





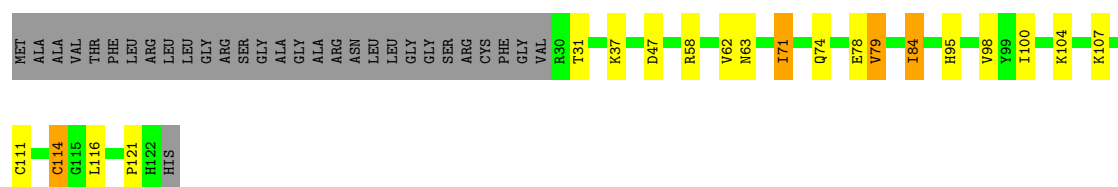
- Molecule 4: NADH dehydrogenase [ubiquinone] flavoprotein 2, mitochondrial

Chain E:



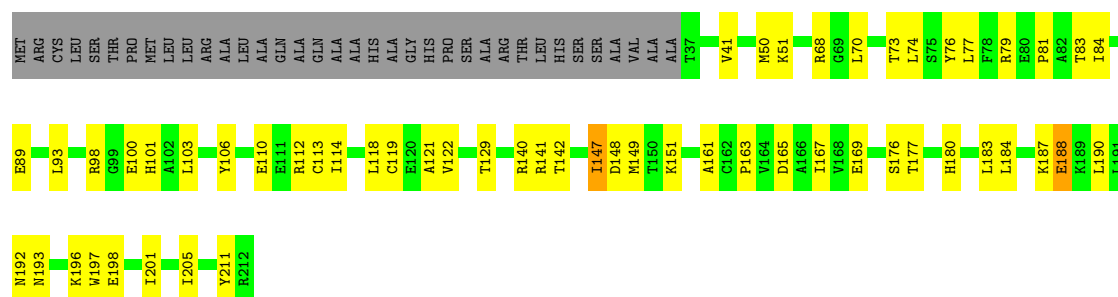
- Molecule 5: NADH dehydrogenase [ubiquinone] iron-sulfur protein 6, mitochondrial

Chain F:



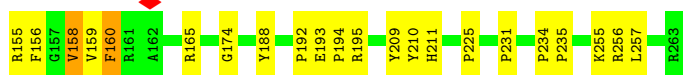
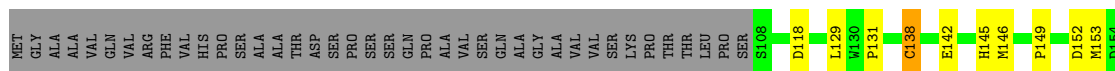
- Molecule 6: NADH dehydrogenase [ubiquinone] iron-sulfur protein 8, mitochondrial

Chain H:

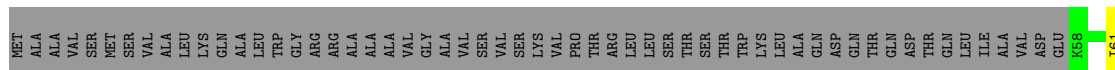


- Molecule 7: NADH dehydrogenase [ubiquinone] iron-sulfur protein 7, mitochondrial

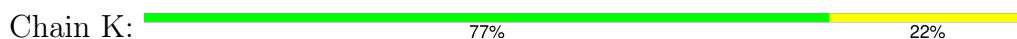
Chain I:



- Molecule 8: NADH dehydrogenase [ubiquinone] iron-sulfur protein 4, mitochondrial



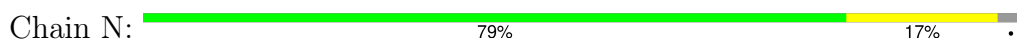
- Molecule 9: NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 12



- Molecule 10: NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 9, mitochondrial



- Molecule 11: NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 5 isoform X1



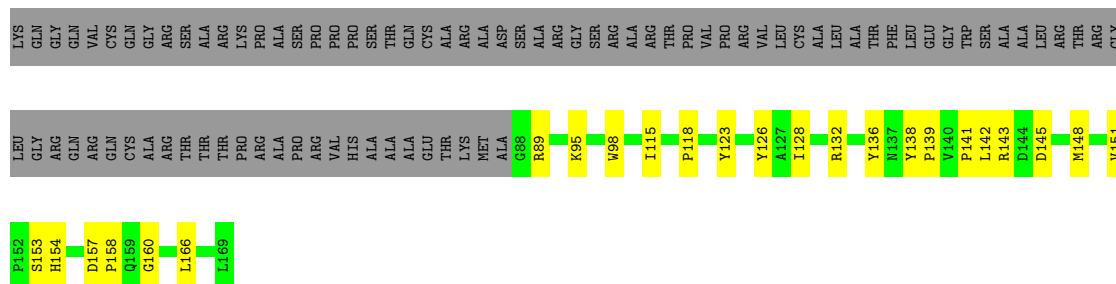
- Molecule 12: Acyl carrier protein

Chain S:  74% 23% .



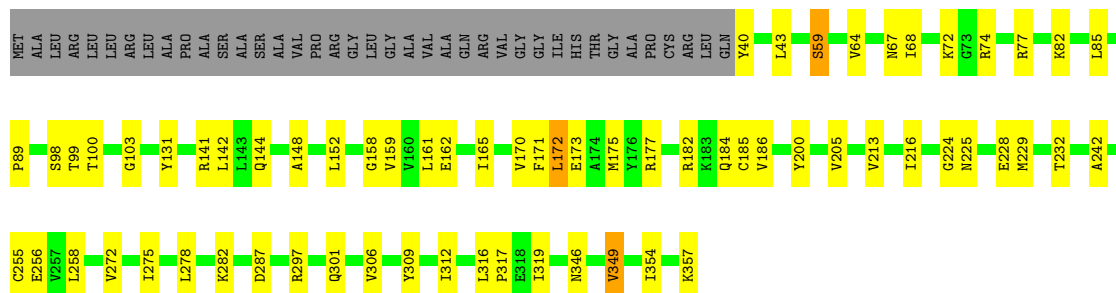
- Molecule 17: NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 3

Chain T:  34% 14% 51%



- Molecule 18: NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 10, mitochondrial

Chain U:  70% 18% 11%



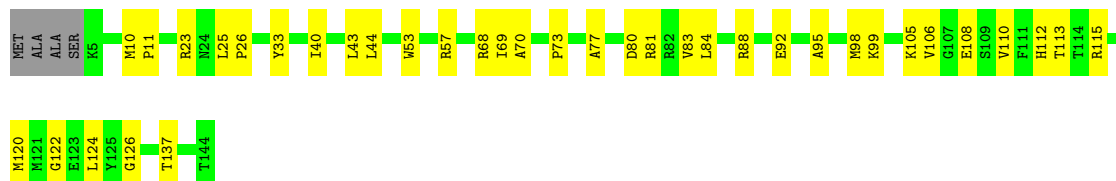
- Molecule 19: NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 11

Chain V:  87% 12% .



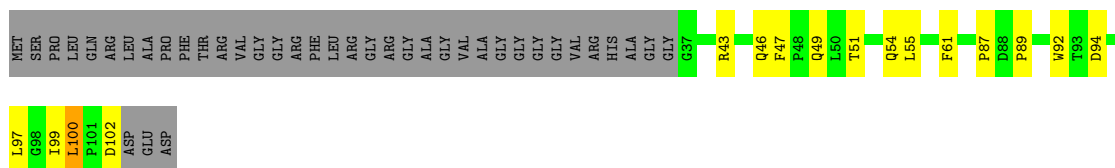
- Molecule 20: NADH:ubiquinone oxidoreductase subunit A13

Chain W:  72% 26% .



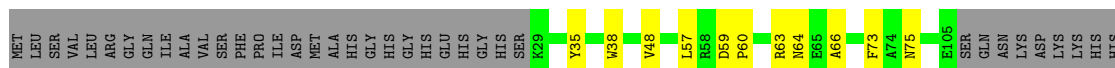
- Molecule 21: NADH:ubiquinone oxidoreductase subunit B2

Chain Y: 



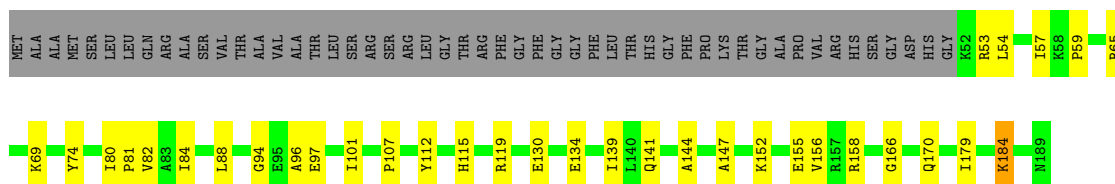
- Molecule 22: NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 3

Chain Z: 



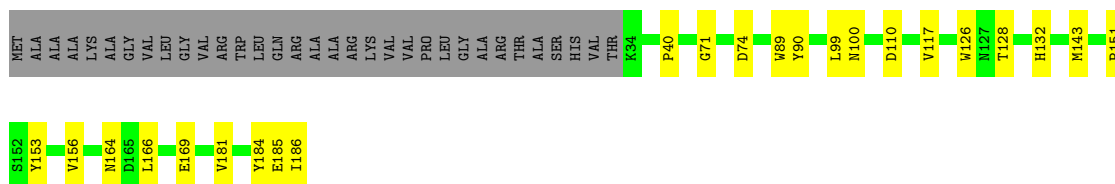
- Molecule 23: NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 5, mitochondrial

Chain a: 



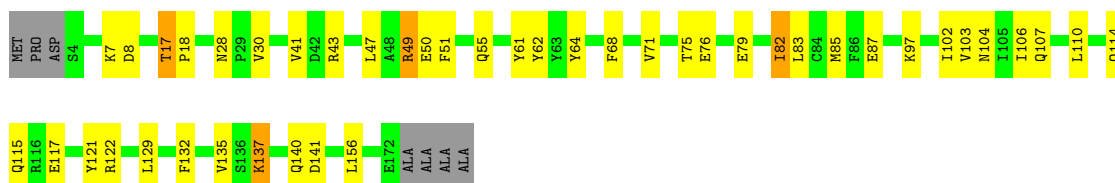
- Molecule 24: NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 8, mitochondrial

Chain c: 



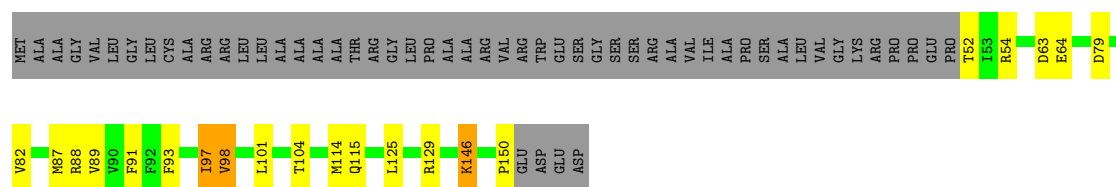
- Molecule 25: NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 10

Chain d: 

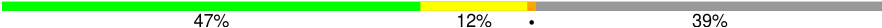


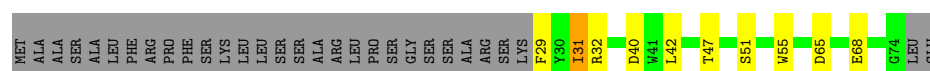
- Molecule 26: NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 11, mitochondrial

Chain e: 



- Molecule 27: NADH dehydrogenase [ubiquinone] 1 subunit C1, mitochondrial

Chain f: 



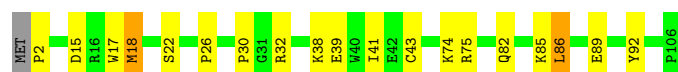
- Molecule 28: NADH dehydrogenase [ubiquinone] 1 subunit C2

Chain g: 



