wwPDB EM Validation Summary Report

Feb 13, 2024 – 09:27 AM EST

PDB ID	:	7TMQ
EMDB ID	:	EMD-25999
Title	:	V1 complex lacking subunit C from Saccharomyces cerevisiae, State 3
Authors	:	Vasanthakumar, T.; Keon, K.A.; Bueler, S.A.; Jaskolka, M.C.; Rubinstein, J.L.
Deposited on	:	2022-01-19
Resolution	:	3.30 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:

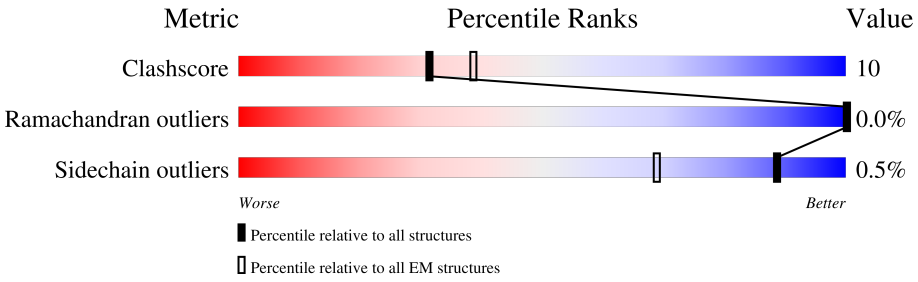
EMDB validation analysis	:	0.0.1.dev70
Mogul	:	1.8.5 (274361), CSD as541be (2020)
MolProbity	:	4.02b-467
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ	:	1.9.9
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 3.30 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)
Clashscore	158937	4297
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826

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

Mol	Chain	Length	Quality of chain
1	A	639	72% 20% 8%
1	C	639	75% 18% 8%
1	E	639	69% 23% 9%
2	B	517	65% 25% 10%
2	D	517	66% 22% 11%
2	F	517	67% 21% 12%
3	G	233	58% 10% 33%
3	I	233	82% 12% 7%

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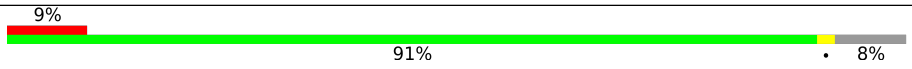
Mol	Chain	Length	Quality of chain
3	K	233	
4	H	114	
4	J	114	
4	L	114	
5	M	256	
6	N	118	
7	P	478	

2 Entry composition (i)

There are 9 unique types of molecules in this entry. The entry contains 31642 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 H(+)-transporting two-sector ATPase.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	591	Total 4241	C 2747	N 743	O 733	S 18	0	0
1	C	590	Total 4043	C 2624	N 711	O 692	S 16	0	0
1	E	583	Total 4117	C 2672	N 719	O 710	S 16	0	0

There are 66 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	617	ASP	-	expression tag	UNP A0A6L0YX77
A	618	TYR	-	expression tag	UNP A0A6L0YX77
A	619	LYS	-	expression tag	UNP A0A6L0YX77
A	620	ASP	-	expression tag	UNP A0A6L0YX77
A	621	HIS	-	expression tag	UNP A0A6L0YX77
A	622	ASP	-	expression tag	UNP A0A6L0YX77
A	623	GLY	-	expression tag	UNP A0A6L0YX77
A	624	ASP	-	expression tag	UNP A0A6L0YX77
A	625	TYR	-	expression tag	UNP A0A6L0YX77
A	626	LYS	-	expression tag	UNP A0A6L0YX77
A	627	ASP	-	expression tag	UNP A0A6L0YX77
A	628	HIS	-	expression tag	UNP A0A6L0YX77
A	629	ASP	-	expression tag	UNP A0A6L0YX77
A	630	ILE	-	expression tag	UNP A0A6L0YX77
A	631	ASP	-	expression tag	UNP A0A6L0YX77
A	632	TYR	-	expression tag	UNP A0A6L0YX77
A	633	LYS	-	expression tag	UNP A0A6L0YX77
A	634	ASP	-	expression tag	UNP A0A6L0YX77
A	635	ASP	-	expression tag	UNP A0A6L0YX77
A	636	ASP	-	expression tag	UNP A0A6L0YX77
A	637	ASP	-	expression tag	UNP A0A6L0YX77
A	638	LYS	-	expression tag	UNP A0A6L0YX77
C	617	ASP	-	expression tag	UNP A0A6L0YX77
C	618	TYR	-	expression tag	UNP A0A6L0YX77

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Chain	Residue	Modelled	Actual	Comment	Reference
C	619	LYS	-	expression tag	UNP A0A6L0YX77
C	620	ASP	-	expression tag	UNP A0A6L0YX77
C	621	HIS	-	expression tag	UNP A0A6L0YX77
C	622	ASP	-	expression tag	UNP A0A6L0YX77
C	623	GLY	-	expression tag	UNP A0A6L0YX77
C	624	ASP	-	expression tag	UNP A0A6L0YX77
C	625	TYR	-	expression tag	UNP A0A6L0YX77
C	626	LYS	-	expression tag	UNP A0A6L0YX77
C	627	ASP	-	expression tag	UNP A0A6L0YX77
C	628	HIS	-	expression tag	UNP A0A6L0YX77
C	629	ASP	-	expression tag	UNP A0A6L0YX77
C	630	ILE	-	expression tag	UNP A0A6L0YX77
C	631	ASP	-	expression tag	UNP A0A6L0YX77
C	632	TYR	-	expression tag	UNP A0A6L0YX77
C	633	LYS	-	expression tag	UNP A0A6L0YX77
C	634	ASP	-	expression tag	UNP A0A6L0YX77
C	635	ASP	-	expression tag	UNP A0A6L0YX77
C	636	ASP	-	expression tag	UNP A0A6L0YX77
C	637	ASP	-	expression tag	UNP A0A6L0YX77
C	638	LYS	-	expression tag	UNP A0A6L0YX77
E	617	ASP	-	expression tag	UNP A0A6L0YX77
E	618	TYR	-	expression tag	UNP A0A6L0YX77
E	619	LYS	-	expression tag	UNP A0A6L0YX77
E	620	ASP	-	expression tag	UNP A0A6L0YX77
E	621	HIS	-	expression tag	UNP A0A6L0YX77
E	622	ASP	-	expression tag	UNP A0A6L0YX77
E	623	GLY	-	expression tag	UNP A0A6L0YX77
E	624	ASP	-	expression tag	UNP A0A6L0YX77
E	625	TYR	-	expression tag	UNP A0A6L0YX77
E	626	LYS	-	expression tag	UNP A0A6L0YX77
E	627	ASP	-	expression tag	UNP A0A6L0YX77
E	628	HIS	-	expression tag	UNP A0A6L0YX77
E	629	ASP	-	expression tag	UNP A0A6L0YX77
E	630	ILE	-	expression tag	UNP A0A6L0YX77
E	631	ASP	-	expression tag	UNP A0A6L0YX77
E	632	TYR	-	expression tag	UNP A0A6L0YX77
E	633	LYS	-	expression tag	UNP A0A6L0YX77
E	634	ASP	-	expression tag	UNP A0A6L0YX77
E	635	ASP	-	expression tag	UNP A0A6L0YX77
E	636	ASP	-	expression tag	UNP A0A6L0YX77
E	637	ASP	-	expression tag	UNP A0A6L0YX77
E	638	LYS	-	expression tag	UNP A0A6L0YX77

- Molecule 2 is a protein called Vacuolar proton pump subunit B.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	464	Total	C	N	O	S	0	0
			3424	2204	602	606	12		
2	D	458	Total	C	N	O	S	0	0
			3320	2154	591	564	11		
2	F	457	Total	C	N	O	S	0	0
			3300	2139	588	563	10		

- Molecule 3 is a protein called V-type proton ATPase subunit E.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	G	157	Total	C	N	O	S	0	0
			1063	692	192	177	2		
3	I	217	Total	C	N	O	S	0	0
			1360	869	250	239	2		
3	K	206	Total	C	N	O	S	0	0
			1322	850	244	226	2		

- Molecule 4 is a protein called V-type proton ATPase subunit G.

Mol	Chain	Residues	Atoms				AltConf	Trace
4	H	51	Total	C	N	O	0	0
			299	190	56	53		
4	J	108	Total	C	N	O	0	0
			581	358	112	111		
4	L	88	Total	C	N	O	0	0
			486	302	93	91		

- Molecule 5 is a protein called V-type proton ATPase subunit D.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	M	214	Total	C	N	O	S	0	0
			1286	800	251	234	1		

- Molecule 6 is a protein called V-type proton ATPase subunit F.

