



## Full wwPDB EM Validation Report ⓘ

Nov 13, 2022 – 08:34 AM EST

PDB ID : 6VMG  
EMDB ID : EMD-21241  
Title : Chloroplast ATP synthase (O3, CF1FO)  
Authors : Yang, J.-H.; Williams, D.; Kandiah, E.; Fromme, P.; Chiu, P.-L.  
Deposited on : 2020-01-27  
Resolution : 6.46 Å(reported)  
Based on initial model : 6FKI

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

We welcome your comments at [validation@mail.wwpdb.org](mailto: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>.

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

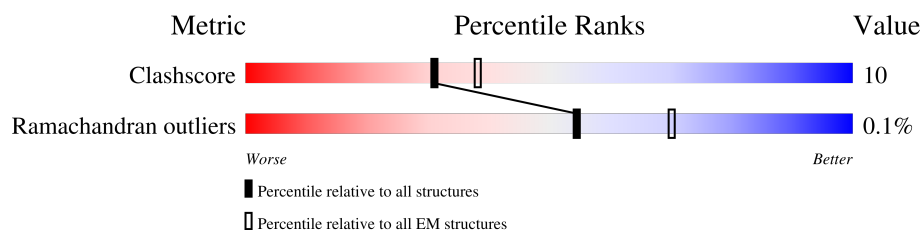
EMDB validation analysis : 0.0.1.dev43  
MolProbity : 4.02b-467  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
MapQ : 1.9.9  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP)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

# 1 Overall quality at a glance

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

The reported resolution of this entry is 6.46 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

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  $\geq 3$ , 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	a	247	<div> <div>9%</div> <div>89%</div> <div>10%</div> </div>
2	I	184	<div> <div>77%</div> <div>20%</div> </div>
3	d	257	<div> <div>68%</div> <div>31%</div> </div>
4	J	222	<div> <div>57%</div> <div>41%</div> </div>
5	e	134	<div> <div>9%</div> <div>96%</div> </div>
6	g	364	<div> <div>88%</div> <div>12%</div> </div>
7	M	81	<div> <div>31%</div> <div>93%</div> <div>5%</div> </div>
7	N	81	<div> <div>32%</div> <div>93%</div> <div>5%</div> </div>
7	O	81	<div> <div>17%</div> <div>93%</div> <div>5%</div> </div>

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Mol	Chain	Length	Quality of chain
7	P	81	
7	Q	81	
7	R	81	
7	S	81	
7	T	81	
7	U	81	
7	V	81	
7	W	81	
7	X	81	
7	Y	81	
7	Z	81	
8	A	507	
8	B	507	
8	C	507	
9	D	498	
9	E	498	
9	F	498	

## 2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 25375 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 ATP synthase subunit a, chloroplastic.

Mol	Chain	Residues	Atoms				AltConf	Trace
1	a	222	Total	C	N	O	0	0
			1091	647	222	222		

- Molecule 2 is a protein called ATP synthase subunit b, chloroplastic.

Mol	Chain	Residues	Atoms				AltConf	Trace
2	I	148	Total	C	N	O	0	0
			731	435	148	148		

- Molecule 3 is a protein called ATP synthase delta chain, chloroplastic.

Mol	Chain	Residues	Atoms				AltConf	Trace
3	d	177	Total	C	N	O	0	0
			877	523	177	177		

- Molecule 4 is a protein called ATP synthase subunit b', chloroplastic.

Mol	Chain	Residues	Atoms				AltConf	Trace
4	J	131	Total	C	N	O	0	0
			651	389	131	131		

- Molecule 5 is a protein called ATP synthase epsilon chain, chloroplastic.

Mol	Chain	Residues	Atoms				AltConf	Trace
5	e	129	Total	C	N	O	0	0
			637	379	129	129		

- Molecule 6 is a protein called ATP synthase gamma chain, chloroplastic.

Mol	Chain	Residues	Atoms				AltConf	Trace
6	g	321	Total	C	N	O	0	0
			1590	948	321	321		

- Molecule 7 is a protein called ATP synthase subunit c, chloroplastic.

Mol	Chain	Residues	Atoms				AltConf	Trace
7	Z	79	Total	C	N	O	0	0
			384	226	79	79		
7	Q	79	Total	C	N	O	0	0
			384	226	79	79		
7	P	79	Total	C	N	O	0	0
			384	226	79	79		
7	O	79	Total	C	N	O	0	0
			384	226	79	79		
7	N	79	Total	C	N	O	0	0
			384	226	79	79		
7	M	79	Total	C	N	O	0	0
			384	226	79	79		
7	R	79	Total	C	N	O	0	0
			384	226	79	79		
7	Y	79	Total	C	N	O	0	0
			384	226	79	79		
7	X	79	Total	C	N	O	0	0
			384	226	79	79		
7	W	79	Total	C	N	O	0	0
			384	226	79	79		
7	S	79	Total	C	N	O	0	0
			384	226	79	79		
7	T	79	Total	C	N	O	0	0
			384	226	79	79		
7	U	79	Total	C	N	O	0	0
			384	226	79	79		
7	V	79	Total	C	N	O	0	0
			384	226	79	79		

- Molecule 8 is a protein called ATP synthase subunit alpha, chloroplastic.

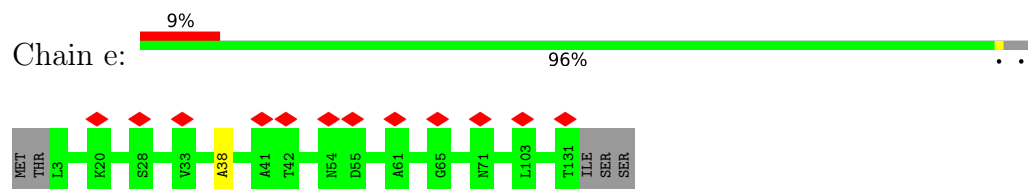
Mol	Chain	Residues	Atoms				AltConf	Trace
8	A	495	Total	C	N	O	0	0
			2436	1446	495	495		
8	C	501	Total	C	N	O	0	0
			2466	1464	501	501		
8	B	504	Total	C	N	O	0	0
			2481	1473	504	504		

- Molecule 9 is a protein called ATP synthase subunit beta, chloroplastic.

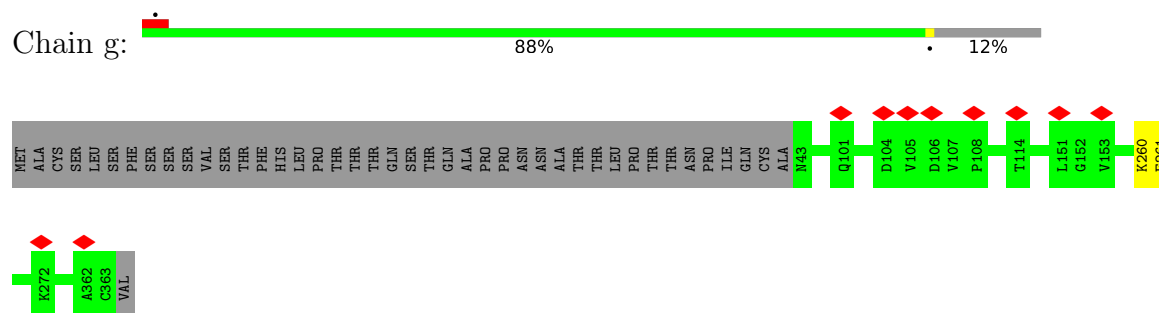
Mol	Chain	Residues	Atoms				AltConf	Trace
9	F	479	Total 2348	C 1390	N 479	O 479	0	0
9	E	478	Total 2343	C 1387	N 478	O 478	0	0
9	D	479	Total 2348	C 1390	N 479	O 479	0	0



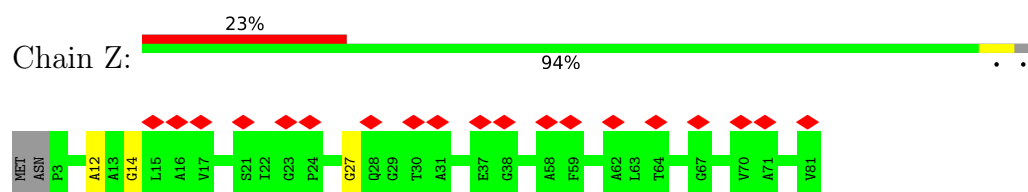
- Molecule 5: ATP synthase epsilon chain, chloroplastic



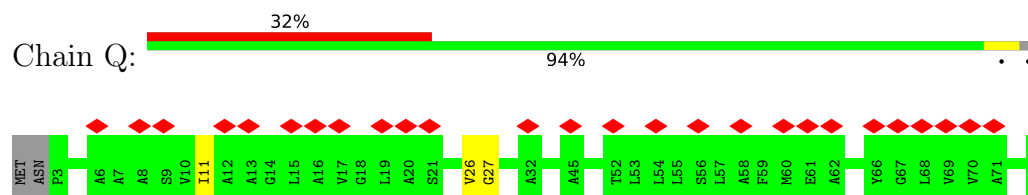
- Molecule 6: ATP synthase gamma chain, chloroplastic



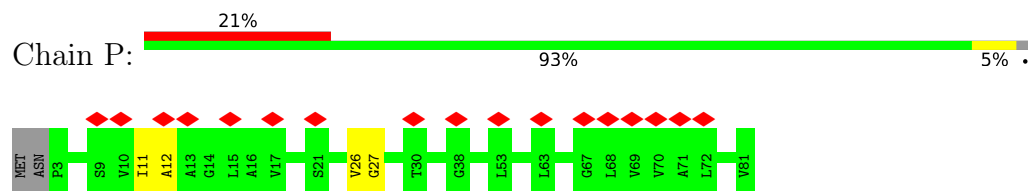
- Molecule 7: ATP synthase subunit c, chloroplastic