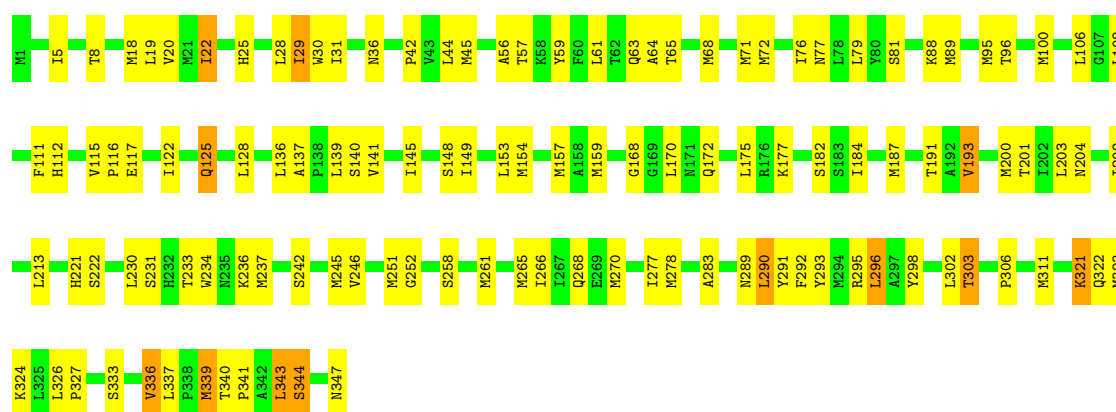
- Molecule 29: NADH dehydrogenase [ubiquinone] iron-sulfur protein 5

Chain h: 



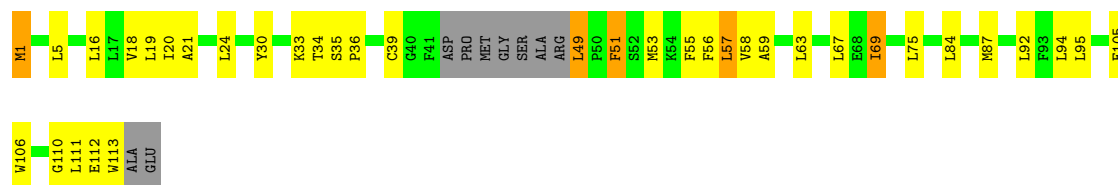
- Molecule 30: NADH-ubiquinone oxidoreductase chain 2

Chain i: 



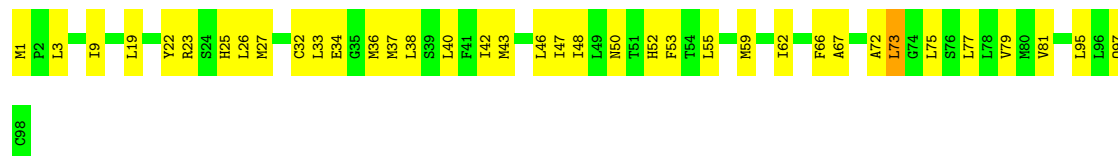
- Molecule 31: NADH-ubiquinone oxidoreductase chain 3

Chain j: 



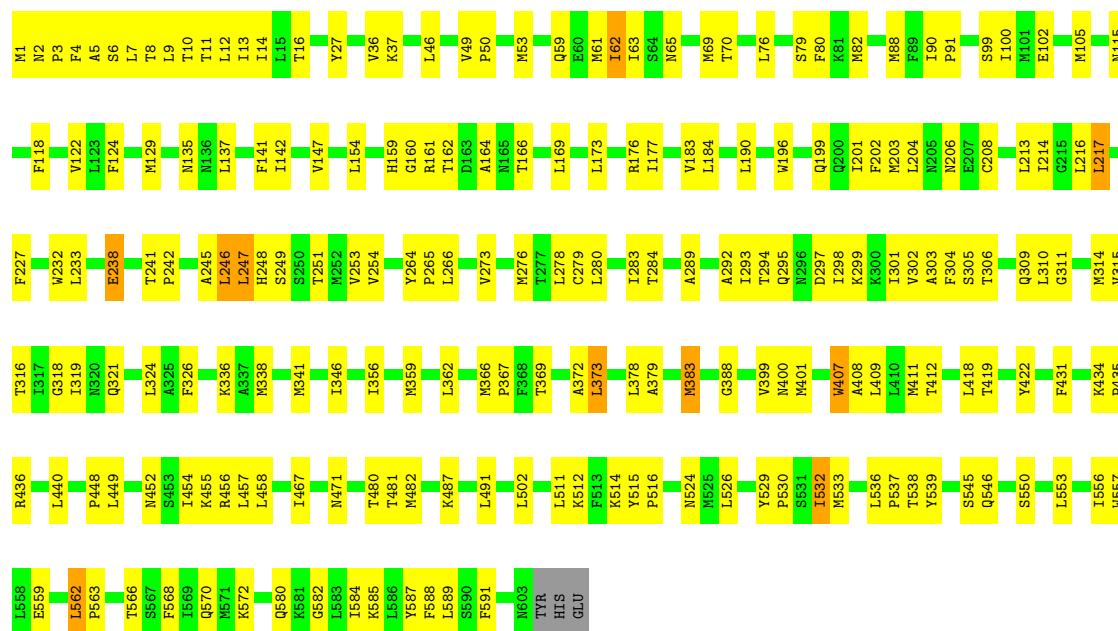
- Molecule 32: NADH-ubiquinone oxidoreductase chain 4L

Chain k: 62% 37%



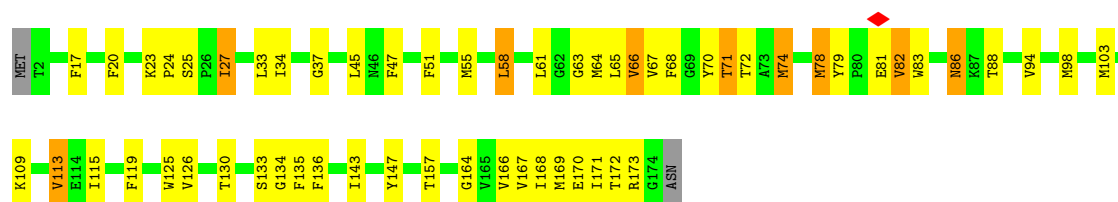
- Molecule 33: NADH-ubiquinone oxidoreductase chain 5

Chain l: 65% 33%



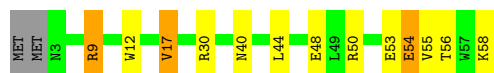
- Molecule 34: NADH-ubiquinone oxidoreductase chain 6

Chain m: 66% 28% 5%



- Molecule 35: NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 1

Chain n:  74% 17% 5% .



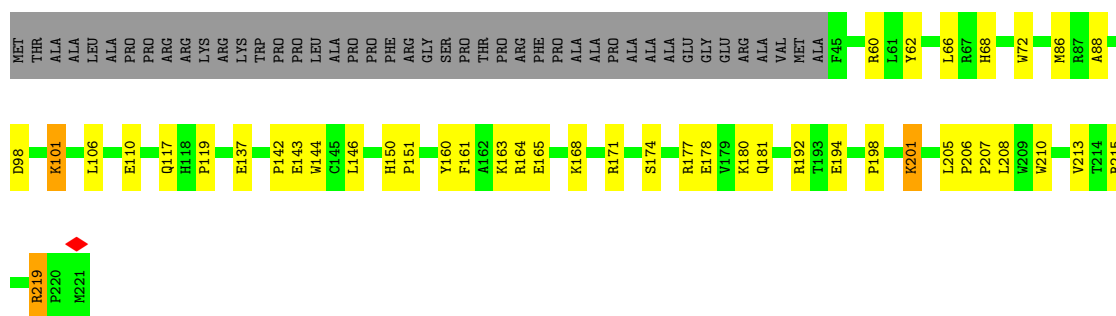
- Molecule 36: NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 4

Chain o:  74% 22% . .



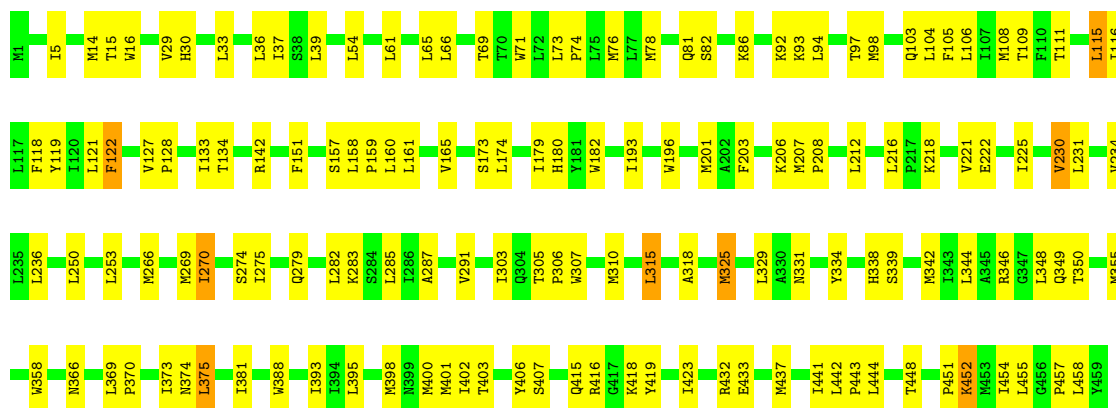
- Molecule 37: NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 9

Chain p:  60% 19% . 20%



- Molecule 38: NADH-ubiquinone oxidoreductase chain 4

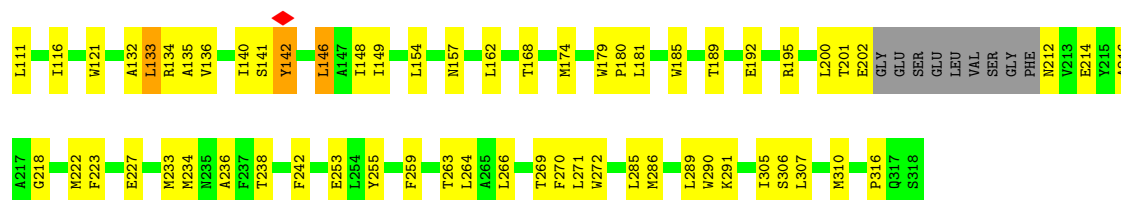
Chain q:  68% 30% .



- Molecule 39: NADH-ubiquinone oxidoreductase chain 1

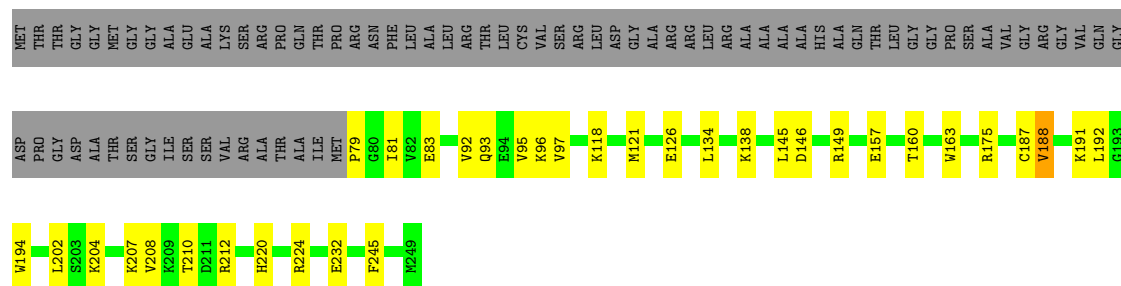
Chain r:  66% 29% . .