Mol	Chain	Residues	Atoms				AltConf	Trace
6	N	111	Total	C	N	O	0	0
			561	339	111	111		

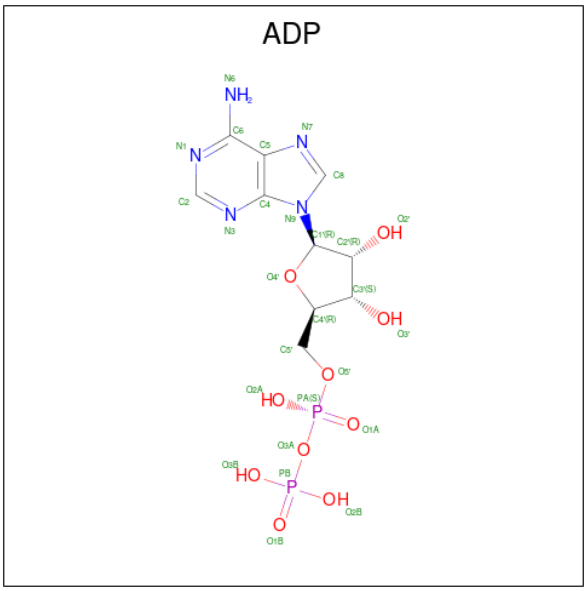
- Molecule 7 is a protein called V-type proton ATPase subunit H.

Mol	Chain	Residues	Atoms				AltConf	Trace
7	P	441	Total	C	N	O	0	0
			2211	1329	441	441		

- Molecule 8 is MAGNESIUM ION (three-letter code: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		AltConf
8	A	1	Total	Mg	0
			1	1	

- Molecule 9 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$) (labeled as "Ligand of Interest" by depositor).



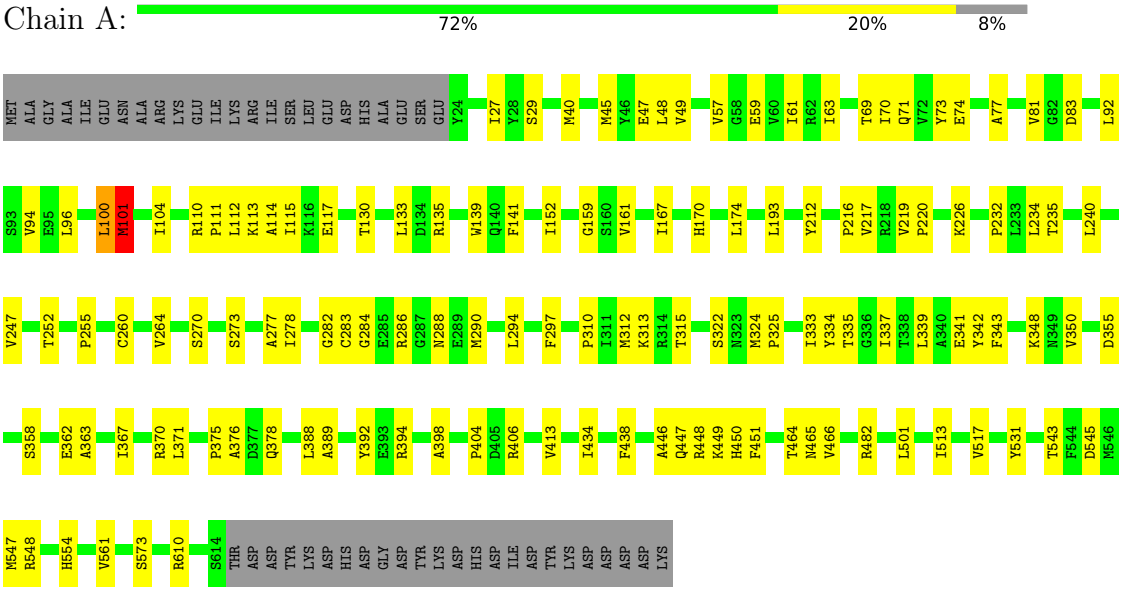
Mol	Chain	Residues	Atoms					AltConf
9	A	1	Total	C	N	O	P	0
			27	10	5	10	2	

3 Residue-property plots (i)

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

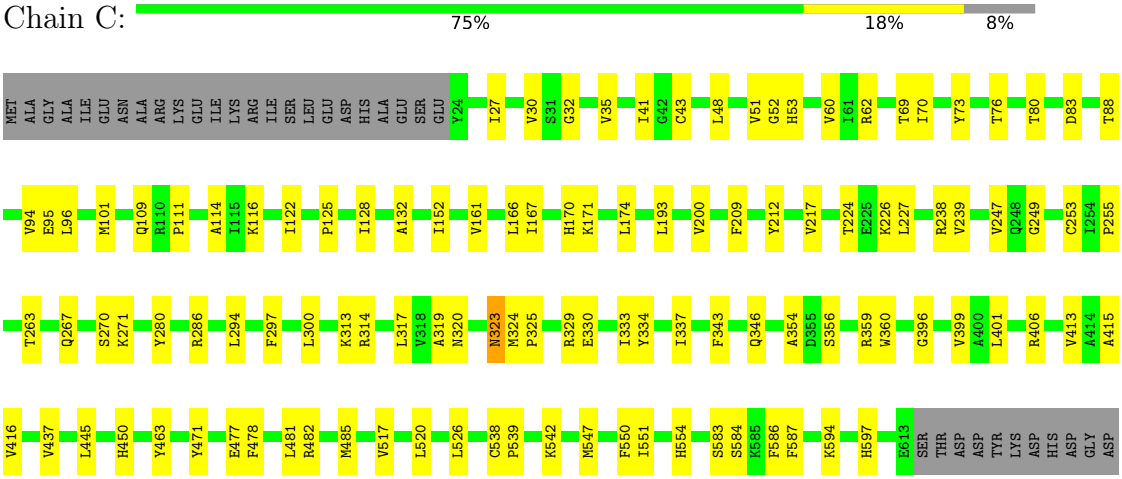
• Molecule 1: H(+)-transporting two-sector ATPase

Chain A:



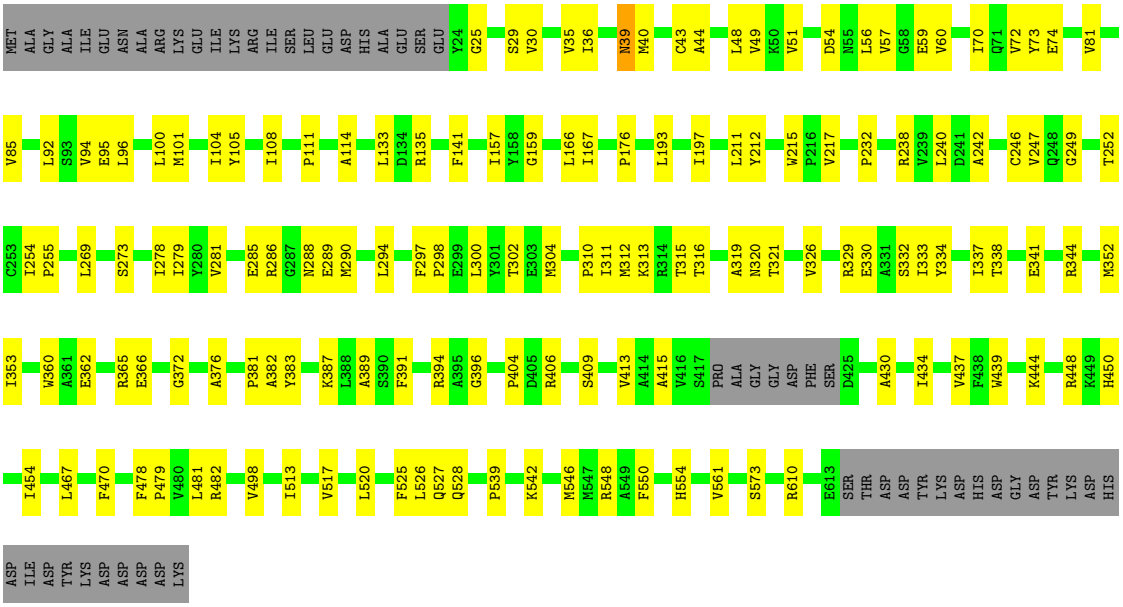
• Molecule 1: H(+)-transporting two-sector ATPase