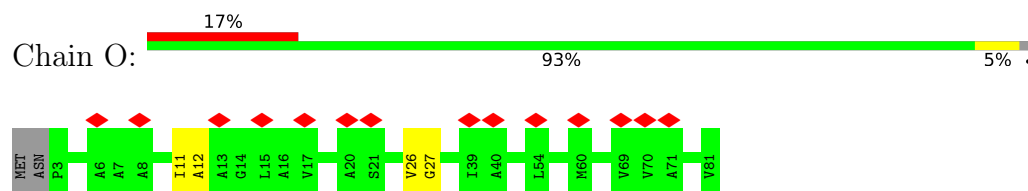
- Molecule 7: ATP synthase subunit c, chloroplastic



- Molecule 7: ATP synthase subunit c, chloroplastic

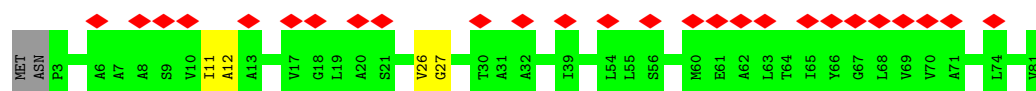


- Molecule 7: ATP synthase subunit c, chloroplastic

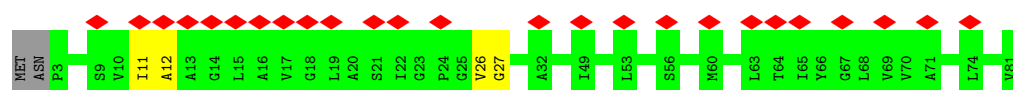


- Molecule 7: ATP synthase subunit c, chloroplastic

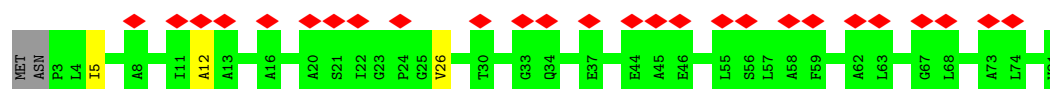




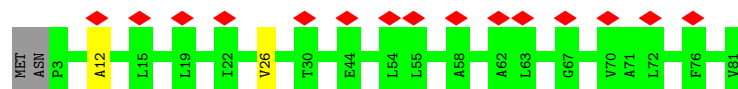
- Molecule 7: ATP synthase subunit c, chloroplastic



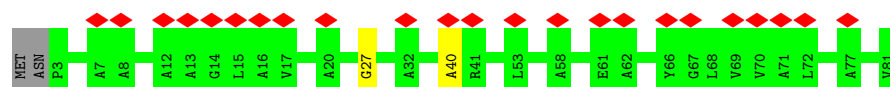
- Molecule 7: ATP synthase subunit c, chloroplastic



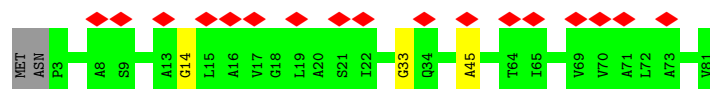
- Molecule 7: ATP synthase subunit c, chloroplastic



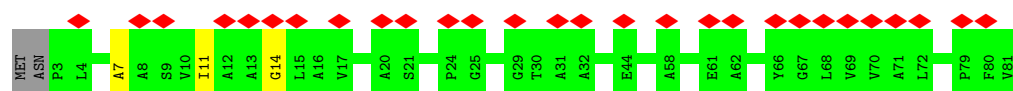
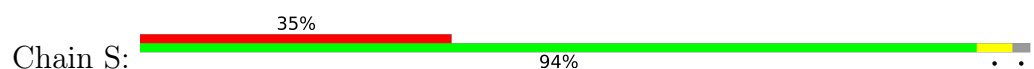
- Molecule 7: ATP synthase subunit c, chloroplastic



- Molecule 7: ATP synthase subunit c, chloroplastic



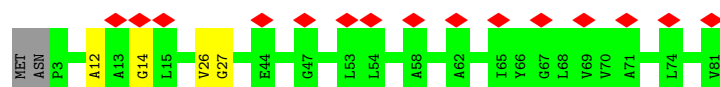
- Molecule 7: ATP synthase subunit c, chloroplastic



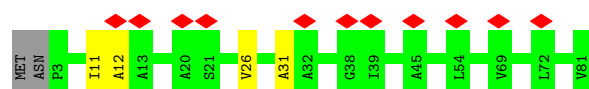
- Molecule 7: ATP synthase subunit c, chloroplastic



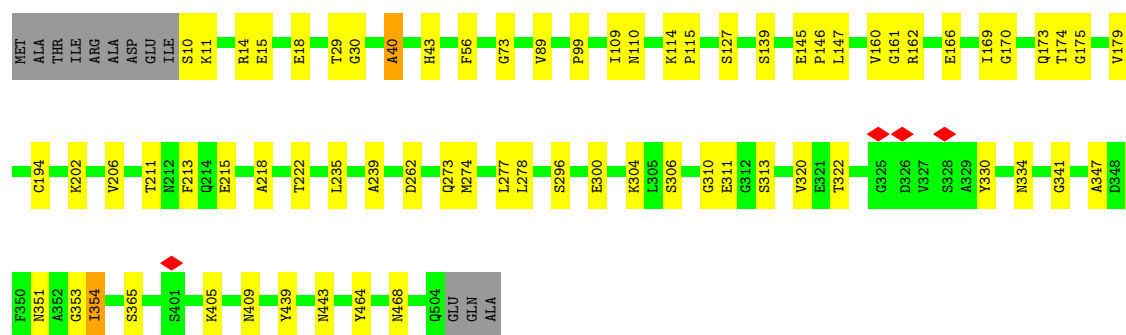
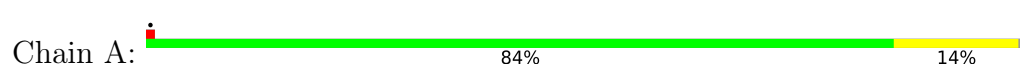
- Molecule 7: ATP synthase subunit c, chloroplastic



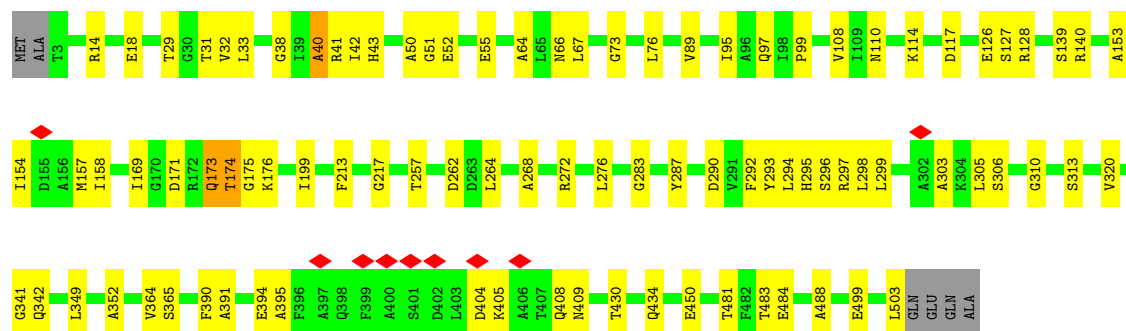
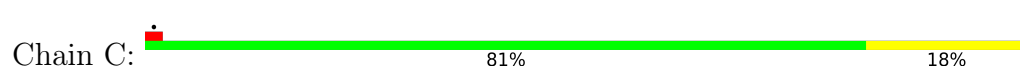
- Molecule 7: ATP synthase subunit c, chloroplastic



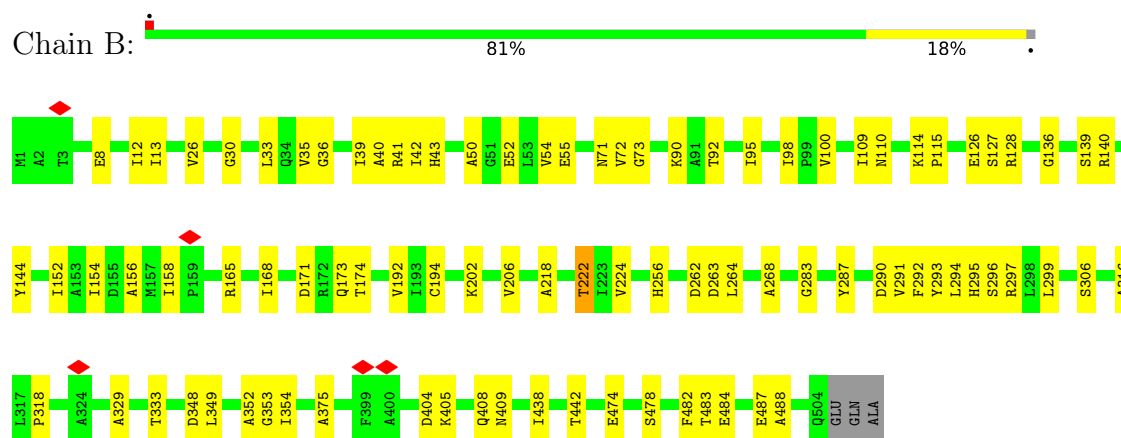
- Molecule 8: ATP synthase subunit alpha, chloroplastic