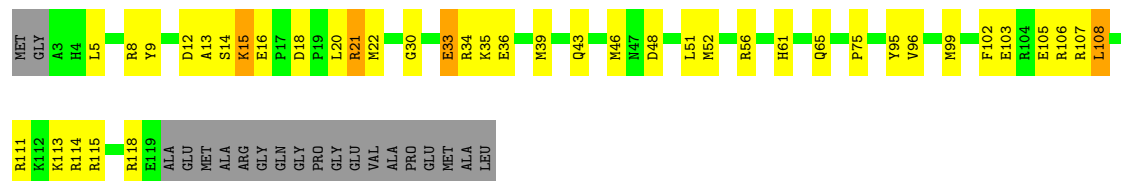
- Molecule 40: NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 8

Chain s:



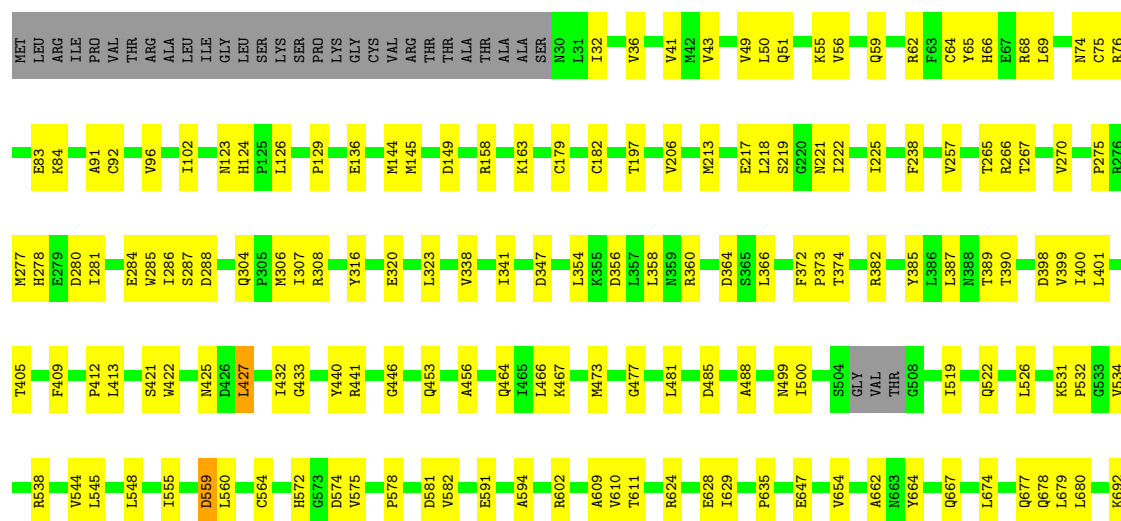
- Molecule 41: NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 7

Chain t:



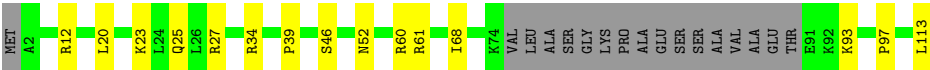
- Molecule 42: NADH-ubiquinone oxidoreductase 75 kDa subunit, mitochondrial

Chain G:

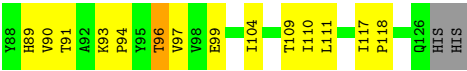
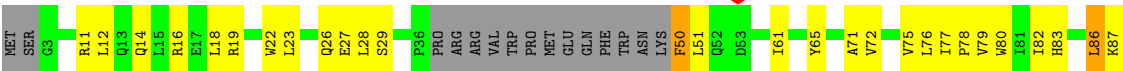




- Molecule 43: NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 7



- Molecule 44: NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 6



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	16982	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 ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	1500	Depositor
Magnification	81000	Depositor
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	2.035	Depositor
Minimum map value	-0.637	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.036	Depositor
Recommended contour level	0.1	Depositor
Map size (Å)	547.84, 547.84, 547.84	wwPDB
Map dimensions	512, 512, 512	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.07, 1.07, 1.07	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SF4, NDP, ZMP, FES, FMN

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	B	0.22	0/3375	0.36	0/4561
2	C	0.27	0/3545	0.39	0/4805
3	D	0.12	0/1788	0.28	0/2435
4	E	0.30	0/1698	0.45	0/2311
5	F	0.11	0/733	0.23	0/988
6	H	0.15	0/1443	0.30	0/1952
7	I	0.28	0/1280	0.45	0/1730
8	J	0.10	0/986	0.25	0/1329
9	K	0.11	0/1244	0.25	0/1693
10	L	0.35	0/2694	0.47	0/3649
11	N	0.13	0/930	0.27	0/1258
12	O	0.11	0/679	0.25	0/916
12	X	0.14	0/701	0.30	0/946
13	P	0.13	0/680	0.33	0/916
14	Q	0.10	0/979	0.23	0/1317
15	R	0.48	0/304	0.58	0/410
16	S	0.14	0/577	0.32	0/777
17	T	0.13	0/659	0.29	0/905
18	U	0.31	0/2633	0.42	0/3565
19	V	0.12	0/1042	0.26	0/1411
20	W	0.14	0/1193	0.28	0/1609
21	Y	0.13	0/597	0.32	0/818
22	Z	0.09	0/639	0.23	0/864
23	a	0.22	0/1184	0.36	0/1603
24	c	0.20	0/1346	0.35	0/1840
25	d	0.13	0/1458	0.27	0/1965
26	e	0.30	0/849	0.45	0/1153
27	f	0.14	0/404	0.29	0/547
28	g	0.12	0/1031	0.27	0/1394
29	h	0.12	0/889	0.25	0/1190
30	i	0.16	0/2773	0.32	0/3768
31	j	0.52	0/866	0.67	0/1183

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
32	k	0.40	0/759	0.61	0/1029
33	l	0.24	0/4914	0.43	0/6683
34	m	0.48	0/1356	0.66	0/1839
35	n	0.55	0/491	0.77	0/663
36	o	0.17	0/1070	0.31	0/1451
37	p	0.20	0/1585	0.32	0/2148
38	q	0.15	0/3721	0.31	0/5073
39	r	0.43	0/2507	0.62	1/3428 (0.0%)
40	s	0.11	0/1436	0.28	0/1938
41	t	0.22	0/1038	0.34	0/1389
42	G	0.12	0/5347	0.28	0/7243
43	M	0.12	0/792	0.30	0/1069
44	b	0.20	0/976	0.44	0/1328
All	All	0.24	0/67191	0.38	1/91089 (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
2	C	0	1
7	I	0	1
26	e	0	1
35	n	0	2
37	p	0	2
39	r	0	1
All	All	0	8

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
39	r	91	MET	N-CA-C	-6.05	99.84	109.04

There are no chirality outliers.

5 of 8 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	C	113	ARG	Sidechain
7	I	165	ARG	Sidechain

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Mol	Chain	Res	Type	Group
26	e	88	ARG	Sidechain
35	n	30	ARG	Sidechain
35	n	9	ARG	Sidechain

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	B	3300	0	3258	70	0
2	C	3453	0	3389	75	0
3	D	1737	0	1691	38	0
4	E	1658	0	1662	49	0
5	F	720	0	683	13	0
6	H	1412	0	1363	45	0
7	I	1249	0	1254	30	0
8	J	963	0	962	23	0
9	K	1203	0	1161	24	0
10	L	2625	0	2640	65	0
11	N	911	0	950	13	0
12	O	668	0	672	18	0
12	X	689	0	687	24	0
13	P	669	0	675	14	0
14	Q	955	0	960	25	0
15	R	295	0	279	13	0
16	S	562	0	557	12	0
17	T	638	0	637	20	0
18	U	2573	0	2534	44	0
19	V	1021	0	1025	9	0
20	W	1162	0	1156	31	0
21	Y	571	0	522	17	0
22	Z	620	0	602	9	0
23	a	1151	0	1164	31	0
24	c	1291	0	1185	21	0
25	d	1426	0	1394	37	0
26	e	826	0	788	13	0
27	f	391	0	392	9	0
28	g	1000	0	994	21	0
29	h	867	0	871	17	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
30	i	2710	0	2874	88	0
31	j	845	0	886	42	0
32	k	748	0	799	36	0
33	l	4785	0	4935	150	0
34	m	1321	0	1320	53	0
35	n	479	0	486	9	0
36	o	1041	0	1053	19	0
37	p	1529	0	1465	30	0
38	q	3630	0	3837	91	0
39	r	2437	0	2542	83	0
40	s	1398	0	1374	29	0
41	t	1014	0	983	33	0
42	G	5260	0	5287	106	0
43	M	774	0	801	16	0
44	b	946	0	963	34	0
45	B	31	0	19	0	0
46	B	8	0	0	2	0
46	G	16	0	0	0	0
46	H	16	0	0	1	0
46	I	8	0	0	2	0
47	E	4	0	0	1	0
47	G	4	0	0	0	0
48	L	48	0	26	5	0
49	Q	30	0	30	2	0
All	All	65688	0	65787	1310	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 10.

The worst 5 of 1310 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:B:385:CYS:HB2	46:B:502:SF4:S3	2.08	0.93
6:H:101:HIS:H	6:H:149:MET:HE1	1.42	0.85
10:L:357:LEU:HD13	10:L:357:LEU:H	1.42	0.84
1:B:385:CYS:HB3	46:B:502:SF4:S1	2.19	0.83
23:a:130:GLU:HB2	35:n:58:LYS:HD2	1.59	0.83

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	B	427/464 (92%)	414 (97%)	13 (3%)	0	100	100
2	C	428/469 (91%)	417 (97%)	11 (3%)	0	100	100
3	D	206/264 (78%)	200 (97%)	6 (3%)	0	100	100
4	E	212/249 (85%)	204 (96%)	7 (3%)	1 (0%)	24	59
5	F	91/123 (74%)	90 (99%)	1 (1%)	0	100	100
6	H	174/212 (82%)	172 (99%)	2 (1%)	0	100	100
7	I	154/196 (79%)	149 (97%)	5 (3%)	0	100	100
8	J	116/175 (66%)	114 (98%)	2 (2%)	0	100	100
9	K	142/145 (98%)	141 (99%)	1 (1%)	0	100	100
10	L	324/372 (87%)	312 (96%)	12 (4%)	0	100	100
11	N	110/116 (95%)	109 (99%)	1 (1%)	0	100	100
12	O	81/156 (52%)	80 (99%)	1 (1%)	0	100	100
12	X	83/156 (53%)	83 (100%)	0	0	100	100
13	P	81/99 (82%)	79 (98%)	2 (2%)	0	100	100
14	Q	110/154 (71%)	106 (96%)	3 (3%)	1 (1%)	14	47
15	R	33/110 (30%)	32 (97%)	1 (3%)	0	100	100
16	S	68/70 (97%)	68 (100%)	0	0	100	100
17	T	80/169 (47%)	78 (98%)	2 (2%)	0	100	100
18	U	316/357 (88%)	311 (98%)	4 (1%)	1 (0%)	36	68
19	V	138/141 (98%)	137 (99%)	1 (1%)	0	100	100
20	W	138/144 (96%)	137 (99%)	1 (1%)	0	100	100
21	Y	64/105 (61%)	61 (95%)	3 (5%)	0	100	100
22	Z	75/114 (66%)	75 (100%)	0	0	100	100
23	a	136/189 (72%)	134 (98%)	2 (2%)	0	100	100
24	c	151/186 (81%)	147 (97%)	4 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
25	d	167/176 (95%)	167 (100%)	0	0	100	100
26	e	97/154 (63%)	90 (93%)	6 (6%)	1 (1%)	12	45
27	f	44/76 (58%)	43 (98%)	1 (2%)	0	100	100
28	g	119/122 (98%)	117 (98%)	2 (2%)	0	100	100
29	h	103/106 (97%)	99 (96%)	4 (4%)	0	100	100
30	i	345/347 (99%)	334 (97%)	11 (3%)	0	100	100
31	j	102/115 (89%)	96 (94%)	5 (5%)	1 (1%)	12	45
32	k	96/98 (98%)	94 (98%)	2 (2%)	0	100	100
33	l	601/606 (99%)	574 (96%)	26 (4%)	1 (0%)	43	73
34	m	171/175 (98%)	159 (93%)	12 (7%)	0	100	100
35	n	54/58 (93%)	52 (96%)	2 (4%)	0	100	100
36	o	124/129 (96%)	121 (98%)	3 (2%)	0	100	100
37	p	175/221 (79%)	173 (99%)	2 (1%)	0	100	100
38	q	457/459 (100%)	450 (98%)	7 (2%)	0	100	100
39	r	305/318 (96%)	294 (96%)	10 (3%)	1 (0%)	36	68
40	s	169/249 (68%)	162 (96%)	7 (4%)	0	100	100
41	t	115/137 (84%)	115 (100%)	0	0	100	100
42	G	680/727 (94%)	666 (98%)	14 (2%)	0	100	100
43	M	92/113 (81%)	90 (98%)	2 (2%)	0	100	100
44	b	107/128 (84%)	96 (90%)	11 (10%)	0	100	100
All	All	8061/9449 (85%)	7842 (97%)	212 (3%)	7 (0%)	49	79