Chain C:



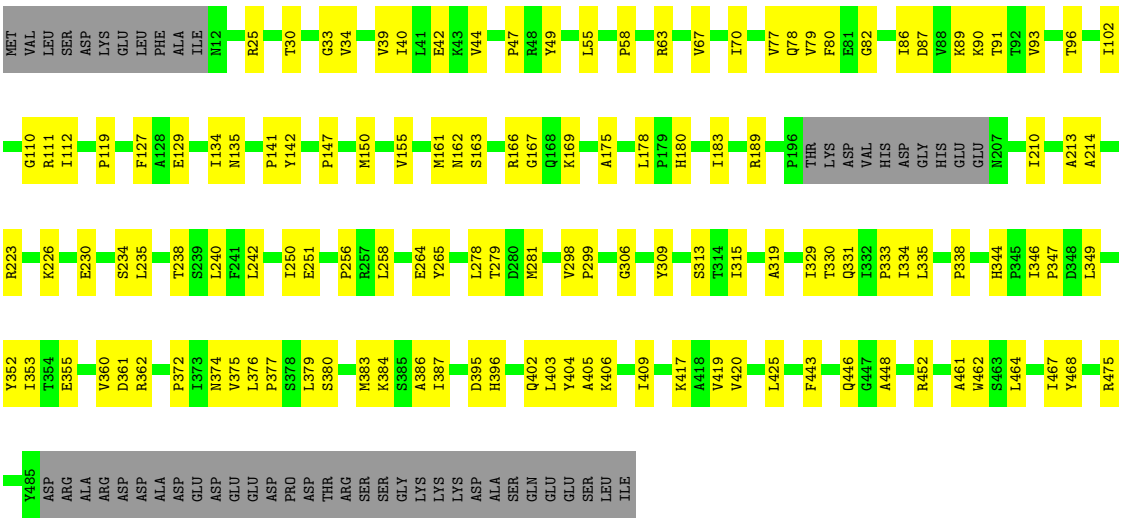
• Molecule 1: H(+)-transporting two-sector ATPase

Chain E:



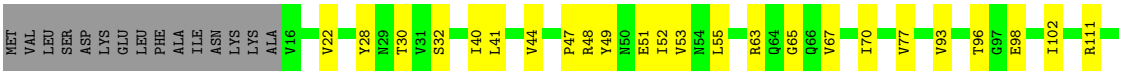
• Molecule 2: Vacuolar proton pump subunit B

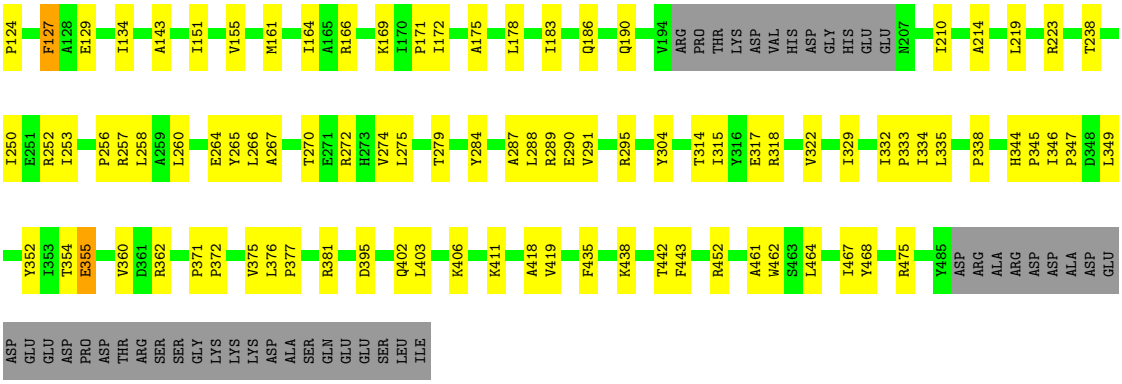
Chain B:



• Molecule 2: Vacuolar proton pump subunit B

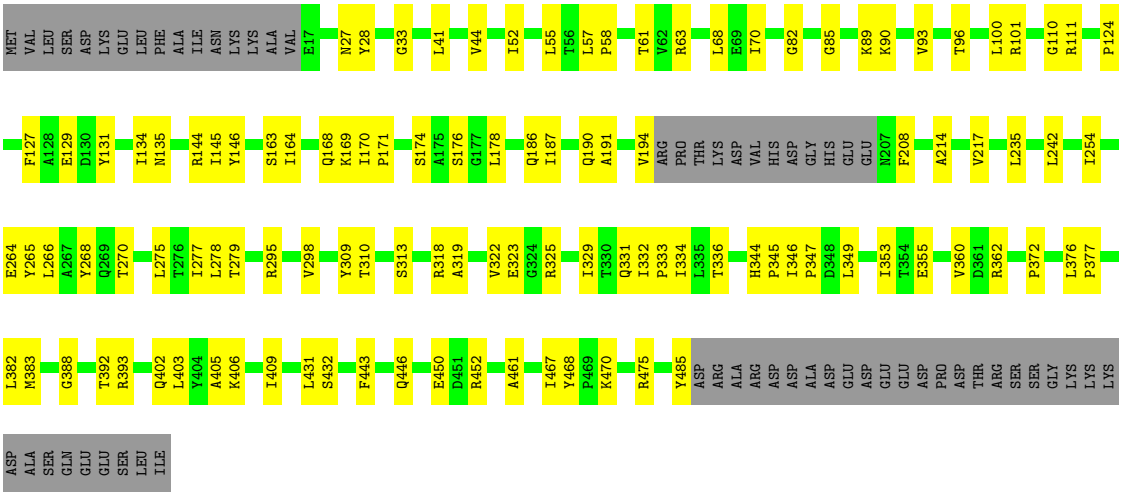
Chain D:





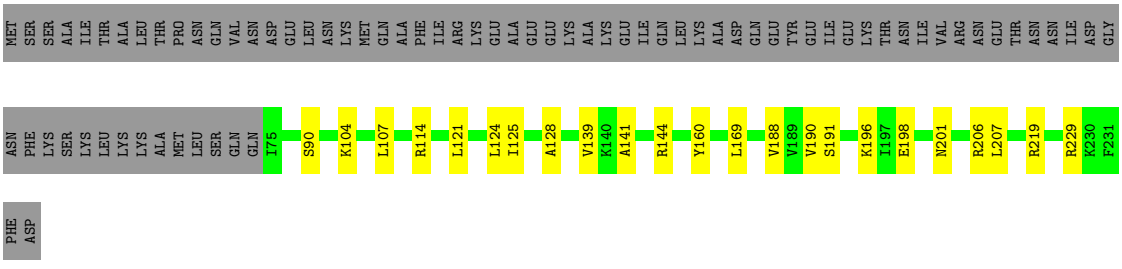
• Molecule 2: Vacuolar proton pump subunit B

Chain F:



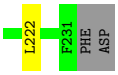
• Molecule 3: V-type proton ATPase subunit E

Chain G:



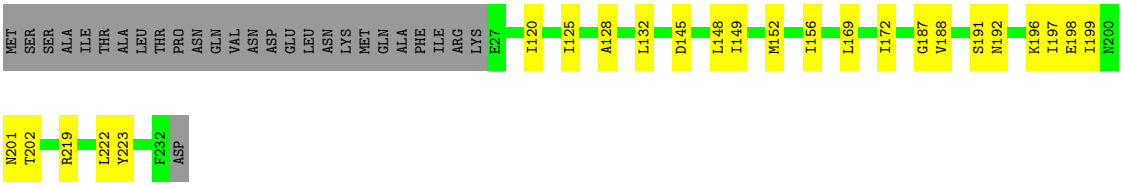
• Molecule 3: V-type proton ATPase subunit E

Chain I:



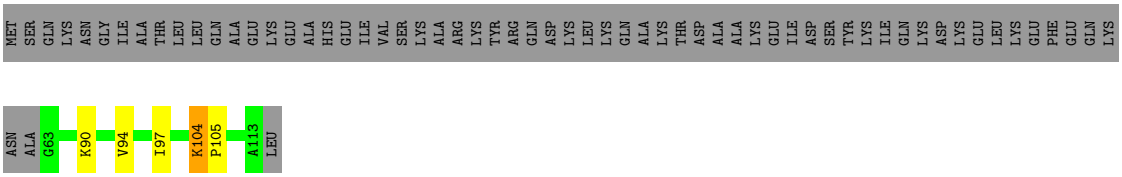
• Molecule 3: V-type proton ATPase subunit E