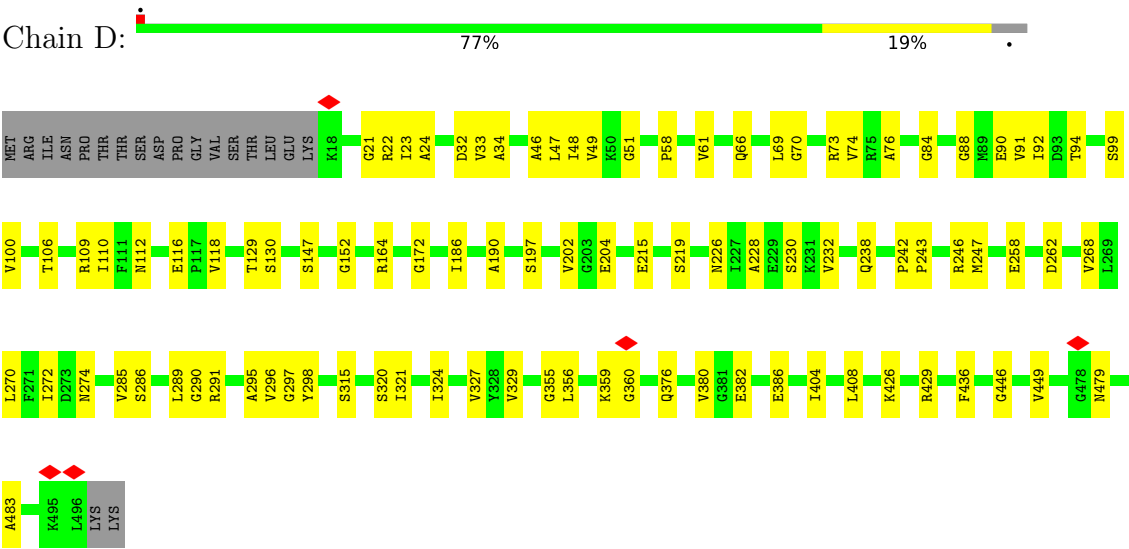


- Molecule 8: ATP synthase subunit alpha, chloroplastic



- Molecule 8: ATP synthase subunit alpha, chloroplastic





## 4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	29090	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 ( $e^-/\text{\AA}^2$ )	43.5	Depositor
Minimum defocus (nm)	-1500	Depositor
Maximum defocus (nm)	-4000	Depositor
Magnification	48077	Depositor
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.137	Depositor
Minimum map value	-0.057	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.006	Depositor
Recommended contour level	0.035	Depositor
Map size (Å)	374.4, 374.4, 374.4	wwPDB
Map dimensions	180, 180, 180	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	2.08, 2.08, 2.08	Depositor

## 5 Model quality

### 5.1 Standard geometry

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	0.32	0/1090	0.44	0/1513
2	I	0.32	0/730	0.36	0/1015
3	d	0.38	0/876	0.47	0/1220
4	J	0.30	0/649	0.36	0/903
5	e	0.39	0/636	0.61	0/884
6	g	0.38	0/1589	0.49	0/2214
7	M	0.31	0/383	0.38	0/528
7	N	0.31	0/383	0.38	0/528
7	O	0.31	0/383	0.38	0/528
7	P	0.31	0/383	0.38	0/528
7	Q	0.31	0/383	0.38	0/528
7	R	0.31	0/383	0.38	0/528
7	S	0.31	0/383	0.38	0/528
7	T	0.31	0/383	0.38	0/528
7	U	0.31	0/383	0.38	0/528
7	V	0.31	0/383	0.38	0/528
7	W	0.31	0/383	0.38	0/528
7	X	0.31	0/383	0.38	0/528
7	Y	0.31	0/383	0.38	0/528
7	Z	0.31	0/383	0.38	0/528
8	A	0.42	0/2435	0.53	0/3384
8	B	0.43	1/2480 (0.0%)	0.57	0/3447
8	C	0.42	0/2465	0.56	0/3426
9	D	0.42	0/2347	0.57	0/3256
9	E	0.41	0/2342	0.56	0/3249
9	F	0.42	0/2347	0.56	0/3256
All	All	0.39	1/25348 (0.0%)	0.51	0/35159

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	a	0	1
3	d	0	1
5	e	0	1
6	g	0	1
8	A	0	5
8	B	0	4
8	C	0	6
9	D	0	1
9	E	0	2
9	F	0	4
All	All	0	26

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	B	144	TYR	C-N	-6.93	1.18	1.34

There are no bond angle outliers.

There are no chirality outliers.

All (26) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
8	A	173	GLN	Peptide
8	A	174	THR	Peptide
8	A	262	ASP	Peptide
8	A	354	ILE	Peptide
8	A	40	ALA	Peptide
8	B	173	GLN	Peptide
8	B	222	THR	Peptide
8	B	26	VAL	Peptide
8	B	262	ASP	Peptide
8	C	171	ASP	Peptide
8	C	173	GLN	Peptide
8	C	174	THR	Peptide
8	C	303	ALA	Peptide
8	C	40	ALA	Peptide
8	C	450	GLU	Peptide
9	D	228	ALA	Peptide
9	E	221	VAL	Peptide
9	E	52	ARG	Peptide
9	F	221	VAL	Peptide
9	F	228	ALA	Peptide

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Mol	Chain	Res	Type	Group
9	F	242	PRO	Peptide
9	F	295	ALA	Peptide
1	a	228	ALA	Peptide
3	d	158	VAL	Peptide
5	e	38	ALA	Peptide
6	g	260	LYS	Peptide

## 5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	a	1091	0	499	0	0
2	I	731	0	331	4	0
3	d	877	0	402	0	0
4	J	651	0	304	3	0
5	e	637	0	303	0	0
6	g	1590	0	734	0	0
7	M	384	0	216	4	0
7	N	384	0	216	4	0
7	O	384	0	216	4	0
7	P	384	0	216	4	0
7	Q	384	0	216	3	0
7	R	384	0	216	4	0
7	S	384	0	216	3	0
7	T	384	0	216	2	0
7	U	384	0	216	4	0
7	V	384	0	216	4	0
7	W	384	0	216	3	0
7	X	384	0	216	2	0
7	Y	384	0	216	2	0
7	Z	384	0	216	3	0
8	A	2436	0	1164	39	0
8	B	2481	0	1190	52	0
8	C	2466	0	1179	50	0
9	D	2348	0	1105	51	0
9	E	2343	0	1103	48	0
9	F	2348	0	1105	44	0
All	All	25375	0	12443	309	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.