5 of 7 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
31	j	34	THR
33	l	562	LEU
39	r	92	PRO
4	E	183	ALA
18	U	59	SER

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	B	343/368 (93%)	339 (99%)	4 (1%)	63	79
2	C	370/398 (93%)	361 (98%)	9 (2%)	43	70
3	D	190/228 (83%)	187 (98%)	3 (2%)	55	75
4	E	183/207 (88%)	173 (94%)	10 (6%)	19	52
5	F	77/97 (79%)	72 (94%)	5 (6%)	15	47
6	H	151/176 (86%)	144 (95%)	7 (5%)	24	58
7	I	132/163 (81%)	128 (97%)	4 (3%)	36	66
8	J	107/152 (70%)	105 (98%)	2 (2%)	50	73
9	K	130/131 (99%)	128 (98%)	2 (2%)	57	76
10	L	283/320 (88%)	262 (93%)	21 (7%)	13	43
11	N	99/101 (98%)	98 (99%)	1 (1%)	68	80
12	O	77/132 (58%)	77 (100%)	0	100	100
12	X	79/132 (60%)	71 (90%)	8 (10%)	7	30
13	P	74/82 (90%)	72 (97%)	2 (3%)	39	68
14	Q	105/134 (78%)	103 (98%)	2 (2%)	50	73
15	R	34/92 (37%)	33 (97%)	1 (3%)	37	67
16	S	58/58 (100%)	53 (91%)	5 (9%)	10	36
17	T	69/134 (52%)	65 (94%)	4 (6%)	18	51
18	U	281/307 (92%)	275 (98%)	6 (2%)	47	71
19	V	101/102 (99%)	99 (98%)	2 (2%)	48	72
20	W	122/124 (98%)	120 (98%)	2 (2%)	55	75
21	Y	60/84 (71%)	58 (97%)	2 (3%)	33	65
22	Z	59/90 (66%)	57 (97%)	2 (3%)	32	64
23	a	121/158 (77%)	120 (99%)	1 (1%)	73	82
24	c	138/160 (86%)	138 (100%)	0	100	100
25	d	153/156 (98%)	143 (94%)	10 (6%)	15	47
26	e	91/129 (70%)	84 (92%)	7 (8%)	12	41
27	f	42/66 (64%)	41 (98%)	1 (2%)	43	70
28	g	108/109 (99%)	101 (94%)	7 (6%)	15	47

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
29	h	93/94 (99%)	90 (97%)	3 (3%)	34	65
30	i	311/311 (100%)	293 (94%)	18 (6%)	18	51
31	j	93/100 (93%)	85 (91%)	8 (9%)	10	36
32	k	85/85 (100%)	79 (93%)	6 (7%)	13	44
33	l	537/540 (99%)	506 (94%)	31 (6%)	18	51
34	m	139/141 (99%)	124 (89%)	15 (11%)	6	27
35	n	53/55 (96%)	50 (94%)	3 (6%)	18	51
36	o	110/114 (96%)	103 (94%)	7 (6%)	16	48
37	p	159/190 (84%)	152 (96%)	7 (4%)	25	59
38	q	409/409 (100%)	393 (96%)	16 (4%)	28	62
39	r	267/275 (97%)	247 (92%)	20 (8%)	12	42
40	s	153/206 (74%)	148 (97%)	5 (3%)	33	65
41	t	108/120 (90%)	101 (94%)	7 (6%)	15	47
42	G	576/610 (94%)	569 (99%)	7 (1%)	63	79
43	M	86/98 (88%)	86 (100%)	0	100	100
44	b	104/121 (86%)	98 (94%)	6 (6%)	18	51
All	All	7120/8059 (88%)	6831 (96%)	289 (4%)	28	60

5 of 289 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
38	q	115	LEU
44	b	86	LEU
38	q	315	LEU
39	r	227	GLU
12	X	124	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 138 such sidechains are listed below:

Mol	Chain	Res	Type
39	r	38	ASN
39	r	212	ASN
42	G	278	HIS
16	S	40	HIS
14	Q	121	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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](#)

11 ligands 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	$\# Z > 2$
46	SF4	I	301	7	0,12,12	-	-	-		
48	NDP	L	401	-	51,52,52	0.49	0	71,80,80	0.80	1 (1%)
45	FMN	B	501	-	33,33,33	1.02	2 (6%)	48,50,50	1.21	8 (16%)
46	SF4	H	302	6	0,12,12	-	-	-		
46	SF4	G	802	42	0,12,12	-	-	-		
49	ZMP	Q	201	-	27,29,36	1.72	4 (14%)	34,38,45	1.67	5 (14%)
47	FES	G	803	42	0,4,4	-	-	-		
46	SF4	H	301	6	0,12,12	-	-	-		
46	SF4	G	801	42	0,12,12	-	-	-		
47	FES	E	301	4	0,4,4	-	-	-		
46	SF4	B	502	1	0,12,12	-	-	-		

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
46	SF4	I	301	7	-	-	0/6/5/5
48	NDP	L	401	-	-	7/34/77/77	0/5/5/5
45	FMN	B	501	-	-	7/18/18/18	0/3/3/3
46	SF4	H	302	6	-	-	0/6/5/5
46	SF4	G	802	42	-	-	0/6/5/5
49	ZMP	Q	201	-	-	9/36/36/43	-
47	FES	G	803	42	-	-	0/1/1/1
46	SF4	H	301	6	-	-	0/6/5/5
46	SF4	G	801	42	-	-	0/6/5/5
47	FES	E	301	4	-	-	0/1/1/1
46	SF4	B	502	1	-	-	0/6/5/5

The worst 5 of 6 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
49	Q	201	ZMP	C13-N1	5.39	1.46	1.33
49	Q	201	ZMP	C16-N2	4.92	1.45	1.33
45	B	501	FMN	C4A-N5	3.30	1.37	1.30
45	B	501	FMN	C10-N1	2.33	1.37	1.33
49	Q	201	ZMP	C10-S1	2.31	1.81	1.76

The worst 5 of 14 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
49	Q	201	ZMP	C9-C10-S1	5.46	119.91	113.40
48	L	401	NDP	P2B-O2B-C2B	-5.39	109.03	123.43
49	Q	201	ZMP	O1-C10-C9	-3.57	119.86	123.98
45	B	501	FMN	C4-N3-C2	-3.29	119.79	125.64
45	B	501	FMN	O4-C4-C4A	-2.64	119.55	126.53

There are no chirality outliers.

5 of 23 torsion outliers are listed below:

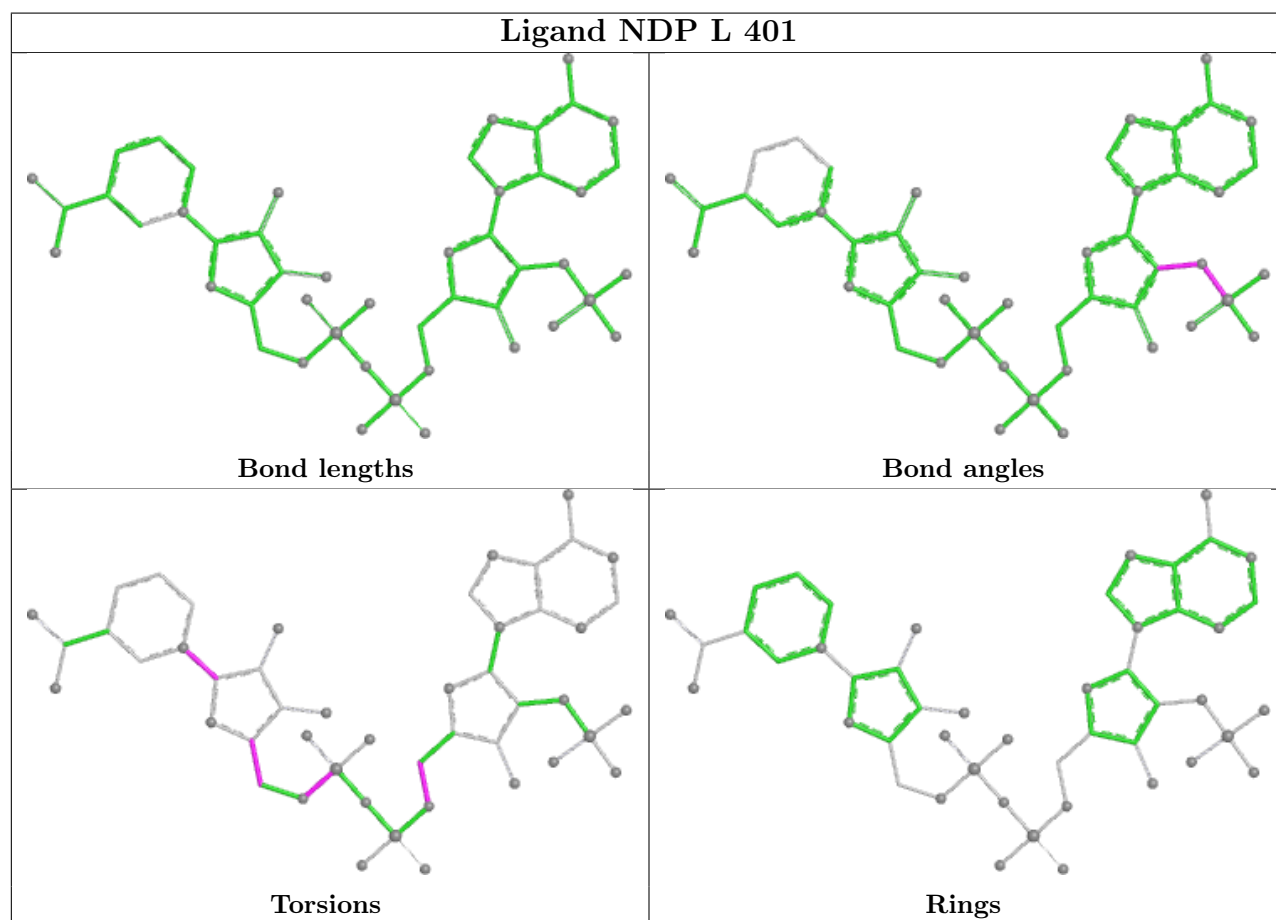
Mol	Chain	Res	Type	Atoms
45	B	501	FMN	N10-C1'-C2'-O2'
45	B	501	FMN	C3'-C4'-C5'-O5'
45	B	501	FMN	O4'-C4'-C5'-O5'
45	B	501	FMN	C5'-O5'-P-O2P
45	B	501	FMN	C5'-O5'-P-O3P