Chain K:



• Molecule 4: V-type proton ATPase subunit G

Chain H:

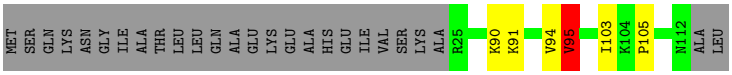


• Molecule 4: V-type proton ATPase subunit G

Chain J:

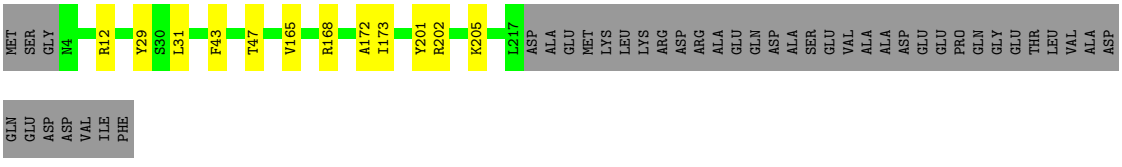
• Molecule 4: V-type proton ATPase subunit G

Chain L:



• Molecule 5: V-type proton ATPase subunit D

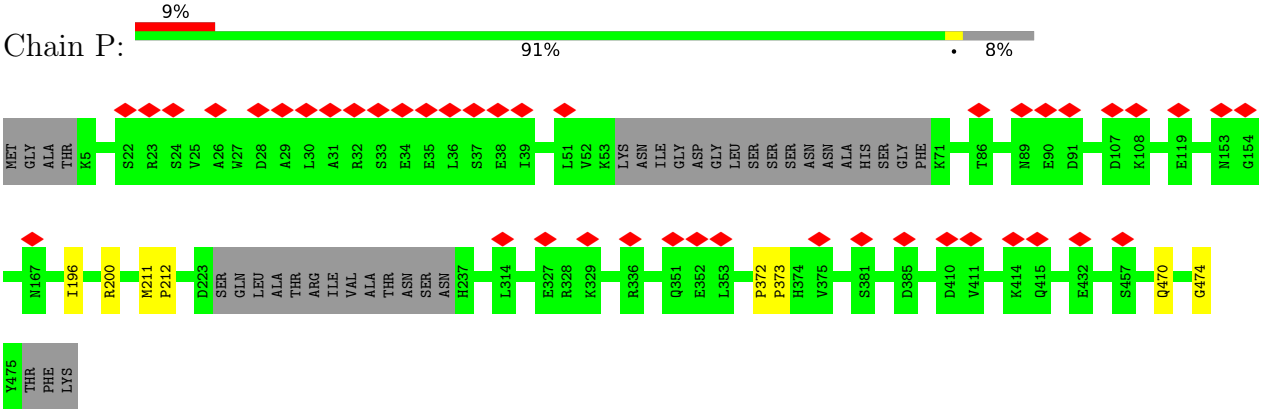
Chain M:



• Molecule 6: V-type proton ATPase subunit F

Chain N:

• Molecule 7: V-type proton ATPase subunit H



4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	156541	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/Å^2$)	41	Depositor
Minimum defocus (nm)	500	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	Not provided	
Image detector	FEI FALCON IV (4k x 4k)	Depositor
Maximum map value	8.316	Depositor
Minimum map value	-5.249	Depositor
Average map value	0.005	Depositor
Map value standard deviation	0.147	Depositor
Recommended contour level	0.3	Depositor
Map size (Å)	412.0, 412.0, 412.0	wwPDB
Map dimensions	400, 400, 400	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.03, 1.03, 1.03	Depositor

5 Model quality

5.1 Standard geometry

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

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	A	0.31	0/4339	0.50	0/5922
1	C	0.28	0/4136	0.47	0/5670
1	E	0.27	0/4212	0.48	0/5762
2	B	0.35	0/3493	0.53	0/4758
2	D	0.35	0/3388	0.53	0/4624
2	F	0.32	0/3368	0.52	0/4601
3	G	0.24	0/1073	0.47	0/1464
3	I	0.24	0/1370	0.44	0/1880
3	K	0.24	0/1333	0.44	0/1825
4	H	0.23	0/300	0.40	0/410
4	J	0.23	0/582	0.33	0/806
4	L	0.25	0/487	0.43	0/672
5	M	0.24	0/1295	0.44	0/1778
6	N	0.23	0/565	0.41	0/791
7	P	0.22	0/2217	0.34	0/3099
All	All	0.29	0/32158	0.48	0/44062

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	A	4241	0	3998	98	0
1	C	4043	0	3652	75	0
1	E	4117	0	3805	100	0
2	B	3424	0	3326	100	0
2	D	3320	0	3190	89	0
2	F	3300	0	3131	74	0
3	G	1063	0	1017	18	0
3	I	1360	0	1152	21	0
3	K	1322	0	1150	16	0
4	H	299	0	259	4	0
4	J	581	0	383	4	0
4	L	486	0	341	6	0
5	M	1286	0	970	10	0
6	N	561	0	277	2	0
7	P	2211	0	993	4	0
8	A	1	0	0	0	0
9	A	27	0	12	2	0
All	All	31642	0	27656	567	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 567 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:E:39:ASN:O	1:E:39:ASN:ND2	2.11	0.83
1:A:517:VAL:HG21	1:A:554:HIS:HB2	1.64	0.79
2:D:266:LEU:O	2:D:270:THR:HB	1.84	0.77
3:G:114:ARG:HH22	3:G:144:ARG:HG2	1.51	0.76
2:B:175:ALA:H	2:B:178:LEU:HD12	1.51	0.75

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	589/639 (92%)	569 (97%)	19 (3%)	1 (0%)	47 77
1	C	588/639 (92%)	569 (97%)	19 (3%)	0	100 100
1	E	579/639 (91%)	562 (97%)	17 (3%)	0	100 100
2	B	460/517 (89%)	445 (97%)	15 (3%)	0	100 100
2	D	454/517 (88%)	434 (96%)	20 (4%)	0	100 100
2	F	453/517 (88%)	444 (98%)	9 (2%)	0	100 100
3	G	155/233 (66%)	152 (98%)	3 (2%)	0	100 100
3	I	215/233 (92%)	209 (97%)	6 (3%)	0	100 100
3	K	204/233 (88%)	202 (99%)	2 (1%)	0	100 100
4	H	49/114 (43%)	48 (98%)	1 (2%)	0	100 100
4	J	106/114 (93%)	106 (100%)	0	0	100 100
4	L	86/114 (75%)	83 (96%)	2 (2%)	1 (1%)	13 42
5	M	212/256 (83%)	207 (98%)	5 (2%)	0	100 100
6	N	109/118 (92%)	107 (98%)	2 (2%)	0	100 100
7	P	435/478 (91%)	430 (99%)	5 (1%)	0	100 100
All	All	4694/5361 (88%)	4567 (97%)	125 (3%)	2 (0%)	100 100

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	101	MET
4	L	95	VAL

5.3.2 Protein sidechains (i)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	374/537 (70%)	372 (100%)	2 (0%)	88 93

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	C	320/537 (60%)	318 (99%)	2 (1%)	86 91
1	E	349/537 (65%)	347 (99%)	2 (1%)	86 91
2	B	327/444 (74%)	325 (99%)	2 (1%)	86 91
2	D	297/444 (67%)	295 (99%)	2 (1%)	84 90
2	F	290/444 (65%)	290 (100%)	0	100 100
3	G	83/208 (40%)	83 (100%)	0	100 100
3	I	85/208 (41%)	85 (100%)	0	100 100
3	K	86/208 (41%)	86 (100%)	0	100 100
4	H	17/94 (18%)	16 (94%)	1 (6%)	19 49
4	J	16/94 (17%)	16 (100%)	0	100 100
4	L	17/94 (18%)	16 (94%)	1 (6%)	19 49
5	M	63/221 (28%)	63 (100%)	0	100 100
6	N	5/104 (5%)	5 (100%)	0	100 100
7	P	9/439 (2%)	9 (100%)	0	100 100
All	All	2338/4613 (51%)	2326 (100%)	12 (0%)	89 93

5 of 12 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	D	355	GLU
1	E	39	ASN
4	L	95	VAL
1	E	406	ARG
2	B	330	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 7 such sidechains are listed below:

Mol	Chain	Res	Type
1	E	320	ASN
2	F	135	ASN
2	F	344	HIS
2	F	331	GLN
2	D	344	HIS

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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 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	$\#\|Z\| > 2$
9	ADP	A	702	8	24,29,29	0.95	1 (4%)	29,45,45	1.41	4 (13%)

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
9	ADP	A	702	8	-	3/12/32/32	0/3/3/3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	A	702	ADP	C5-C4	2.32	1.47	1.40

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
9	A	702	ADP	PA-O3A-PB	-3.54	120.67	132.83
9	A	702	ADP	C3'-C2'-C1'	3.19	105.78	100.98
9	A	702	ADP	N3-C2-N1	-2.99	124.01	128.68
9	A	702	ADP	C4-C5-N7	-2.76	106.52	109.40

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
9	A	702	ADP	C5'-O5'-PA-O1A
9	A	702	ADP	C5'-O5'-PA-O3A
9	A	702	ADP	C5'-O5'-PA-O2A

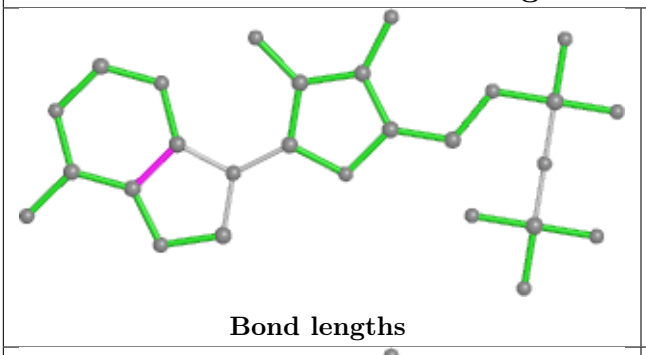
There are no ring outliers.

1 monomer is involved in 2 short contacts:

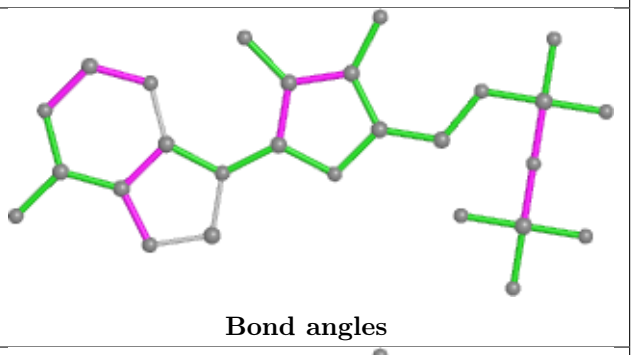
Mol	Chain	Res	Type	Clashes	Symm-Clashes
9	A	702	ADP	2	0

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 then 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.

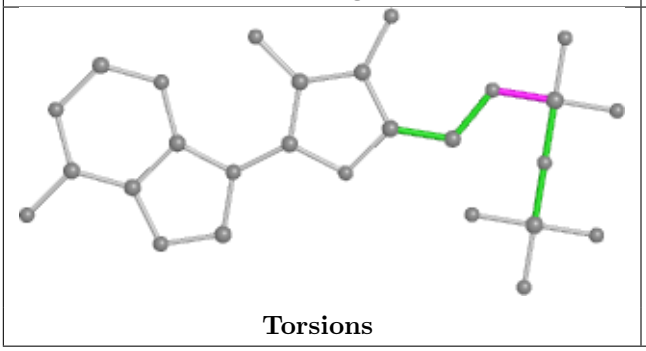
Ligand ADP A 702



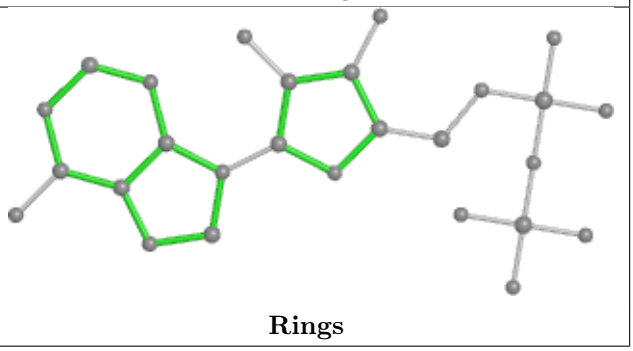
Bond lengths



Bond angles



Torsions



Rings

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

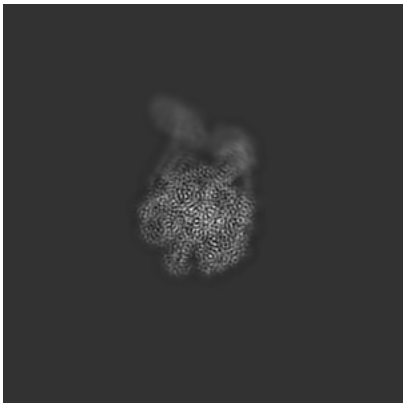
6 Map visualisation

This section contains visualisations of the EMDB entry EMD-25999. These allow visual inspection of the internal detail of the map and identification of artifacts.

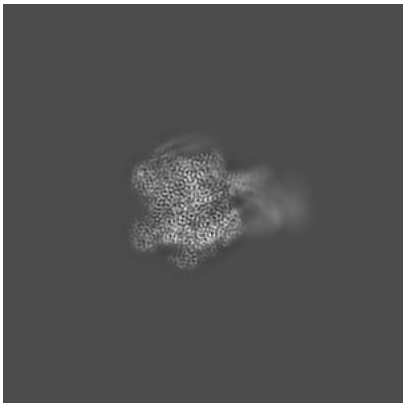
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections

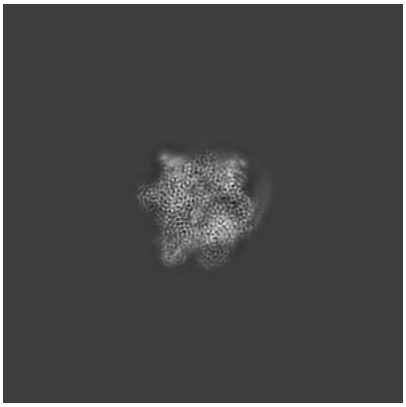
6.1.1 Primary map



X



Y



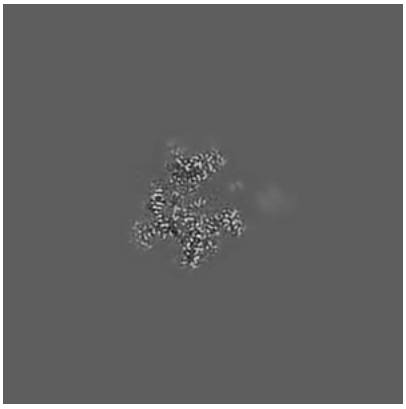
Z

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

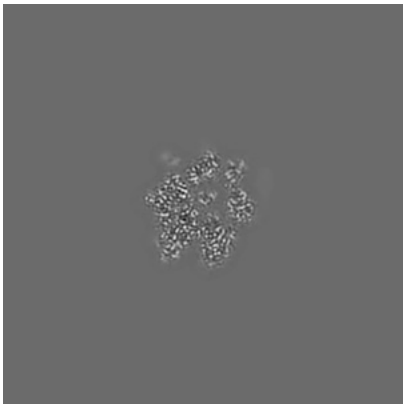
6.2 Central slices

6.2.1 Primary map

X Index: 200



Y Index: 200

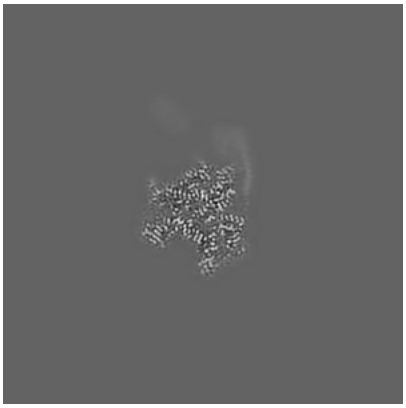


Z Index: 200

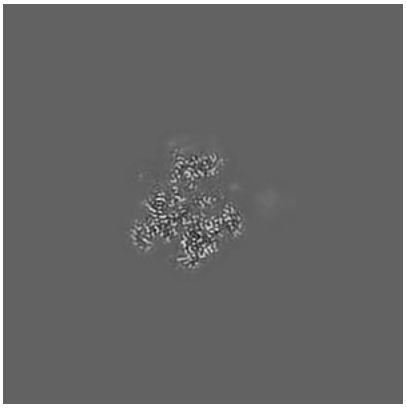
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices (i)