All (309) 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)
9:E:48:ILE:HA	9:E:61:VAL:O	1.62	1.00
9:E:47:LEU:O	9:E:62:THR:HA	1.65	0.95
8:A:169:ILE:HA	8:A:320:VAL:O	1.69	0.93
9:E:100:VAL:O	9:E:129:THR:HA	1.72	0.89
9:E:99:SER:HA	9:E:130:SER:O	1.73	0.87
9:E:271:PHE:HA	9:E:324:ILE:O	1.76	0.86
9:D:47:LEU:HA	9:D:92:ILE:O	1.76	0.86
8:C:40:ALA:O	8:C:73:GLY:HA2	1.75	0.85
9:D:51:GLY:O	9:D:58:PRO:HA	1.76	0.85
9:D:21:GLY:O	9:D:90:GLU:HA	1.76	0.83
8:B:40:ALA:O	8:B:73:GLY:HA2	1.79	0.83
8:A:194:CYS:O	8:A:222:THR:HA	1.78	0.82
8:B:194:CYS:O	8:B:222:THR:HA	1.78	0.81
8:C:484:GLU:O	8:C:488:ALA:HB2	1.82	0.79
9:E:168:ILE:O	9:E:324:ILE:HA	1.83	0.78
9:E:167:LYS:HA	9:E:323:SER:O	1.86	0.76
9:F:99:SER:HA	9:F:130:SER:O	1.85	0.75
9:F:200:GLY:HA2	9:F:235:VAL:O	1.86	0.74
8:B:192:VAL:HA	8:B:256:HIS:O	1.89	0.72
9:D:33:VAL:O	9:D:73:ARG:HA	1.92	0.70
9:F:100:VAL:O	9:F:129:THR:HA	1.92	0.69
9:E:46:ALA:HA	9:E:63:CYS:O	1.93	0.68
8:A:29:THR:HA	8:A:89:VAL:O	1.93	0.67
9:D:48:ILE:HA	9:D:61:VAL:O	1.95	0.67
9:D:66:GLN:H	9:D:76:ALA:HA	1.59	0.67
8:C:55:GLU:O	8:C:89:VAL:HA	1.95	0.66
9:F:69:LEU:H	9:F:74:VAL:HA	1.60	0.66
8:C:290:ASP:O	8:C:294:LEU:N	2.28	0.66
8:B:329:ALA:O	8:B:333:THR:N	2.29	0.65
9:D:99:SER:HA	9:D:130:SER:O	1.96	0.65
7:Q:27:GLY:HA3	7:R:26:VAL:HA	1.79	0.64
9:E:51:GLY:HA3	9:E:59:MET:HA	1.80	0.64
8:A:56:PHE:HA	8:A:89:VAL:HA	1.77	0.64
8:A:166:GLU:HA	8:A:341:GLY:O	1.98	0.64
9:F:259:TYR:O	9:F:263:VAL:N	2.29	0.64
7:R:12:ALA:HB2	7:S:11:ILE:HA	1.78	0.64
7:Z:12:ALA:HB2	7:M:11:ILE:HA	1.80	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:D:100:VAL:O	9:D:129:THR:HA	1.98	0.63
8:A:349:LEU:O	8:A:354:ILE:N	2.33	0.62
7:W:14:GLY:HA3	7:V:12:ALA:HB1	1.81	0.62
8:C:50:ALA:O	9:E:84:GLY:N	2.31	0.62
8:C:292:PHE:O	8:C:296:SER:N	2.32	0.62
8:B:154:ILE:O	8:B:158:ILE:N	2.33	0.62
8:C:169:ILE:HA	8:C:320:VAL:O	1.99	0.62
8:B:110:ASN:N	8:B:114:LYS:O	2.33	0.61
8:A:274:MET:O	8:A:278:LEU:N	2.33	0.61
9:F:19:ASN:O	9:F:92:ILE:HA	2.01	0.61
8:C:342:GLN:O	8:C:364:VAL:N	2.33	0.61
8:A:341:GLY:HA2	8:A:365:SER:HA	1.82	0.61
9:F:436:PHE:HA	9:F:446:GLY:HA3	1.83	0.60
7:Q:26:VAL:HA	7:P:27:GLY:HA3	1.82	0.60
9:D:355:GLY:O	9:D:359:LYS:N	2.32	0.60
8:C:33:LEU:N	8:C:41:ARG:O	2.35	0.60
9:F:53:ASP:N	9:F:57:GLN:O	2.36	0.59
9:E:152:GLY:O	9:E:449:VAL:N	2.36	0.59
9:E:149:PHE:N	9:E:162:TYR:O	2.32	0.59
8:C:391:ALA:O	8:C:395:ALA:N	2.36	0.59
8:C:405:LYS:O	8:C:409:ASN:CB	2.50	0.59
9:F:315:SER:HA	9:F:320:SER:HA	1.85	0.58
8:C:32:VAL:HA	8:C:42:ILE:HA	1.85	0.58
8:C:51:GLY:H	8:C:64:ALA:HB3	1.67	0.58
8:B:165:ARG:HA	8:B:316:ALA:HB3	1.84	0.58
9:D:197:SER:O	9:D:232:VAL:HA	2.03	0.58
9:E:30:VAL:HA	9:E:76:ALA:O	2.04	0.58
9:F:46:ALA:N	9:F:96:ALA:O	2.36	0.58
9:F:470:PRO:O	9:F:474:PHE:N	2.37	0.58
8:C:293:TYR:O	8:C:297:ARG:N	2.32	0.58
8:A:110:ASN:N	8:A:114:LYS:O	2.34	0.57
8:B:35:VAL:HA	8:B:40:ALA:HA	1.86	0.57
7:R:12:ALA:HB1	7:S:14:GLY:HA3	1.86	0.57
9:E:23:ILE:O	9:E:88:GLY:N	2.37	0.57
7:Q:11:ILE:HA	7:P:12:ALA:HB2	1.86	0.57
9:E:352:LEU:HA	9:E:365:VAL:HA	1.86	0.57
9:E:52:ARG:HA	9:E:57:GLN:HA	1.86	0.57
8:A:175:GLY:O	8:A:179:VAL:N	2.38	0.57
9:F:356:LEU:O	9:F:361:ILE:N	2.36	0.57
8:B:41:ARG:HA	8:B:72:VAL:O	2.04	0.57
9:F:223:ASN:H	9:F:230:SER:HA	1.70	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:C:14:ARG:O	8:C:18:GLU:N	2.35	0.57
8:C:99:PRO:HA	8:C:126:GLU:O	2.04	0.56
8:C:404:ASP:O	8:C:408:GLN:N	2.27	0.56
7:Z:27:GLY:HA3	7:M:26:VAL:HA	1.88	0.56
9:F:111:PHE:HA	9:F:117:PRO:HA	1.87	0.56
8:C:484:GLU:O	8:C:488:ALA:CB	2.53	0.56
9:D:23:ILE:O	9:D:88:GLY:N	2.30	0.56
9:D:172:GLY:N	9:D:327:VAL:O	2.38	0.56
7:U:27:GLY:HA3	7:V:26:VAL:HA	1.86	0.56
9:E:166:GLY:H	9:E:322:THR:HA	1.70	0.56
8:B:291:VAL:O	8:B:295:HIS:N	2.30	0.56
9:D:479:ASN:O	9:D:483:ALA:N	2.38	0.55
7:P:26:VAL:HA	7:O:27:GLY:HA3	1.88	0.55
8:A:30:GLY:CA	8:A:43:HIS:O	2.55	0.55
8:A:139:SER:O	8:A:306:SER:N	2.38	0.55
9:E:166:GLY:O	9:E:323:SER:N	2.39	0.55
8:C:154:ILE:O	8:C:158:ILE:N	2.39	0.55
8:C:390:PHE:O	8:C:394:GLU:CB	2.55	0.55
9:E:459:GLN:O	9:E:463:SER:CB	2.54	0.55
8:B:348:ASP:O	8:B:352:ALA:N	2.39	0.55
9:D:382:GLU:O	9:D:386:GLU:CB	2.55	0.55
8:C:97:GLN:HA	8:C:128:ARG:O	2.07	0.54
9:E:148:ILE:HA	9:E:163:ARG:HA	1.88	0.54
8:B:8:GLU:O	8:B:13:ILE:N	2.28	0.54
8:C:264:LEU:O	8:C:268:ALA:N	2.41	0.54
8:B:349:LEU:O	8:B:354:ILE:N	2.39	0.54
8:B:290:ASP:O	8:B:294:LEU:N	2.35	0.54
9:E:147:SER:O	9:E:164:ARG:N	2.39	0.53
8:A:211:THR:O	8:A:215:GLU:CB	2.57	0.53
8:A:464:TYR:O	8:A:468:ASN:CB	2.55	0.53
8:C:110:ASN:N	8:C:114:LYS:O	2.42	0.53
8:A:235:LEU:O	8:A:239:ALA:HB2	2.09	0.53
8:C:342:GLN:N	8:C:364:VAL:O	2.39	0.53
8:B:202:LYS:O	8:B:206:VAL:N	2.37	0.53
9:E:172:GLY:O	9:E:329:VAL:N	2.37	0.53
8:A:306:SER:O	8:A:310:GLY:N	2.42	0.53
8:B:264:LEU:O	8:B:268:ALA:N	2.40	0.53
8:C:66:ASN:O	8:C:73:GLY:N	2.40	0.52
8:B:152:ILE:O	8:B:156:ALA:HB2	2.09	0.52
8:B:405:LYS:O	8:B:409:ASN:CB	2.56	0.52
9:E:66:GLN:H	9:E:76:ALA:HA	1.74	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:A:213:PHE:O	8:A:218:ALA:N	2.43	0.52
8:A:109:ILE:HA	8:A:115:PRO:HA	1.92	0.52
8:A:439:TYR:O	8:A:443:ASN:CB	2.58	0.52
8:B:349:LEU:HA	8:B:352:ALA:HB3	1.92	0.52
9:E:25:GLN:O	9:E:32:ASP:N	2.38	0.52
9:D:243:PRO:O	9:D:247:MET:N	2.41	0.52
8:A:273:GLN:O	8:A:277:LEU:CB	2.58	0.52
8:C:31:THR:O	8:C:43:HIS:N	2.42	0.52
9:D:268:VAL:N	9:D:320:SER:O	2.42	0.52
9:F:353:SER:N	9:F:364:ALA:O	2.40	0.51
9:E:69:LEU:H	9:E:74:VAL:HA	1.75	0.51
9:E:100:VAL:H	9:E:130:SER:H	1.57	0.51
8:B:98:ILE:O	8:B:127:SER:HA	2.10	0.51
9:E:102:VAL:N	9:E:128:THR:O	2.44	0.51
8:B:484:GLU:O	8:B:488:ALA:CB	2.59	0.51
8:A:30:GLY:HA2	8:A:43:HIS:O	2.11	0.51
9:F:294:SER:N	9:F:298:TYR:O	2.43	0.51
8:B:218:ALA:O	8:B:222:THR:N	2.43	0.51
7:P:11:ILE:HA	7:O:12:ALA:HB2	1.91	0.51
8:A:347:ALA:O	8:A:351:ASN:N	2.34	0.51
9:F:314:THR:O	9:F:321:ILE:N	2.39	0.51
8:A:11:LYS:O	8:A:15:GLU:CB	2.58	0.51
9:F:376:GLN:O	9:F:380:VAL:N	2.34	0.51
9:D:147:SER:O	9:D:164:ARG:N	2.33	0.51
9:D:272:ILE:N	9:D:324:ILE:O	2.44	0.51
8:B:55:GLU:N	8:B:90:LYS:O	2.44	0.50
9:D:186:ILE:O	9:D:190:ALA:CB	2.59	0.50
9:E:439:GLU:HA	9:E:443:GLY:H	1.77	0.50
9:D:295:ALA:O	9:D:298:TYR:N	2.42	0.50
8:C:272:ARG:O	8:C:276:LEU:CB	2.60	0.50
9:F:215:GLU:O	9:F:219:SER:N	2.33	0.50
9:E:243:PRO:O	9:E:247:MET:N	2.45	0.50
9:D:376:GLN:O	9:D:380:VAL:N	2.41	0.50
8:C:153:ALA:O	8:C:157:MET:CB	2.60	0.49
8:C:257:THR:N	8:C:313:SER:O	2.43	0.49
9:D:286:SER:O	9:D:291:ARG:N	2.45	0.49
8:C:52:GLU:HA	8:C:95:ILE:HA	1.93	0.49
8:C:108:VAL:O	8:C:117:ASP:N	2.44	0.49
9:F:417:ASP:O	9:F:421:VAL:N	2.31	0.49
8:B:139:SER:O	8:B:306:SER:N	2.34	0.49
9:F:382:GLU:O	9:F:386:GLU:CB	2.61	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:E:353:SER:N	9:E:364:ALA:O	2.44	0.49
8:B:295:HIS:O	8:B:299:LEU:N	2.44	0.49
8:B:52:GLU:HA	8:B:95:ILE:HA	1.95	0.49
7:X:40:ALA:HA	7:W:45:ALA:HB3	1.93	0.49
8:C:306:SER:O	8:C:310:GLY:N	2.44	0.49
9:D:110:ILE:O	9:D:118:VAL:N	2.46	0.49
9:D:285:VAL:O	9:D:289:LEU:N	2.39	0.49
4:J:198:ASP:O	4:J:202:SER:N	2.45	0.49
9:F:419:LEU:O	9:F:423:ARG:CB	2.61	0.49
8:C:199:ILE:N	8:C:262:ASP:O	2.46	0.49
8:B:33:LEU:H	8:B:42:ILE:HA	1.77	0.49
8:A:14:ARG:O	8:A:18:GLU:CB	2.60	0.48
9:F:214:MET:O	9:F:218:GLU:CB	2.61	0.48
9:E:46:ALA:N	9:E:96:ALA:O	2.29	0.48
8:B:283:GLY:N	8:B:287:TYR:O	2.44	0.48
9:F:275:ILE:N	9:F:326:ALA:O	2.38	0.48
9:D:172:GLY:O	9:D:329:VAL:N	2.46	0.48
8:A:304:LYS:HA	8:A:313:SER:HA	1.95	0.48
8:C:295:HIS:O	8:C:299:LEU:N	2.29	0.48
9:D:426:LYS:O	9:D:429:ARG:N	2.47	0.48
9:F:416:GLU:O	9:F:420:THR:N	2.44	0.48
8:C:38:GLY:O	8:C:76:LEU:N	2.45	0.48
9:E:20:LEU:HA	9:E:91:VAL:O	2.13	0.48
8:A:145:GLU:O	8:A:162:ARG:N	2.40	0.48
9:F:381:GLY:O	9:F:385:TYR:N	2.35	0.48
9:D:152:GLY:O	9:D:449:VAL:N	2.46	0.48
8:C:213:PHE:O	8:C:217:GLY:N	2.46	0.48
9:E:169:GLY:HA2	9:E:325:GLN:O	2.14	0.48
8:B:152:ILE:O	8:B:156:ALA:CB	2.62	0.48
8:B:109:ILE:HA	8:B:115:PRO:HA	1.95	0.47
8:B:349:LEU:O	8:B:353:GLY:N	2.47	0.47
7:O:26:VAL:HA	7:N:27:GLY:HA3	1.95	0.47
8:A:10:SER:O	8:A:14:ARG:N	2.37	0.47
9:E:203:GLY:N	9:E:237:GLY:O	2.46	0.47
8:B:54:VAL:HA	8:B:92:THR:H	1.80	0.47
7:Y:26:VAL:HA	7:X:27:GLY:HA3	1.96	0.47
8:C:29:THR:HA	8:C:89:VAL:O	2.15	0.47
8:A:405:LYS:O	8:A:409:ASN:CB	2.63	0.47
9:F:459:GLN:O	9:F:463:SER:N	2.42	0.47
8:B:50:ALA:O	9:D:84:GLY:N	2.48	0.47
8:B:483:THR:O	8:B:487:GLU:N	2.30	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:J:214:ILE:O	4:J:218:VAL:CB	2.63	0.47
8:A:306:SER:O	8:A:311:GLU:N	2.36	0.47
9:D:70:GLY:N	9:D:73:ARG:O	2.43	0.47
9:F:47:LEU:HA	9:F:94:THR:H	1.80	0.47
9:D:69:LEU:H	9:D:74:VAL:HA	1.79	0.47
8:A:40:ALA:O	8:A:73:GLY:HA2	2.15	0.47
8:A:349:LEU:O	8:A:353:GLY:N	2.48	0.47
9:E:21:GLY:CA	9:E:34:ALA:O	2.63	0.47
8:B:171:ASP:O	8:B:174:THR:N	2.44	0.47
2:I:161:ASN:O	2:I:165:HIS:CB	2.64	0.46
9:F:27:ILE:N	9:F:30:VAL:O	2.48	0.46
8:B:404:ASP:O	8:B:408:GLN:N	2.34	0.46
9:D:106:THR:O	9:D:109:ARG:N	2.46	0.46
9:F:221:VAL:O	9:F:231:LYS:N	2.48	0.46
9:F:492:MET:O	9:F:496:LEU:N	2.33	0.46
8:C:349:LEU:HA	8:C:352:ALA:HB3	1.97	0.46
8:C:499:GLU:O	8:C:503:LEU:N	2.44	0.46
8:B:8:GLU:O	8:B:12:ILE:N	2.49	0.46
8:B:50:ALA:HB3	9:D:84:GLY:H	1.81	0.46
9:F:458:PHE:O	9:F:462:LEU:CB	2.65	0.45
9:F:471:GLU:O	9:F:475:TYR:N	2.49	0.45
9:E:186:ILE:O	9:E:190:ALA:CB	2.65	0.45
8:B:100:VAL:O	8:B:126:GLU:CB	2.65	0.45
9:D:46:ALA:O	9:D:94:THR:N	2.47	0.45
8:B:36:GLY:N	8:B:39:ILE:O	2.45	0.45
8:B:136:GLY:O	8:B:140:ARG:N	2.50	0.45
8:A:146:PRO:HA	8:A:161:GLY:HA2	1.99	0.45
8:B:30:GLY:HA2	8:B:43:HIS:O	2.17	0.45
7:W:33:GLY:CA	7:V:31:ALA:HA	2.46	0.45
7:U:12:ALA:HB2	7:V:11:ILE:HA	1.99	0.45
8:C:99:PRO:HA	8:C:127:SER:HA	1.98	0.45
9:D:186:ILE:O	9:D:190:ALA:HB3	2.16	0.45
2:I:148:ALA:HA	4:J:211:SER:HA	1.99	0.45
9:F:201:GLY:O	9:F:237:GLY:N	2.37	0.45
8:B:168:ILE:N	8:B:318:PRO:O	2.41	0.45
8:B:375:ALA:HA	8:B:482:PHE:H	1.81	0.45
9:D:315:SER:HA	9:D:321:ILE:H	1.81	0.45
2:I:42:ILE:O	2:I:46:LYS:CB	2.65	0.44
9:F:205:ARG:O	9:F:209:GLY:N	2.47	0.44
8:A:99:PRO:HA	8:A:127:SER:HA	1.99	0.44
9:E:70:GLY:N	9:E:73:ARG:O	2.48	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:D:204:GLU:H	9:D:238:GLN:HA	1.82	0.44
9:E:189:ILE:HA	9:E:192:ALA:HB3	1.99	0.44
9:D:242:PRO:O	9:D:246:ARG:N	2.46	0.44
9:F:307:GLY:O	9:F:311:GLU:N	2.49	0.44
9:F:396:GLN:O	9:F:400:GLU:CB	2.65	0.44
9:D:215:GLU:O	9:D:219:SER:CB	2.66	0.44
9:E:49:VAL:HA	9:E:91:VAL:HA	1.98	0.44
7:T:27:GLY:HA3	7:U:26:VAL:HA	1.99	0.44
8:C:140:ARG:HA	8:C:305:LEU:HA	2.00	0.44
9:F:48:ILE:HA	9:F:61:VAL:O	2.18	0.44
8:C:481:THR:O	8:C:483:THR:N	2.51	0.44
9:E:110:ILE:O	9:E:118:VAL:N	2.50	0.44
9:E:186:ILE:O	9:E:190:ALA:HB2	2.18	0.44
9:F:295:ALA:O	9:F:297:GLY:N	2.34	0.44
9:E:273:ASP:HA	9:E:274:ASN:HA	1.75	0.44
9:D:22:ARG:O	9:D:34:ALA:N	2.39	0.43
9:D:295:ALA:O	9:D:297:GLY:N	2.34	0.43
7:N:26:VAL:HA	7:M:27:GLY:HA3	2.00	0.43
8:B:98:ILE:CB	8:B:128:ARG:O	2.66	0.43
9:E:223:ASN:H	9:E:230:SER:HA	1.83	0.43
9:D:202:VAL:O	9:D:274:ASN:N	2.45	0.43
7:R:5:ILE:HA	7:S:7:ALA:HB2	2.01	0.43
8:B:109:ILE:N	8:B:224:VAL:O	2.50	0.43
8:A:170:GLY:O	8:A:322:THR:N	2.49	0.43
8:C:174:THR:N	8:C:176:LYS:H	2.17	0.43
9:D:112:ASN:N	9:D:116:GLU:O	2.39	0.43
8:A:147:LEU:N	8:A:160:VAL:O	2.52	0.43
8:B:292:PHE:O	8:B:296:SER:N	2.46	0.43
9:E:47:LEU:HA	9:E:93:ASP:HA	2.01	0.43
9:E:63:CYS:HA	9:E:78:ALA:HA	2.01	0.42
8:B:293:TYR:O	8:B:297:ARG:N	2.37	0.42
9:E:479:ASN:O	9:E:483:ALA:N	2.38	0.42
8:C:283:GLY:N	8:C:287:TYR:O	2.35	0.42
9:F:273:ASP:HA	9:F:274:ASN:HA	1.61	0.42
8:C:67:LEU:N	9:E:26:ILE:O	2.51	0.42
7:N:11:ILE:HA	7:M:12:ALA:HB2	2.01	0.42
8:A:296:SER:O	8:A:300:GLU:CB	2.67	0.42
7:Z:14:GLY:HA3	7:Y:12:ALA:HB1	2.01	0.42
9:F:414:SER:O	9:F:418:ARG:N	2.39	0.42
8:C:173:GLN:O	8:C:175:GLY:N	2.51	0.42
8:B:438:ILE:O	8:B:442:THR:CB	2.67	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:D:49:VAL:HA	9:D:91:VAL:HA	2.01	0.42
2:I:165:HIS:O	2:I:169:ILE:CB	2.68	0.42
8:A:218:ALA:O	8:A:222:THR:N	2.52	0.42
9:F:149:PHE:N	9:F:162:TYR:O	2.50	0.42
8:B:484:GLU:O	8:B:488:ALA:HB2	2.20	0.41
8:C:139:SER:O	8:C:306:SER:N	2.40	0.41
8:C:341:GLY:HA2	8:C:365:SER:HA	2.02	0.41
9:E:294:SER:N	9:E:298:TYR:O	2.53	0.41
8:B:43:HIS:HA	8:B:71:ASN:HA	2.02	0.41
9:D:24:ALA:N	9:D:32:ASP:O	2.53	0.41
9:D:404:ILE:O	9:D:408:LEU:CB	2.69	0.41
9:D:356:LEU:O	9:D:360:GLY:N	2.53	0.41
7:O:11:ILE:HA	7:N:12:ALA:HB2	2.02	0.41
9:D:270:LEU:O	9:D:324:ILE:N	2.53	0.41
8:A:330:TYR:O	8:A:334:ASN:CB	2.68	0.41
8:C:294:LEU:O	8:C:298:LEU:N	2.47	0.41
8:A:202:LYS:O	8:A:206:VAL:N	2.48	0.41
9:D:258:GLU:O	9:D:262:ASP:CB	2.69	0.41
9:D:436:PHE:N	9:D:446:GLY:HA3	2.35	0.41
9:F:49:VAL:O	9:F:60:ASN:HA	2.21	0.41
8:C:430:THR:O	8:C:434:GLN:N	2.45	0.40
8:B:474:GLU:O	8:B:478:SER:CB	2.69	0.40
9:D:286:SER:O	9:D:290:GLY:N	2.54	0.40
7:T:12:ALA:HB1	7:U:14:GLY:HA3	2.02	0.40
9:F:49:VAL:HA	9:F:91:VAL:HA	2.03	0.40
9:D:226:ASN:O	9:D:230:SER:N	2.55	0.40