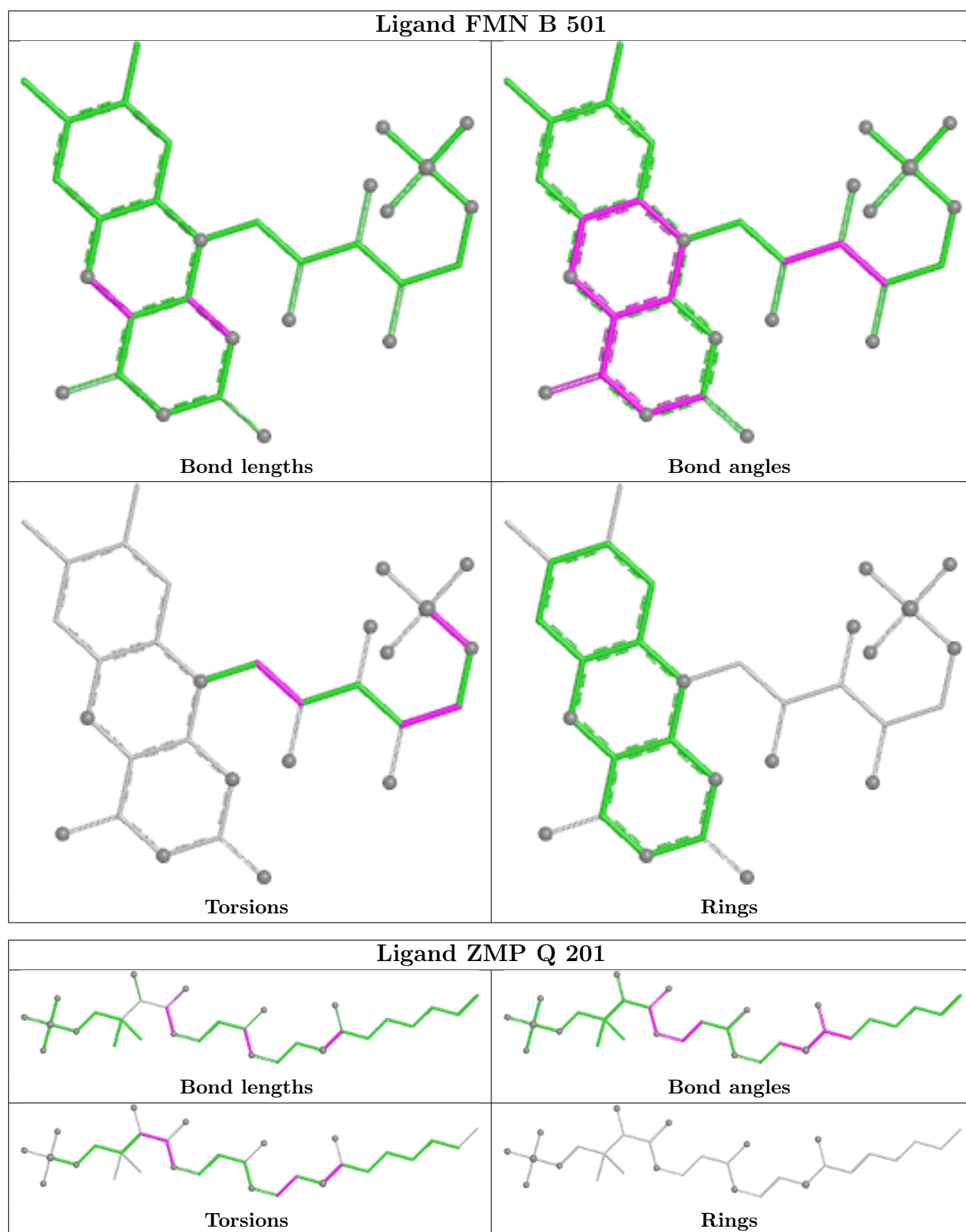
There are no ring outliers.

6 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
46	I	301	SF4	2	0
48	L	401	NDP	5	0
46	H	302	SF4	1	0
49	Q	201	ZMP	2	0
47	E	301	FES	1	0
46	B	502	SF4	2	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 ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

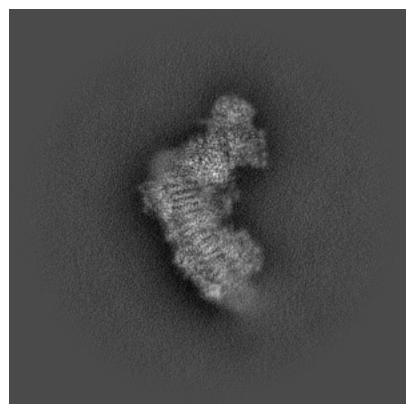
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-45463. These allow visual inspection of the internal detail of the map and identification of artifacts.

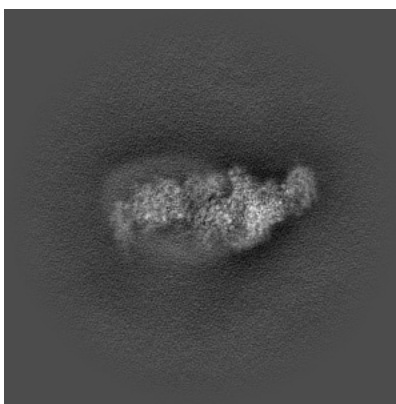
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

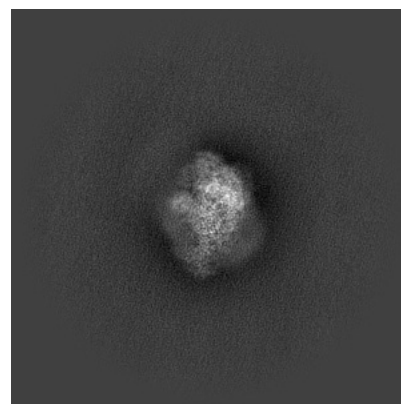
6.1.1 Primary map



X

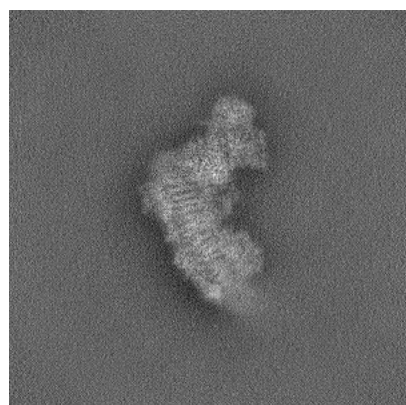


Y

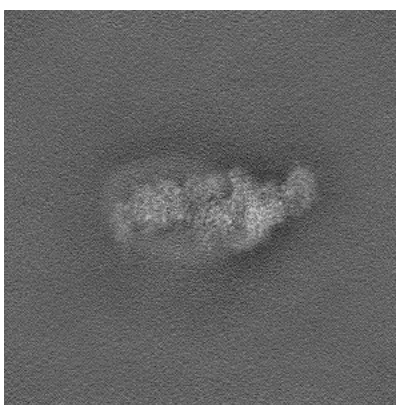


Z

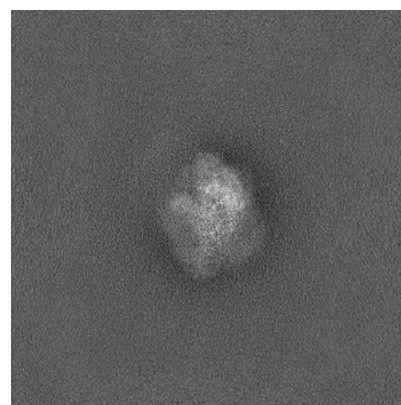
6.1.2 Raw map



X



Y

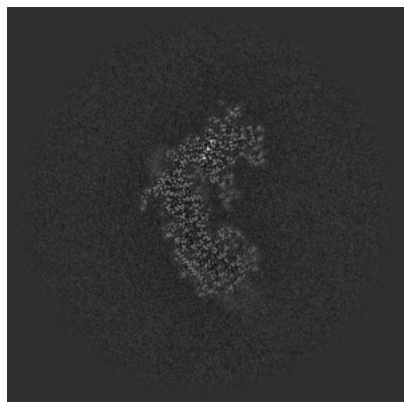


Z

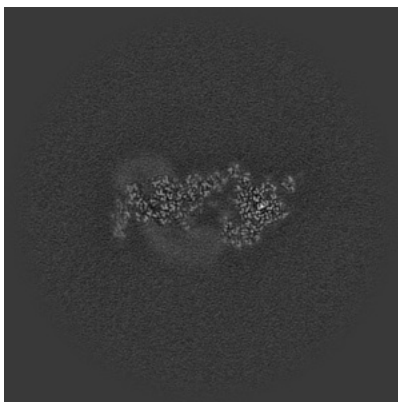
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

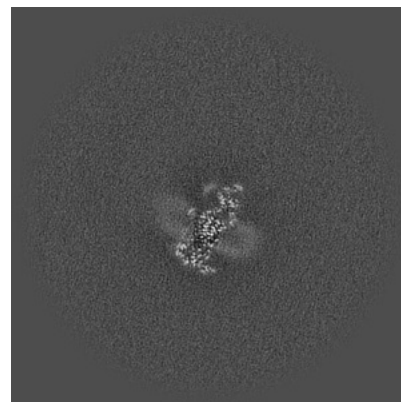
6.2.1 Primary map



X Index: 256

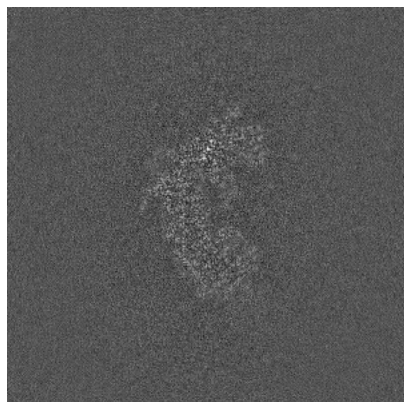


Y Index: 256

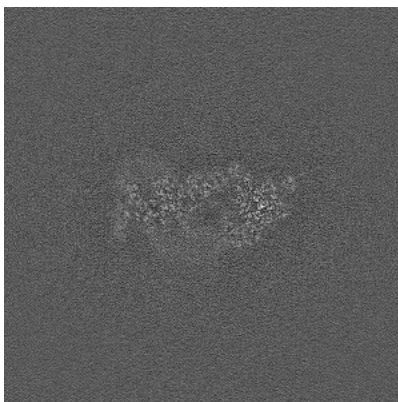


Z Index: 256

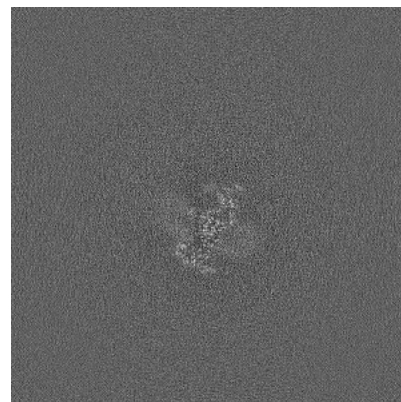
6.2.2 Raw map



X Index: 256



Y Index: 256

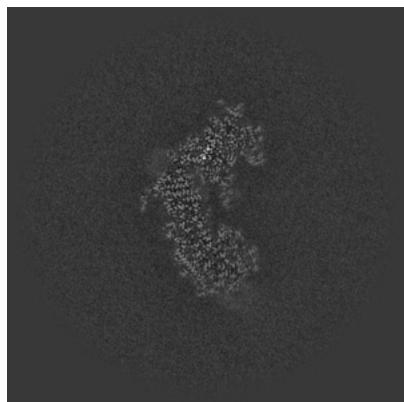


Z Index: 256

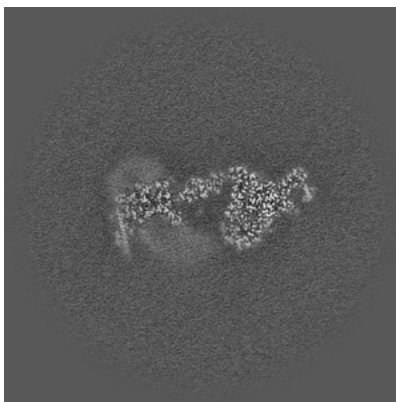
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