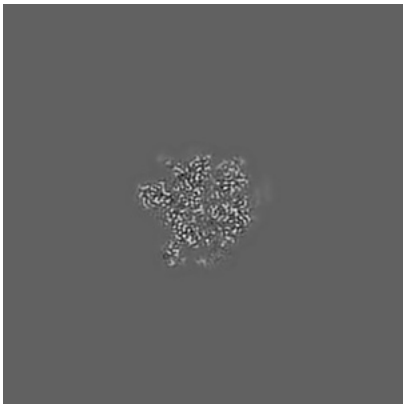
6.3.1 Primary map



X Index: 176



Y Index: 203

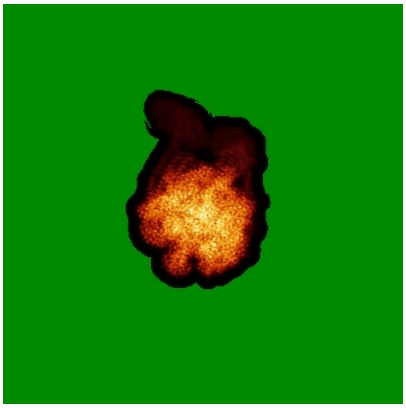


Z Index: 182

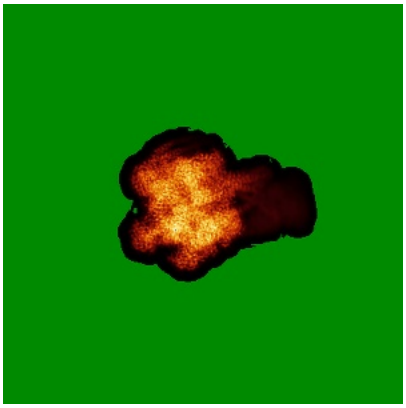
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) (i)

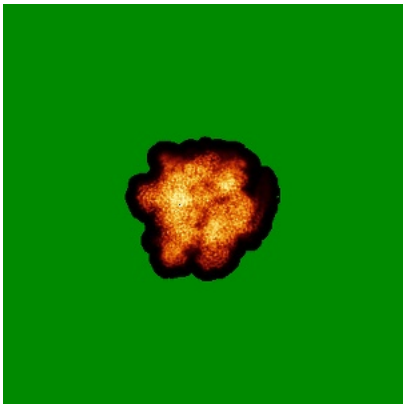
6.4.1 Primary map



X



Y

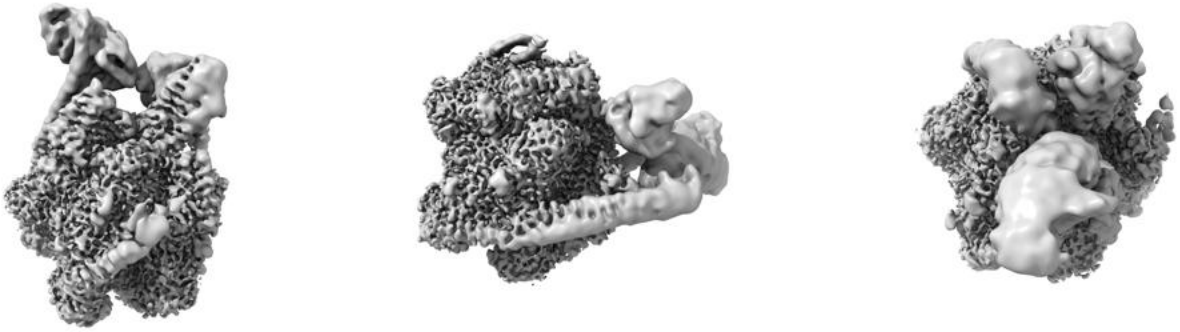


Z

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 0.3. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

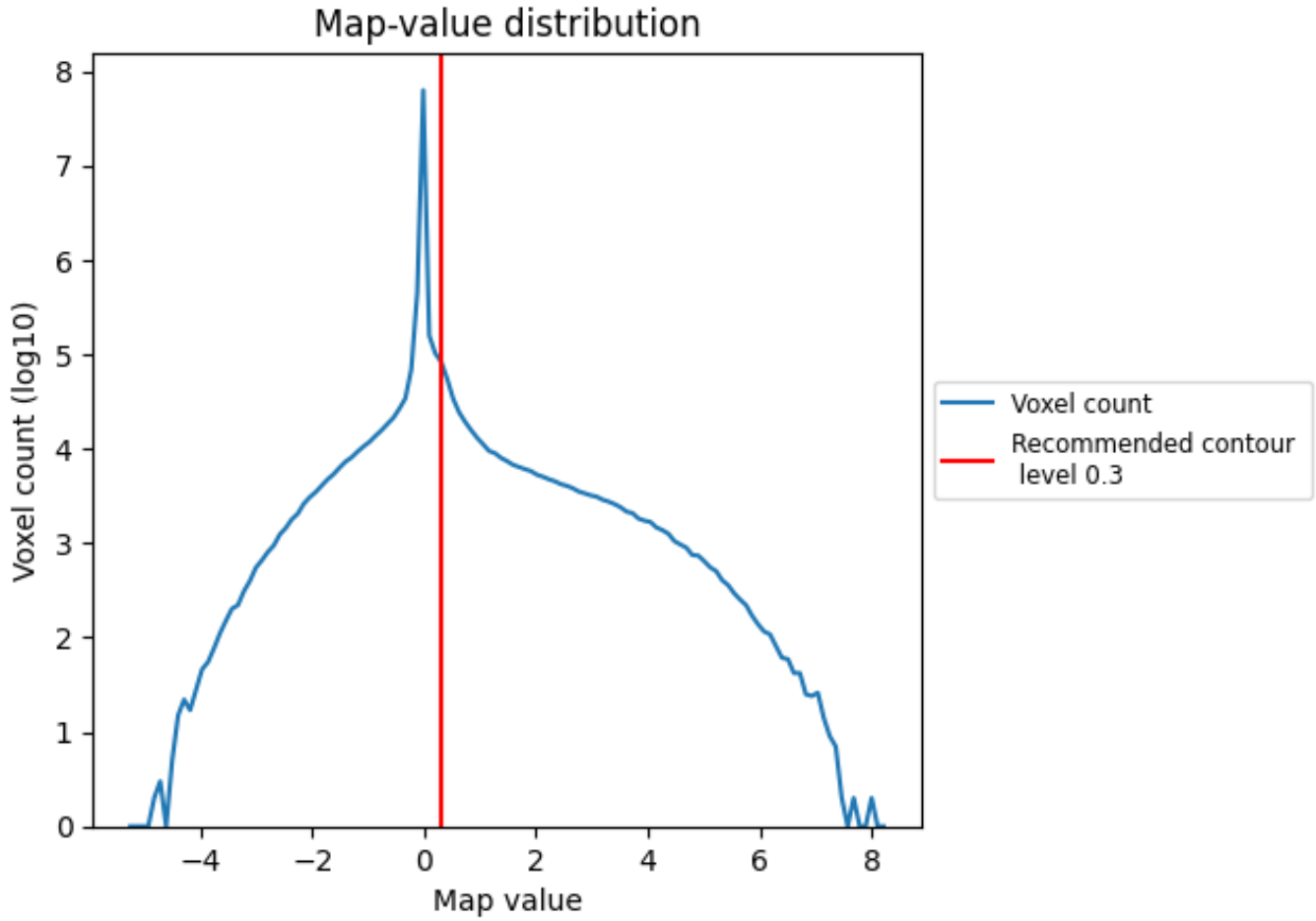
6.6 Mask visualisation (i)