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	220/247 (89%)	194 (88%)	26 (12%)	0	100	100
2	I	146/184 (79%)	143 (98%)	3 (2%)	0	100	100
3	d	175/257 (68%)	158 (90%)	17 (10%)	0	100	100
4	J	128/222 (58%)	126 (98%)	2 (2%)	0	100	100
5	e	127/134 (95%)	109 (86%)	18 (14%)	0	100	100
6	g	319/364 (88%)	294 (92%)	24 (8%)	1 (0%)	41	76
7	M	77/81 (95%)	70 (91%)	7 (9%)	0	100	100
7	N	77/81 (95%)	70 (91%)	7 (9%)	0	100	100
7	O	77/81 (95%)	70 (91%)	7 (9%)	0	100	100
7	P	77/81 (95%)	70 (91%)	7 (9%)	0	100	100
7	Q	77/81 (95%)	70 (91%)	7 (9%)	0	100	100
7	R	77/81 (95%)	70 (91%)	7 (9%)	0	100	100
7	S	77/81 (95%)	70 (91%)	7 (9%)	0	100	100
7	T	77/81 (95%)	70 (91%)	7 (9%)	0	100	100
7	U	77/81 (95%)	70 (91%)	7 (9%)	0	100	100
7	V	77/81 (95%)	70 (91%)	7 (9%)	0	100	100
7	W	77/81 (95%)	70 (91%)	7 (9%)	0	100	100
7	X	77/81 (95%)	70 (91%)	7 (9%)	0	100	100
7	Y	77/81 (95%)	70 (91%)	7 (9%)	0	100	100
7	Z	77/81 (95%)	70 (91%)	7 (9%)	0	100	100
8	A	493/507 (97%)	449 (91%)	44 (9%)	0	100	100
8	B	502/507 (99%)	456 (91%)	45 (9%)	1 (0%)	47	81
8	C	499/507 (98%)	443 (89%)	56 (11%)	0	100	100
9	D	477/498 (96%)	416 (87%)	60 (13%)	1 (0%)	47	81
9	E	476/498 (96%)	422 (89%)	54 (11%)	0	100	100
9	F	477/498 (96%)	431 (90%)	44 (9%)	2 (0%)	34	72
All	All	5117/5557 (92%)	4621 (90%)	491 (10%)	5 (0%)	54	85