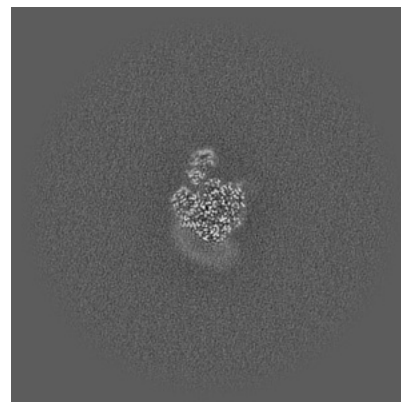
6.3.1 Primary map



X Index: 257

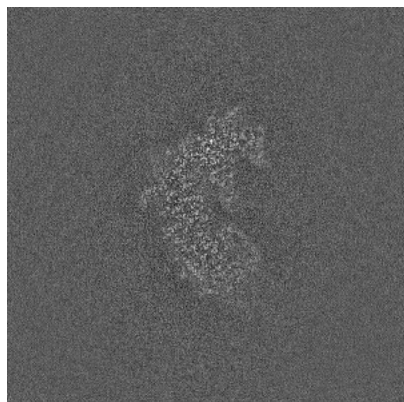


Y Index: 265

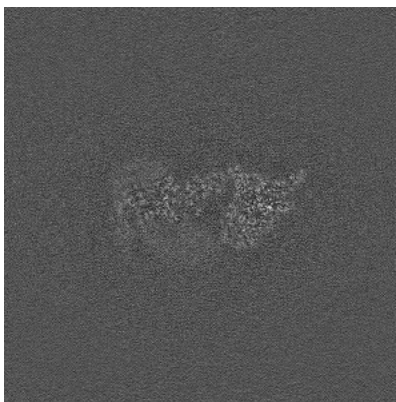


Z Index: 313

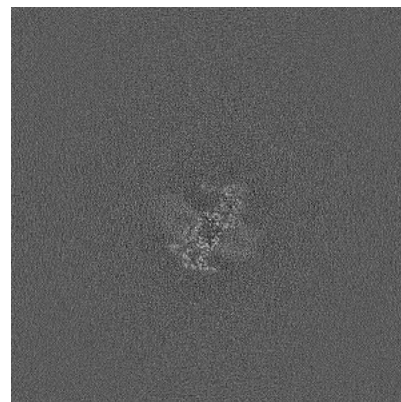
6.3.2 Raw map



X Index: 255



Y Index: 261

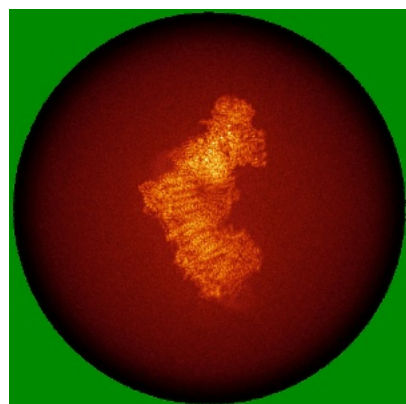


Z Index: 260

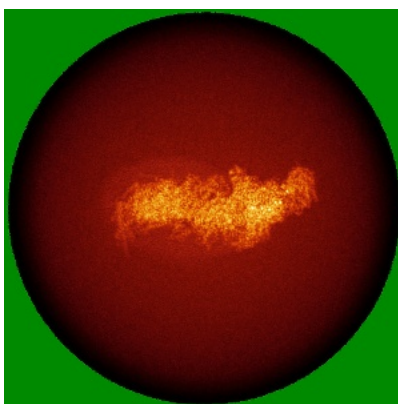
The images above show the largest variance slices of the map in three orthogonal directions.

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

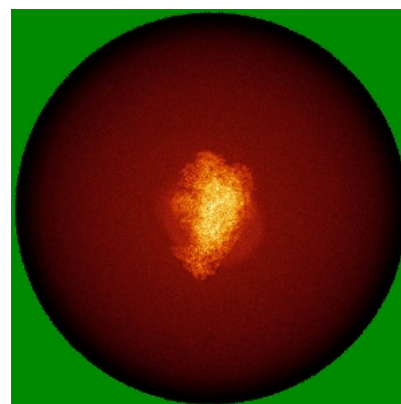
6.4.1 Primary map



X

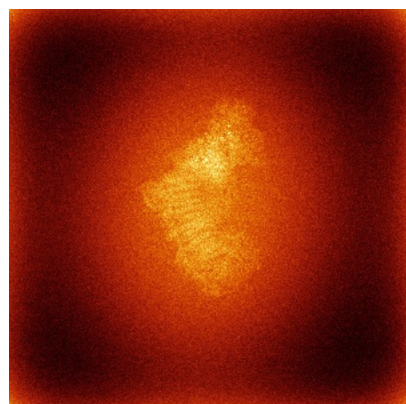


Y

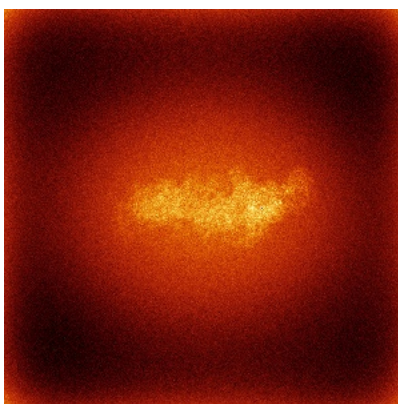


Z

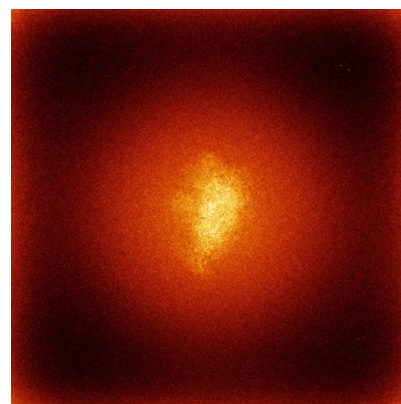
6.4.2 Raw map



X



Y

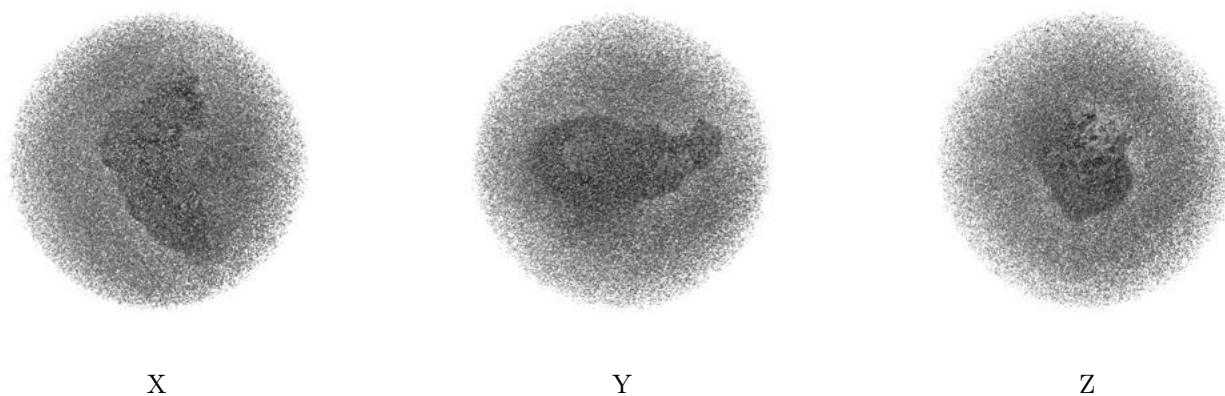


Z

The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

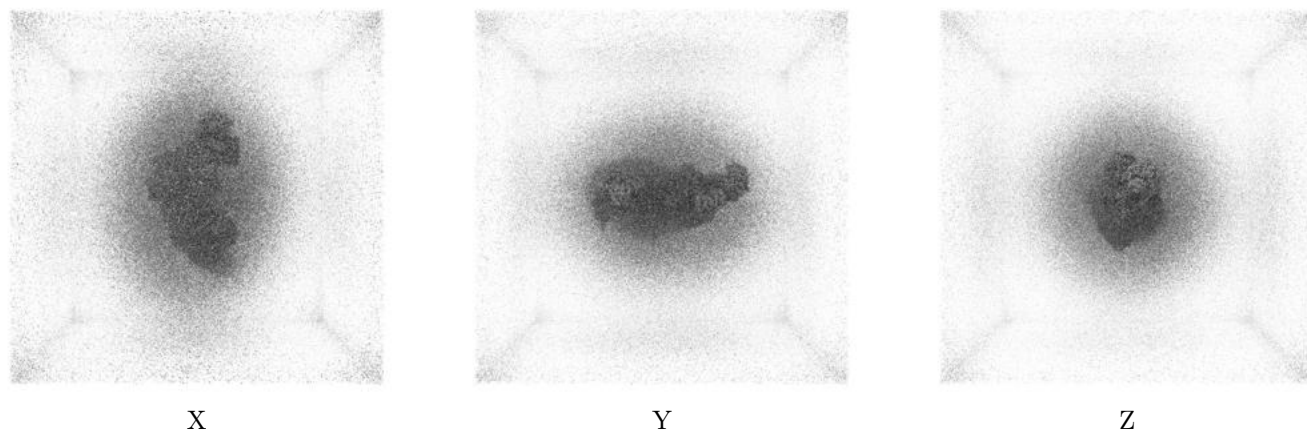
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.1. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



These images show the 3D surface of the raw map. The raw map's contour level was selected so that its surface encloses the same volume as the primary map does at its recommended contour level.

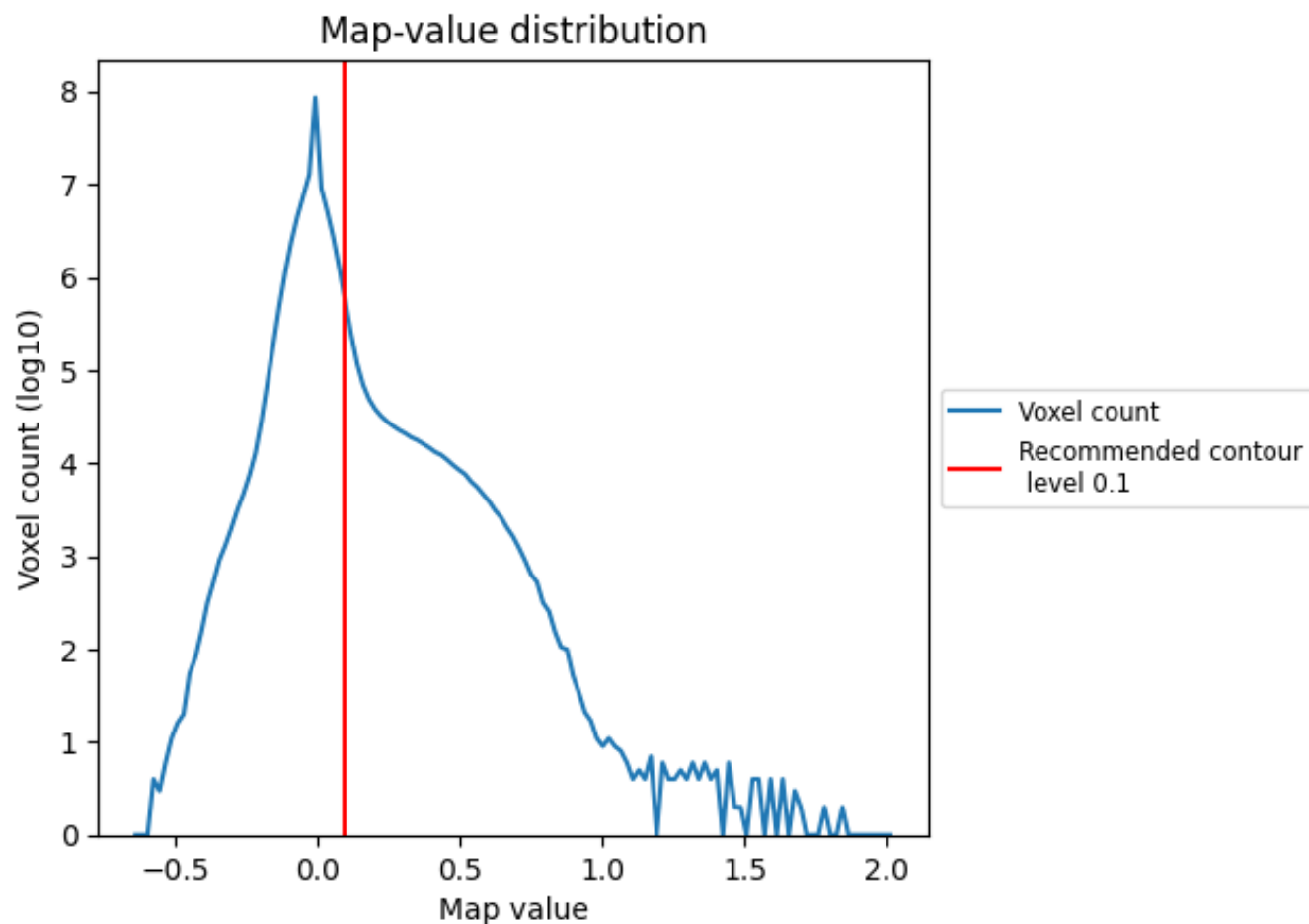
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