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

7 Map analysis

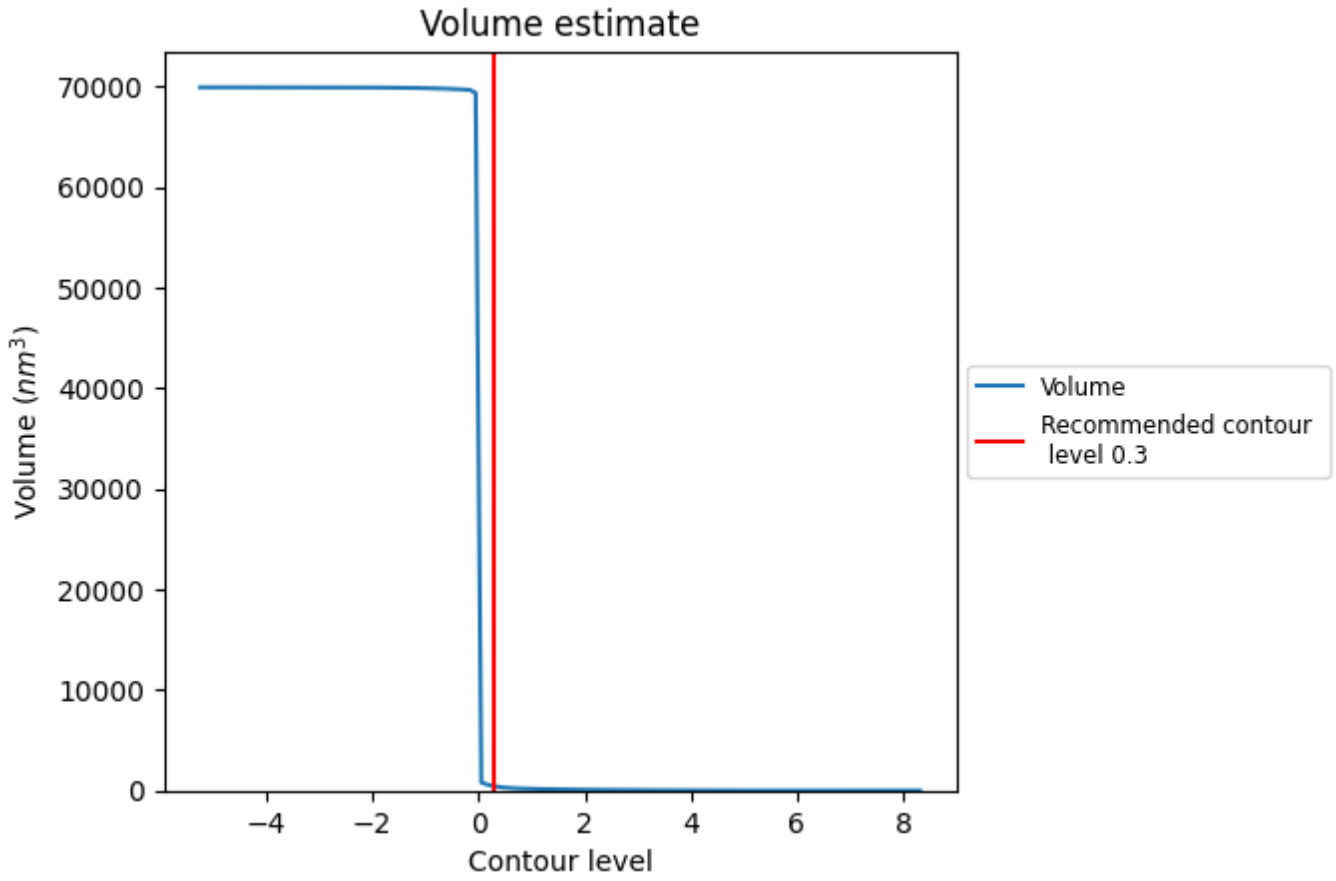
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution



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

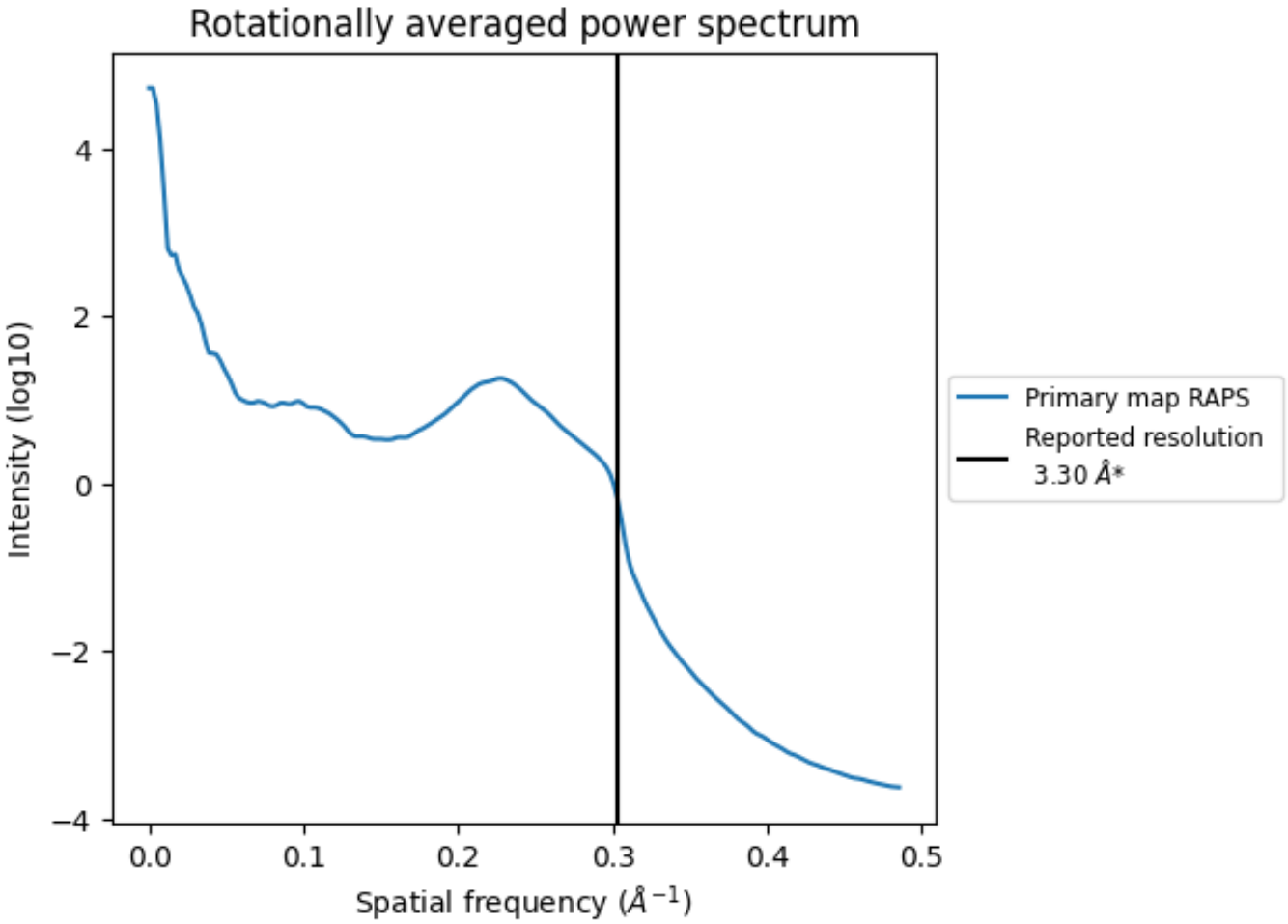
7.2 Volume estimate (i)



The volume at the recommended contour level is 434 nm^3; this corresponds to an approximate mass of 392 kDa.

The volume estimate graph shows how the enclosed volume varies with the contour level. The recommended contour level is shown as a vertical line and the intersection between the line and the curve gives the volume of the enclosed surface at the given level.

7.3 Rotationally averaged power spectrum (i)



*Reported resolution corresponds to spatial frequency of 0.303 $Å^{-1}$

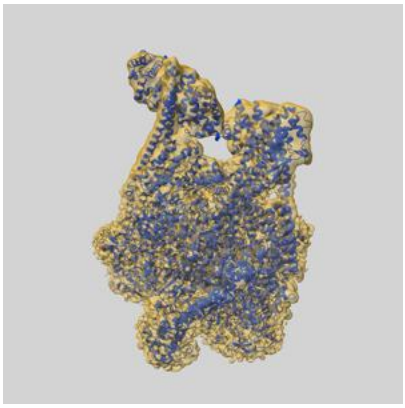
8 Fourier-Shell correlation (i)

This section was not generated. No FSC curve or half-maps provided.

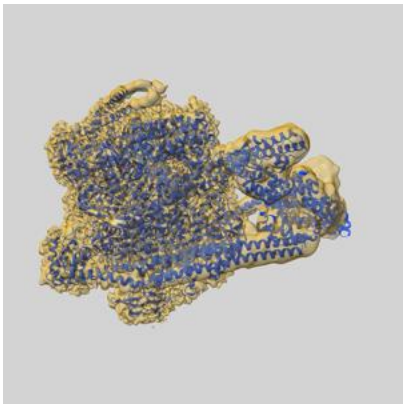
9 Map-model fit

This section contains information regarding the fit between EMDB map EMD-25999 and PDB model 7TMQ. Per-residue inclusion information can be found in section 3 on page 8.