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
6	g	261	GLU
9	F	296	VAL
9	D	296	VAL

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
9	F	295	ALA
8	B	263	ASP

### 5.3.2 Protein sidechains [i](#)

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

### 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](#)

There are no ligands in this entry.

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
4	J	1
8	B	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	J	219:LEU	C	220:PRO	N	3.44
1	B	144:TYR	C	145:GLU	N	1.18

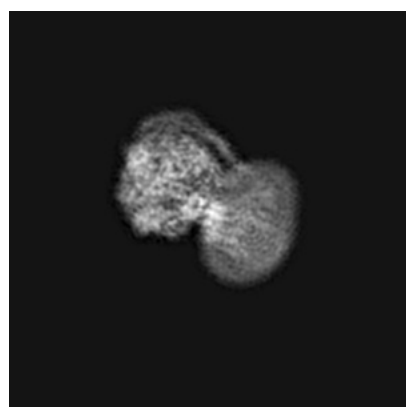
## 6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-21241. These allow visual inspection of the internal detail of the map and identification of artifacts.

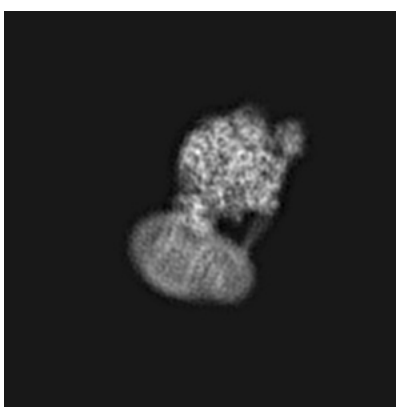
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

### 6.1 Orthogonal projections [i](#)

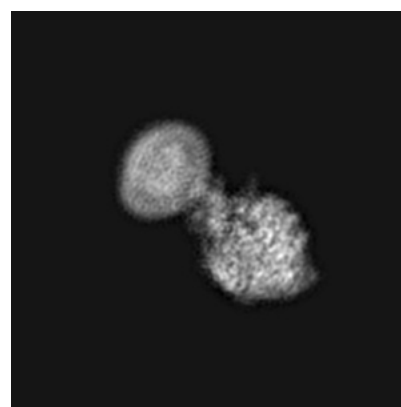
#### 6.1.1 Primary map



X



Y

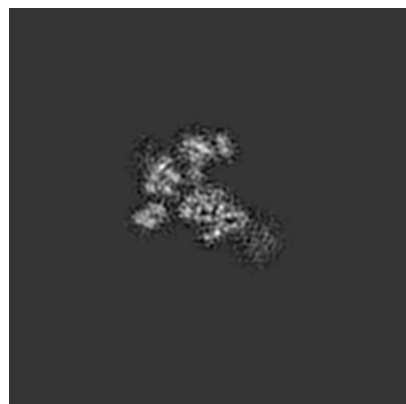


Z

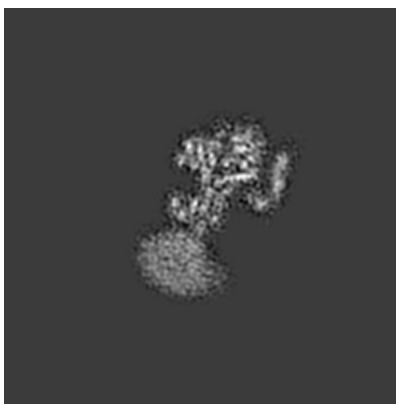
The images above show the map projected in three orthogonal directions.

### 6.2 Central slices [i](#)

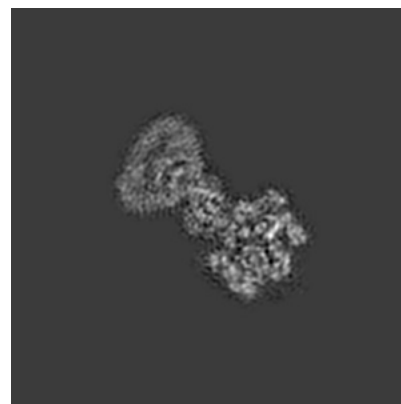
#### 6.2.1 Primary map



X Index: 90



Y Index: 90

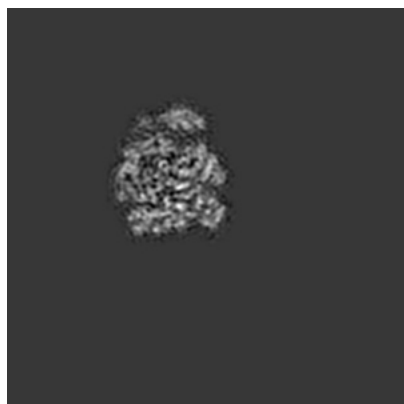


Z Index: 90

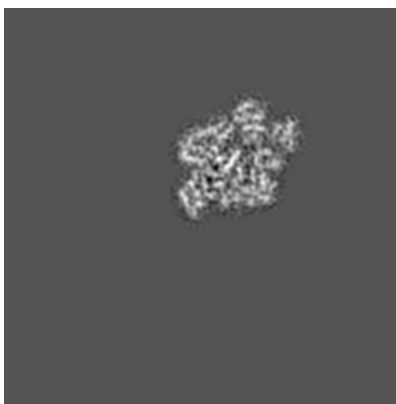
The images above show central slices of the map in three orthogonal directions.