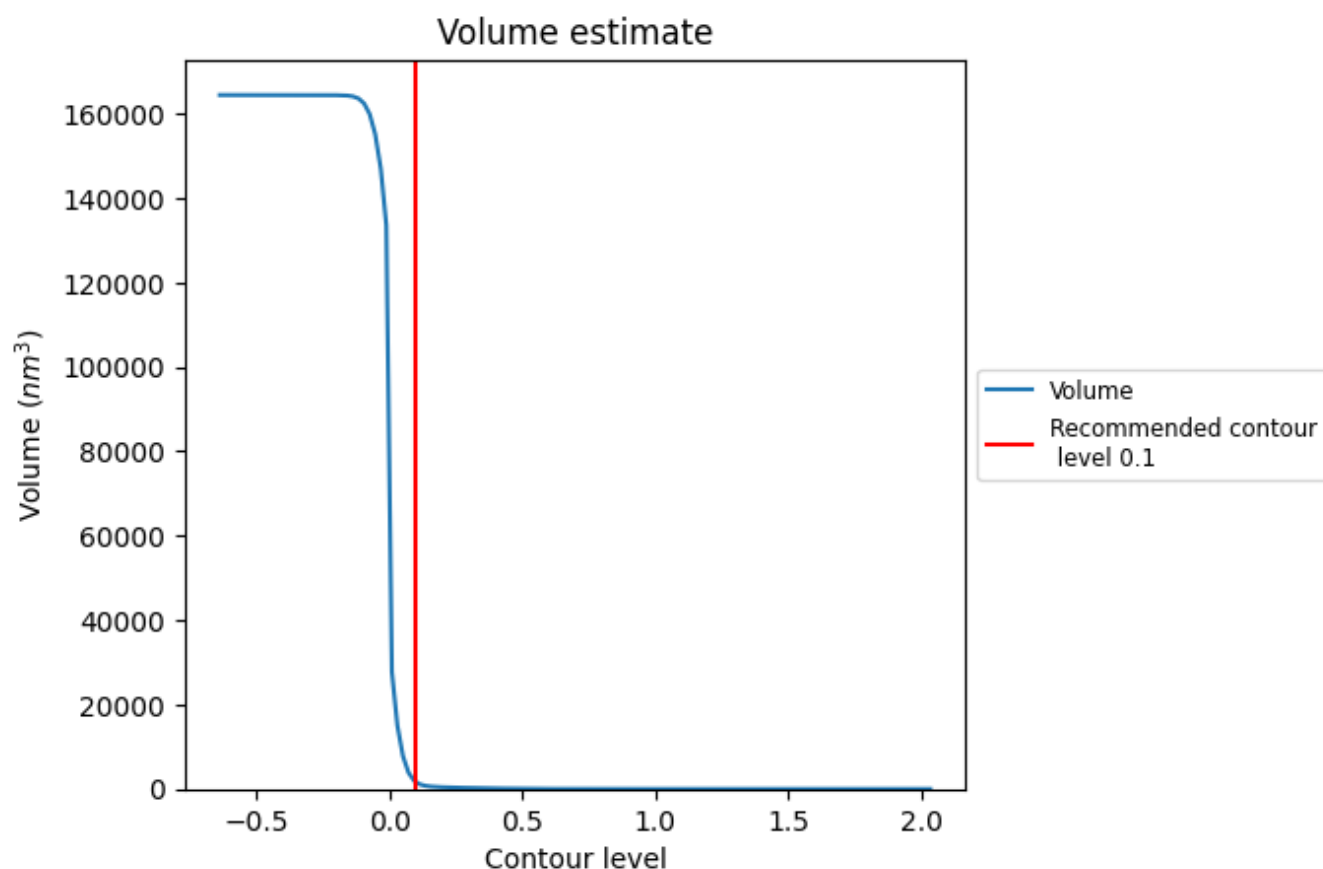
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

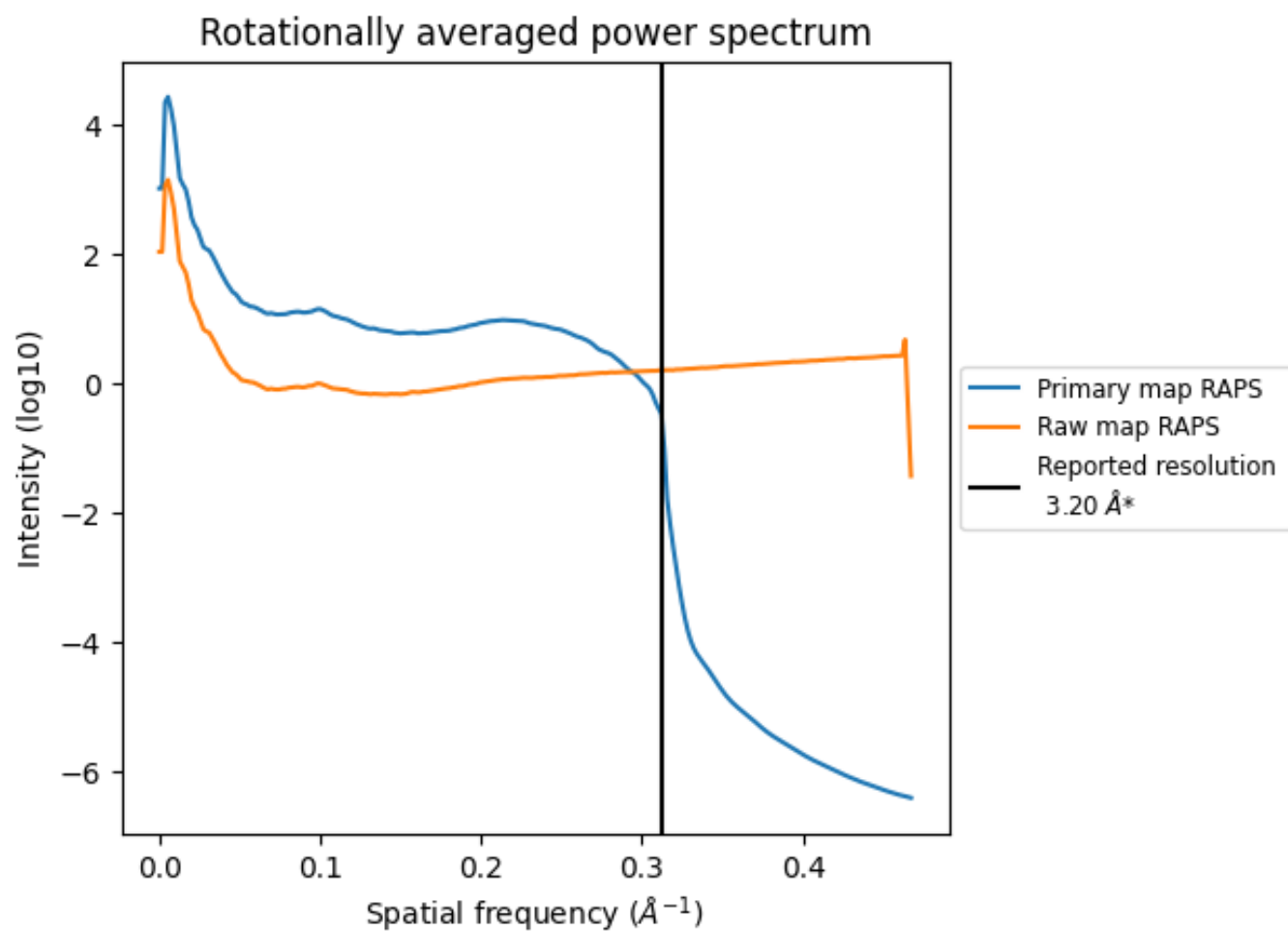
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 1706 nm^3 ; this corresponds to an approximate mass of 1541 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 ⓘ

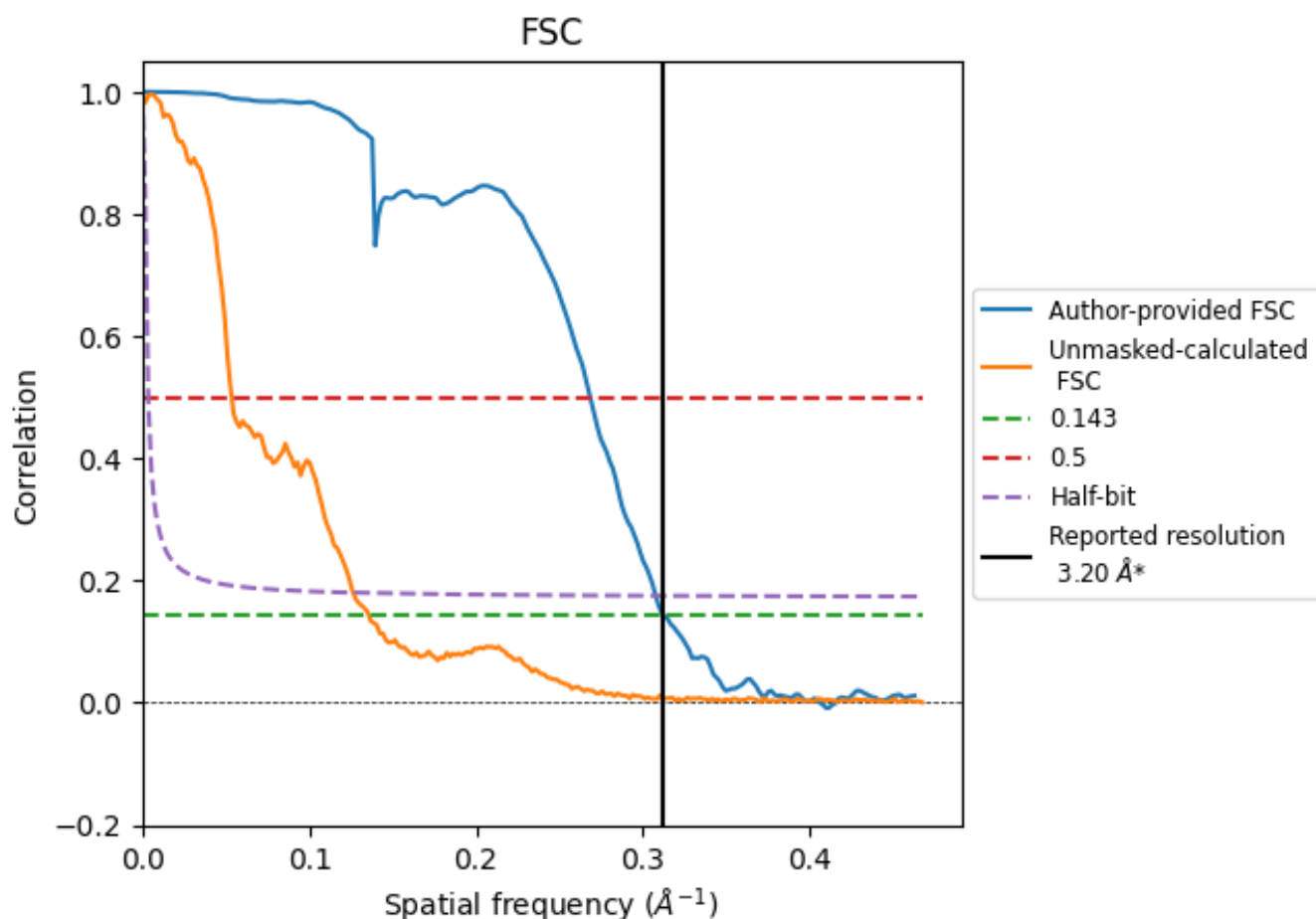


*Reported resolution corresponds to spatial frequency of $0.312 \text{ \AA}^{-1}$