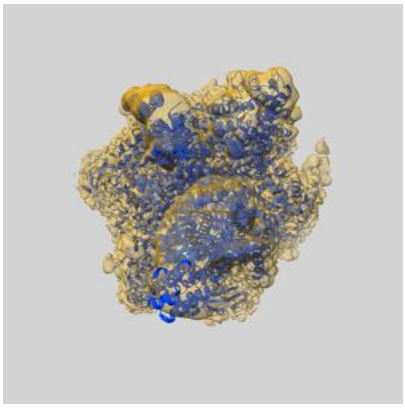
9.1 Map-model overlay



X



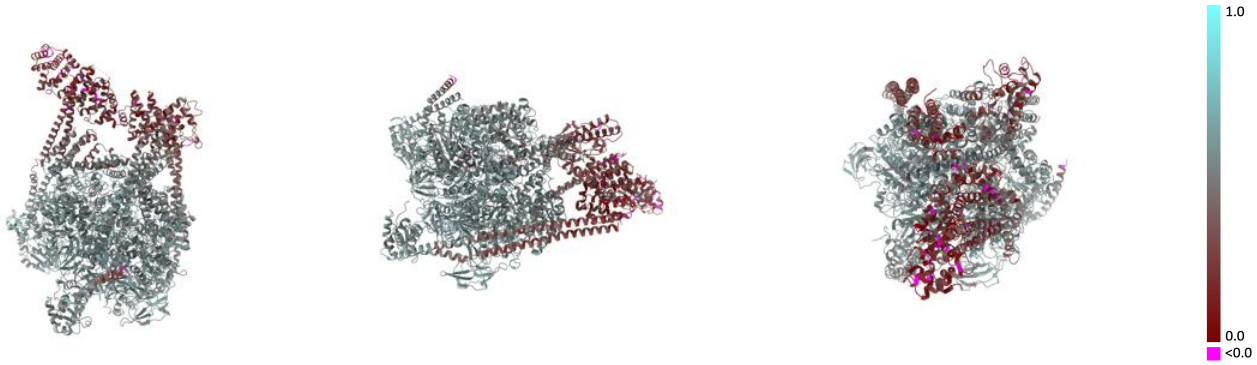
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.3 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.

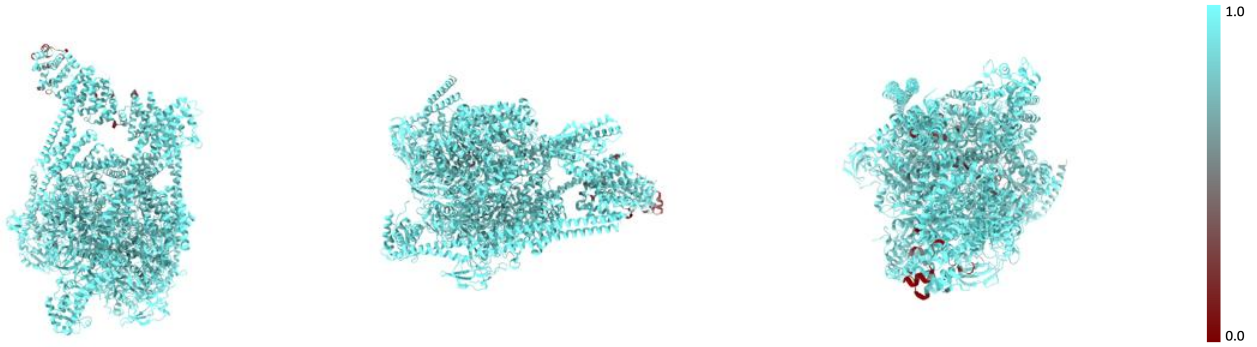
9.2 Q-score mapped to coordinate model



X Y Z

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

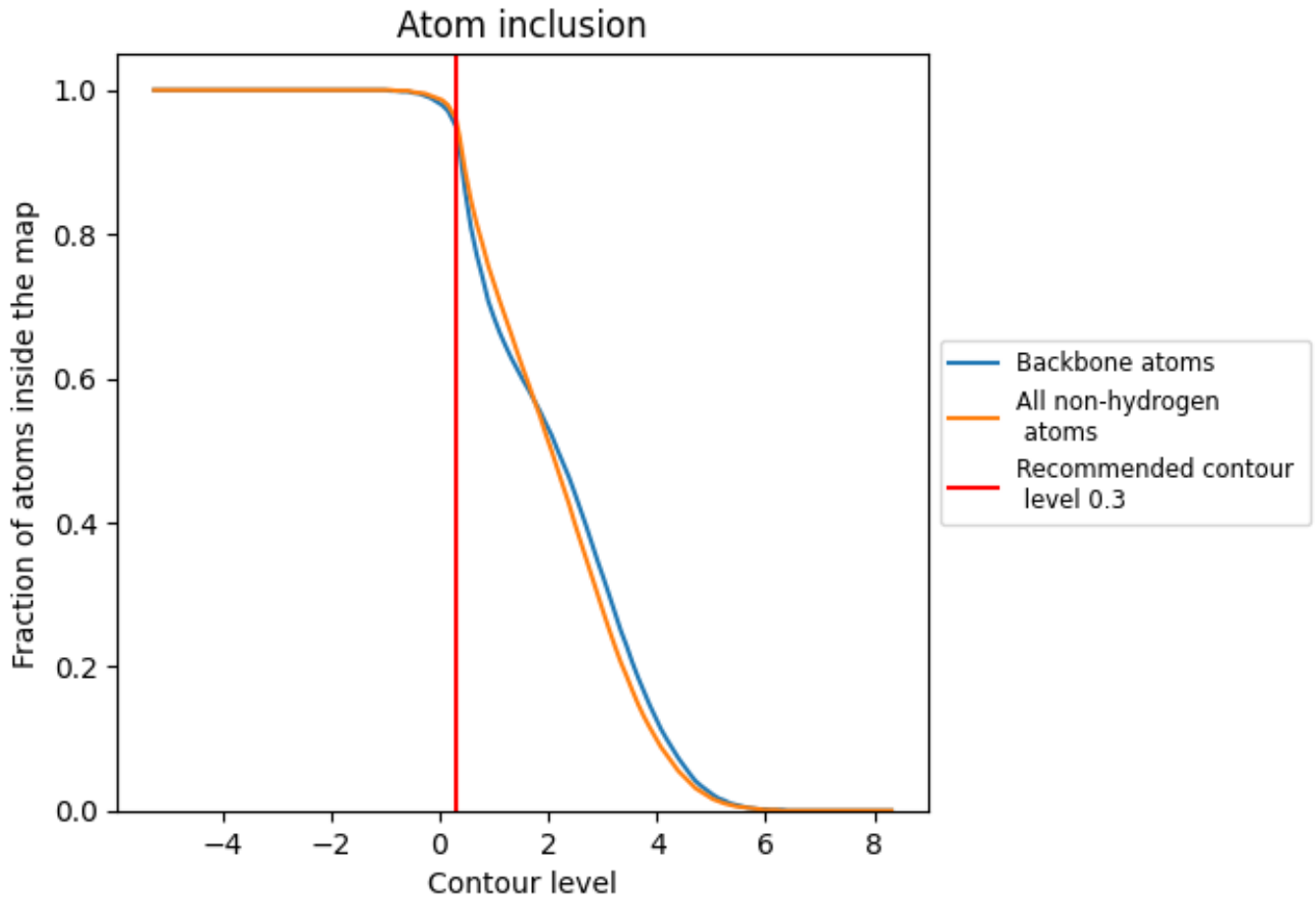
9.3 Atom inclusion mapped to coordinate model



X Y Z

The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.3).

9.4 Atom inclusion (i)



At the recommended contour level, 95% of all backbone atoms, 96% 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 (0.3) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.9610	0.5030
A	0.9690	0.5580
B	0.9630	0.5570
C	0.9690	0.5260
D	0.9620	0.5540
E	0.9590	0.5410
F	0.9690	0.5620
G	0.9730	0.5300
H	0.9570	0.4540
I	0.9810	0.4920
J	0.9790	0.3860
K	0.9800	0.5020
L	0.9750	0.3950
M	0.9580	0.4290
N	0.9910	0.3180
P	0.8780	0.1870