## 6.3 Largest variance slices [i](#)

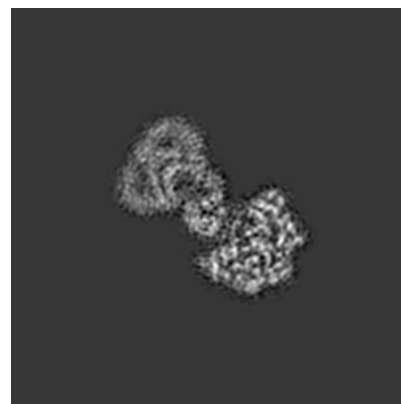
### 6.3.1 Primary map



X Index: 116



Y Index: 61

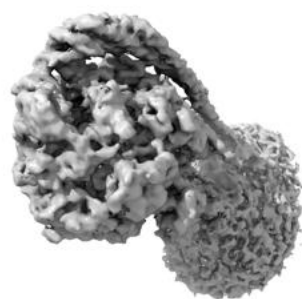


Z Index: 88

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

## 6.4 Orthogonal surface views [i](#)

### 6.4.1 Primary map



X



Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.035. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

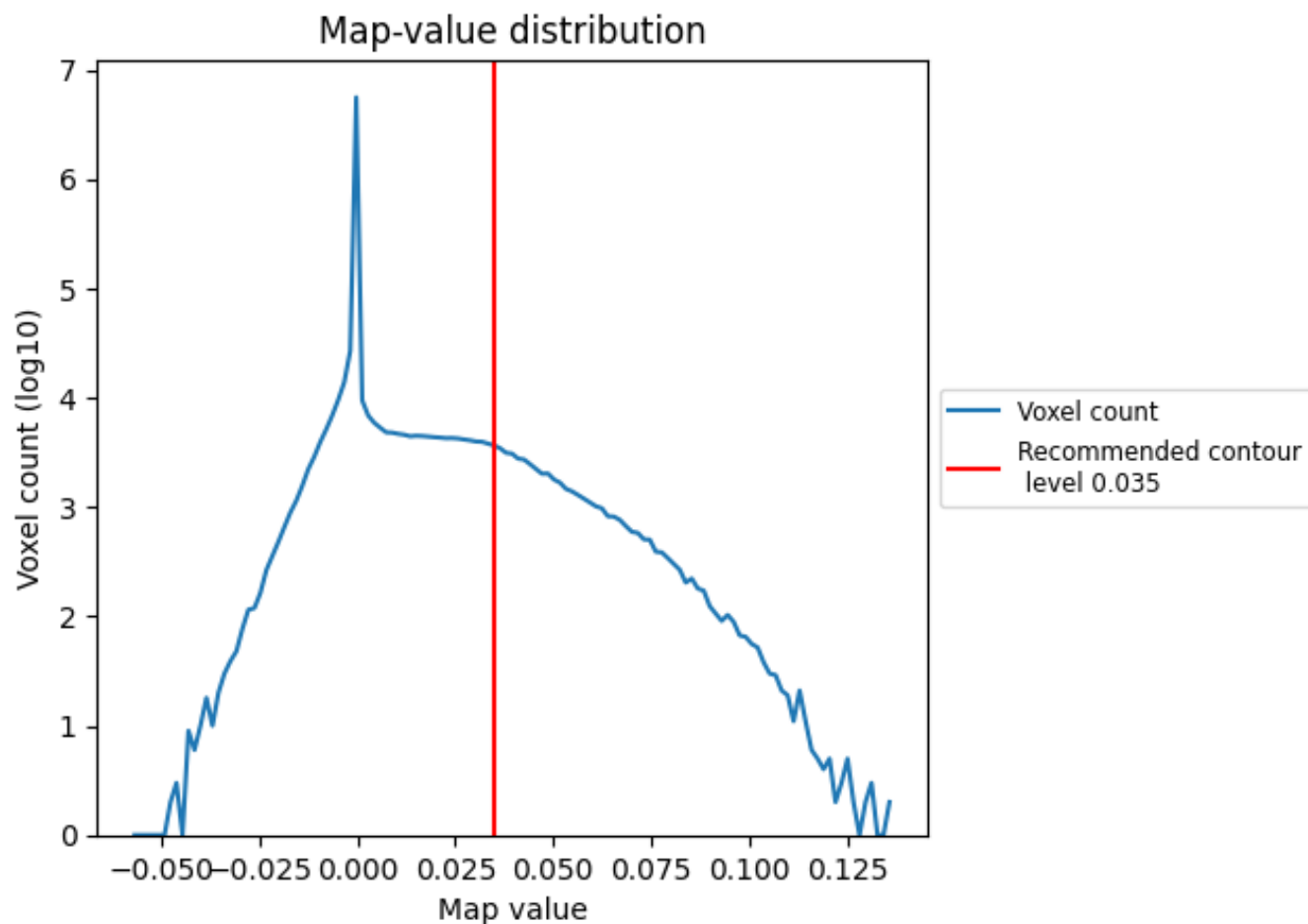
## 6.5 Mask visualisation

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

## 7 Map analysis [i](#)

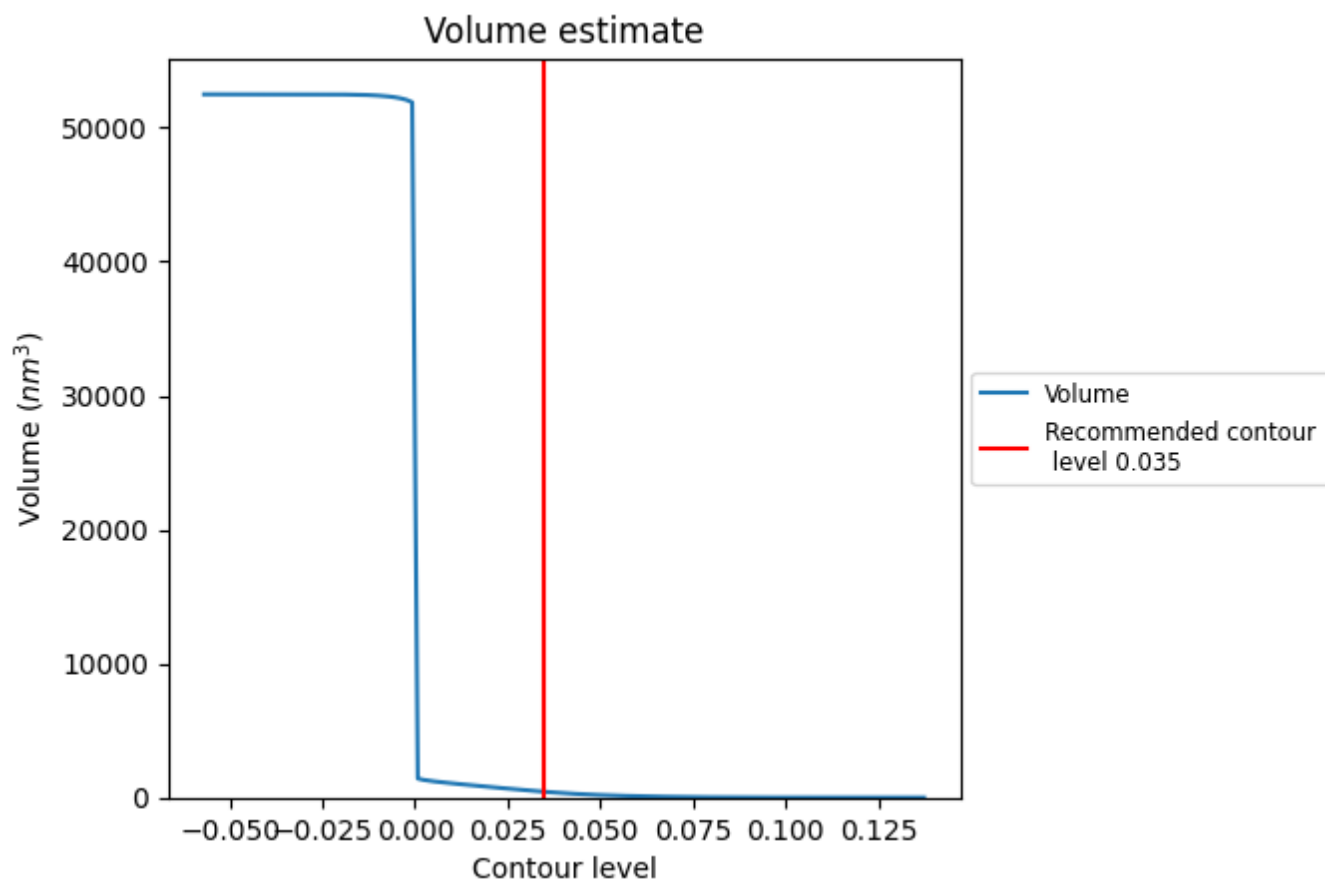
This section contains the results of statistical analysis of the map.