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 $0.312 \text{ \AA}^{-1}$

8.2 Resolution estimates [i](#)

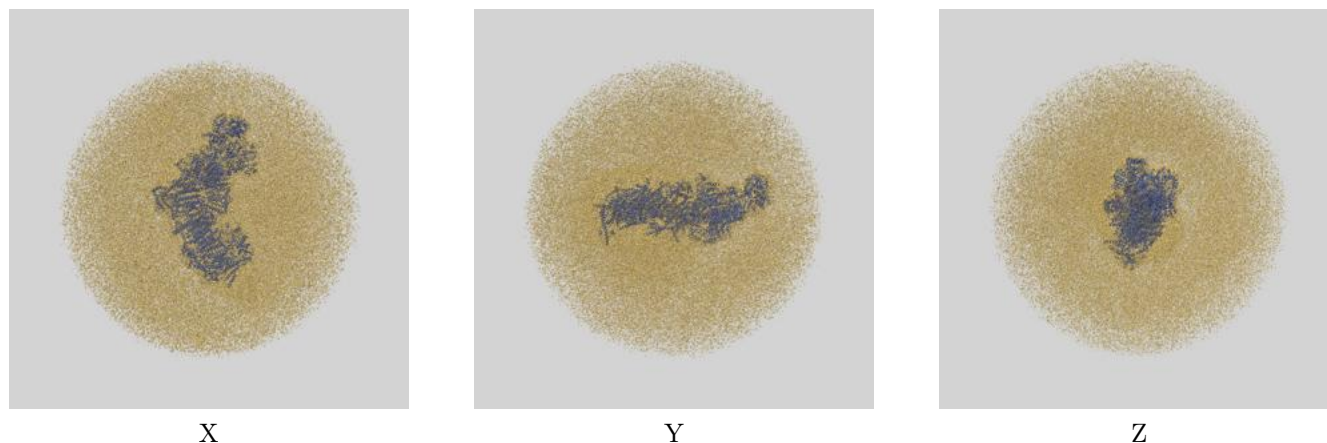
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.20	-	-
Author-provided FSC curve	3.20	3.73	3.25
Unmasked-calculated*	7.37	18.66	7.91

*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 7.37 differs from the reported value 3.2 by more than 10 %

9 Map-model fit [i](#)

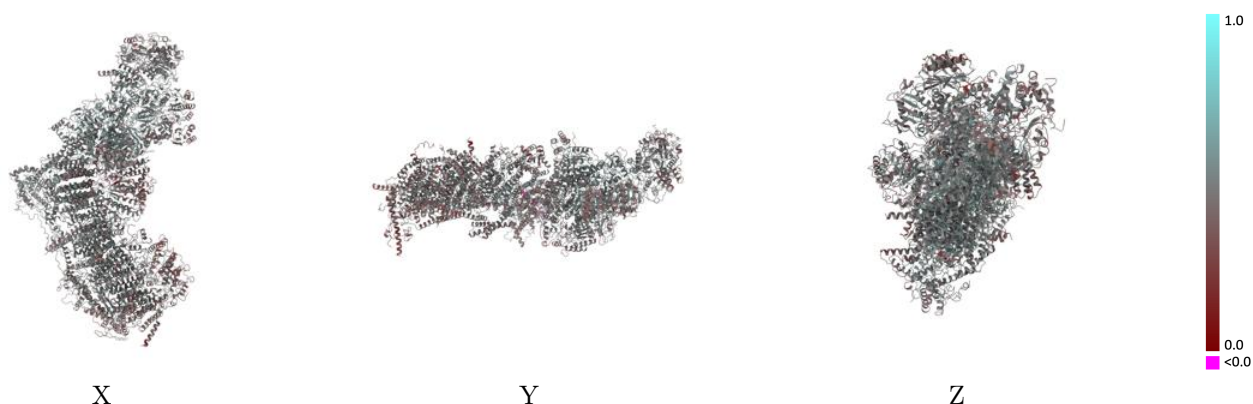
This section contains information regarding the fit between EMDB map EMD-45463 and PDB model 9CD6. Per-residue inclusion information can be found in section [3](#) on page [15](#).

9.1 Map-model overlay [i](#)



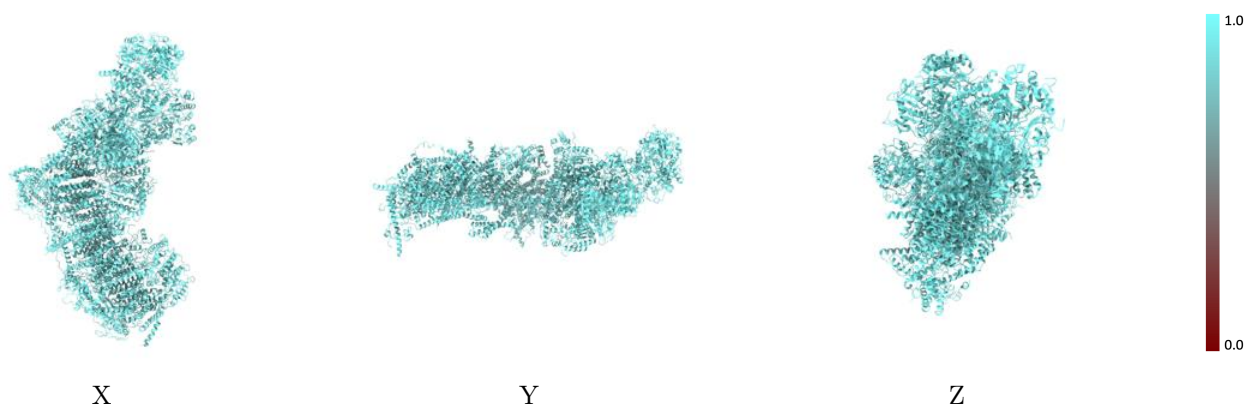
The images above show the 3D surface view of the map at the recommended contour level 0.1 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



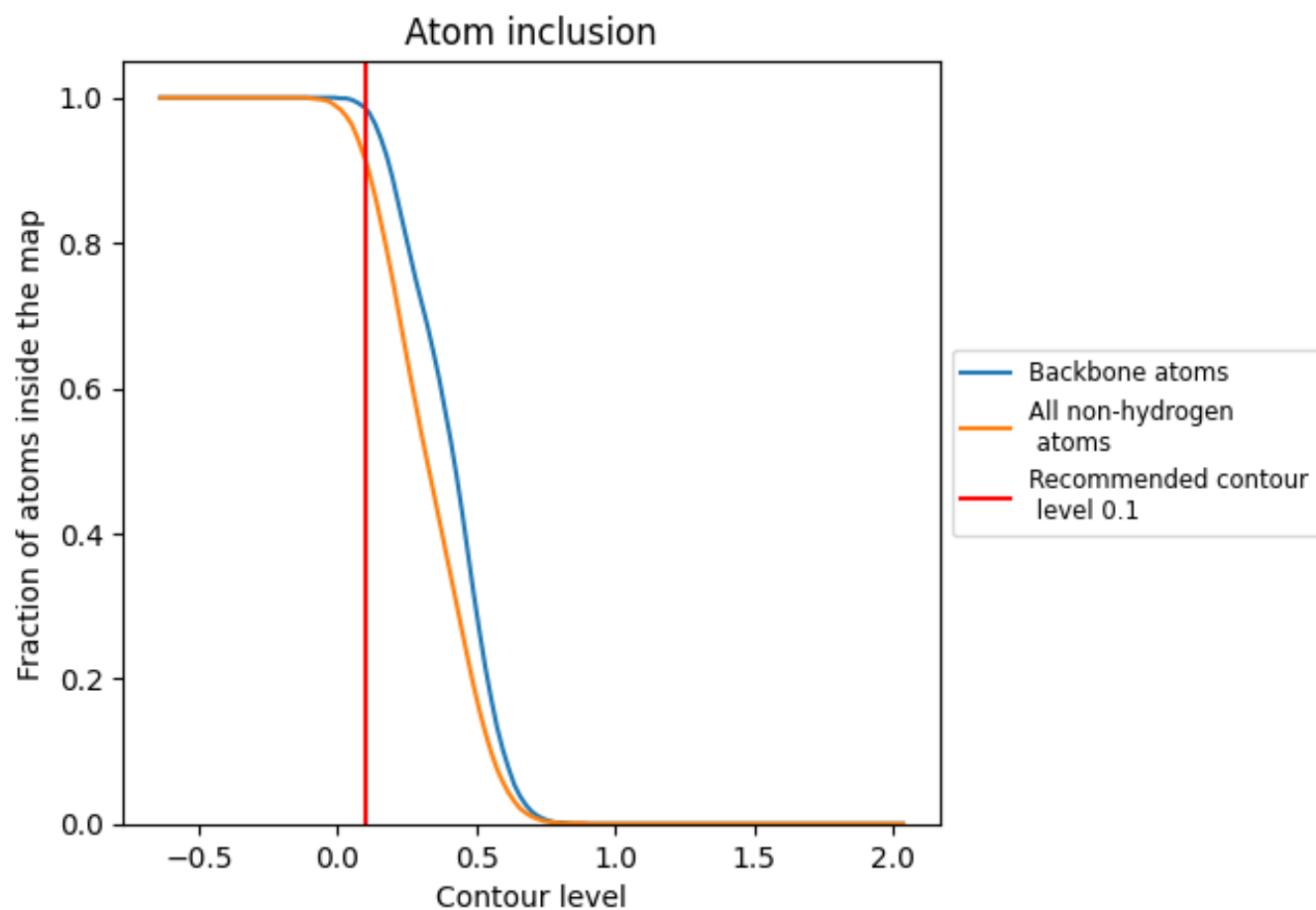
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.1).

9.4 Atom inclusion ⓘ



At the recommended contour level, 99% of all backbone atoms, 92% 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 (0.1) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.9150	 0.4610
B	 0.9450	 0.4660
C	 0.9250	 0.5020
D	 0.9370	 0.5090
E	 0.9330	 0.4420
F	 0.9240	 0.4740
G	 0.9350	 0.4810
H	 0.9510	 0.5110
I	 0.9350	 0.4930
J	 0.9270	 0.4960
K	 0.9350	 0.4890
L	 0.9080	 0.4360
M	 0.9310	 0.4990
N	 0.9120	 0.4360
O	 0.8310	 0.3470
P	 0.9110	 0.4140
Q	 0.8970	 0.4530
R	 0.9370	 0.4320
S	 0.9560	 0.5030
T	 0.9310	 0.4610
U	 0.8900	 0.4300
V	 0.9220	 0.4420
W	 0.9430	 0.4850
X	 0.8680	 0.3850
Y	 0.8780	 0.3720
Z	 0.8790	 0.3700
a	 0.9360	 0.4770
b	 0.8720	 0.3600
c	 0.9020	 0.4440
d	 0.9170	 0.4390
e	 0.8910	 0.4220
f	 0.9000	 0.4230
g	 0.9370	 0.4800
h	 0.9280	 0.4620
i	 0.9240	 0.5010



Continued on next page...

Continued from previous page...

Chain	Atom inclusion	Q-score
j	 0.8680	 0.4440
k	 0.8980	 0.4880
l	 0.8940	 0.4560
m	 0.8800	 0.4300
n	 0.8810	 0.4060
o	 0.8940	 0.4350
p	 0.9080	 0.4170
q	 0.9160	 0.4930
r	 0.9120	 0.4900
s	 0.9250	 0.4760
t	 0.8920	 0.3650