### 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](#)

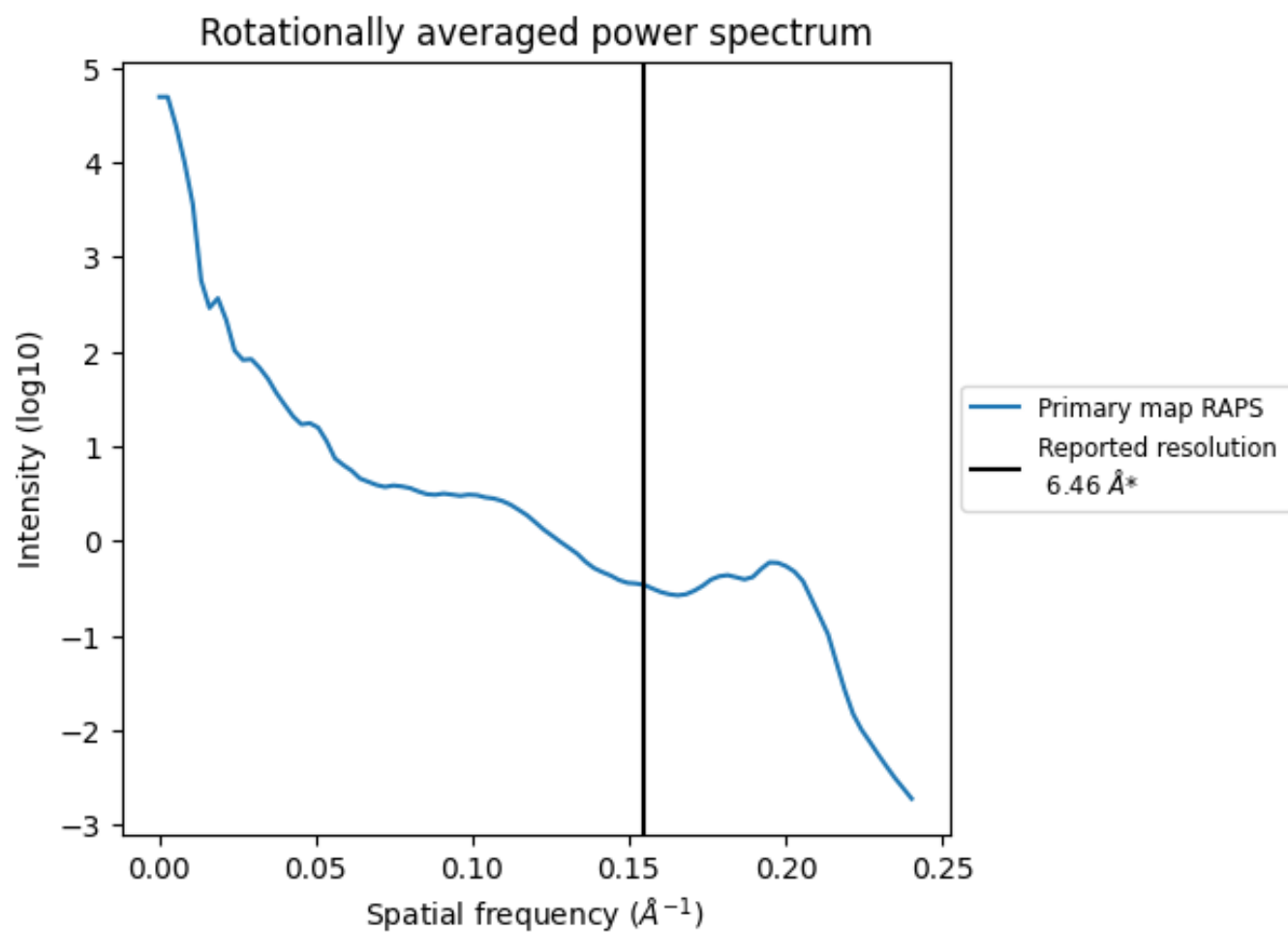


The volume at the recommended contour level is 432 nm<sup>3</sup>; this corresponds to an approximate mass of 390 kDa.

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 ⓘ

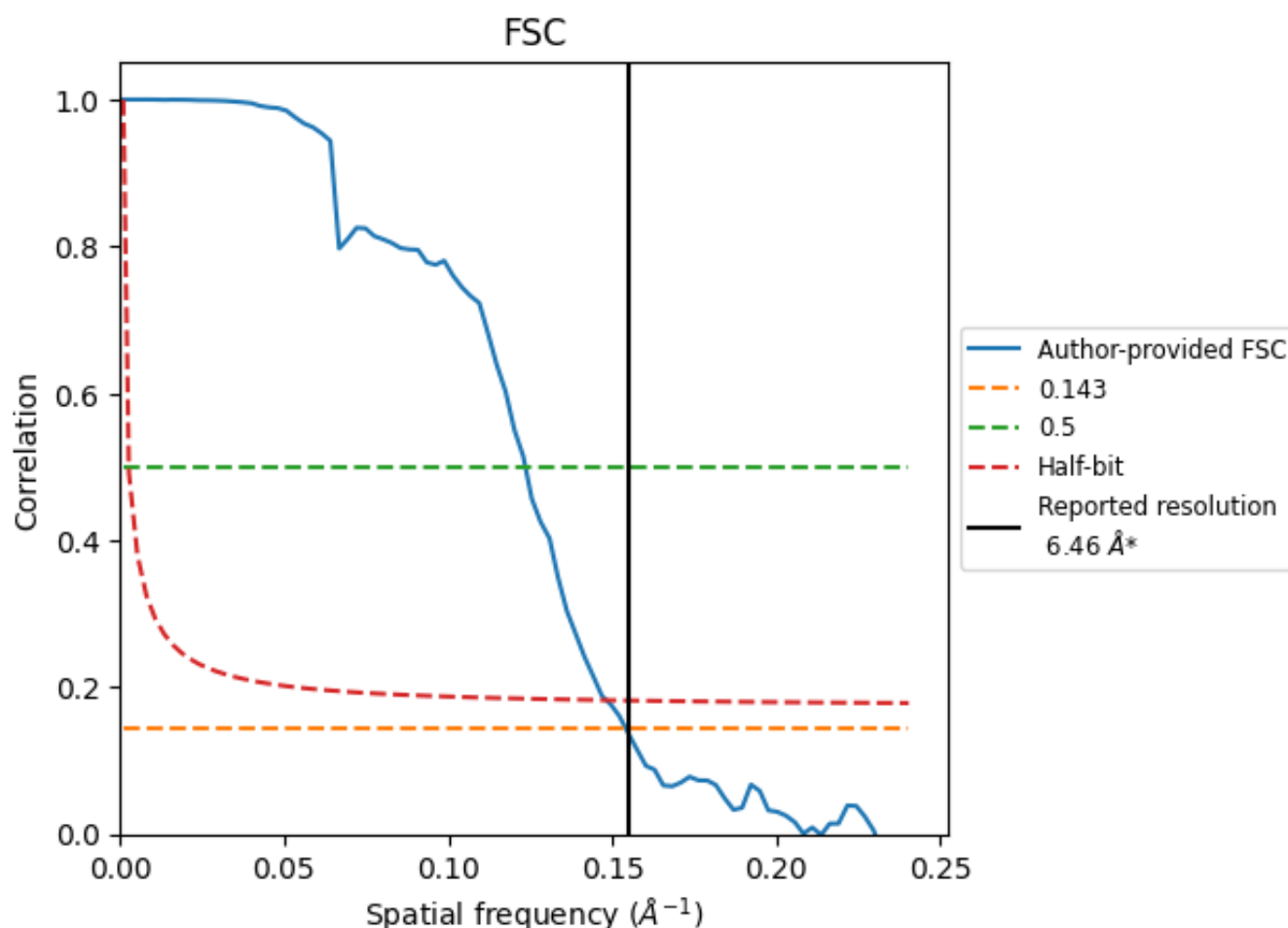


\*Reported resolution corresponds to spatial frequency of 0.155 Å<sup>-1</sup>

## 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 0.155  $\text{\AA}^{-1}$

## 8.2 Resolution estimates [i](#)

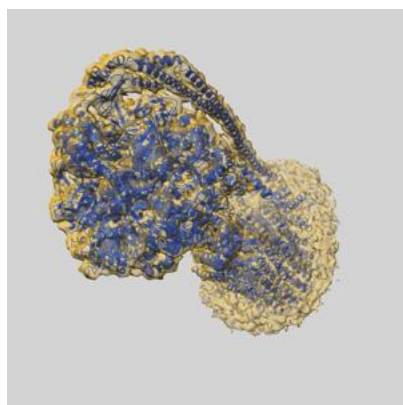
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	6.46	-	-
Author-provided FSC curve	6.49	8.10	6.75
Unmasked-calculated*	-	-	-

\*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps.

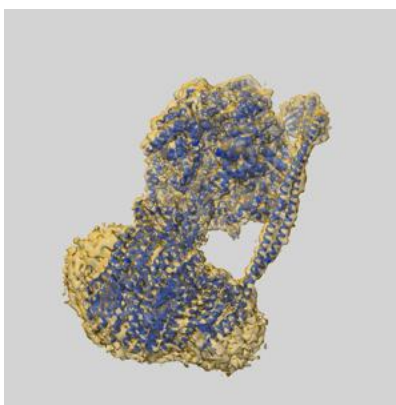
## 9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-21241 and PDB model 6VMG. Per-residue inclusion information can be found in section [3](#) on page [7](#).

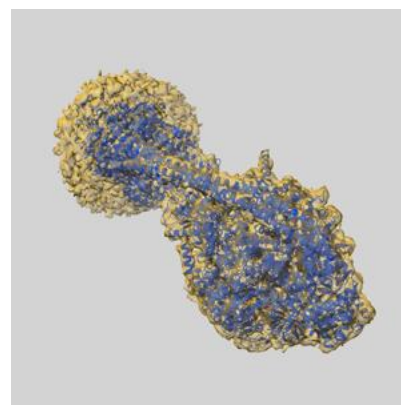
### 9.1 Map-model overlay [i](#)



X



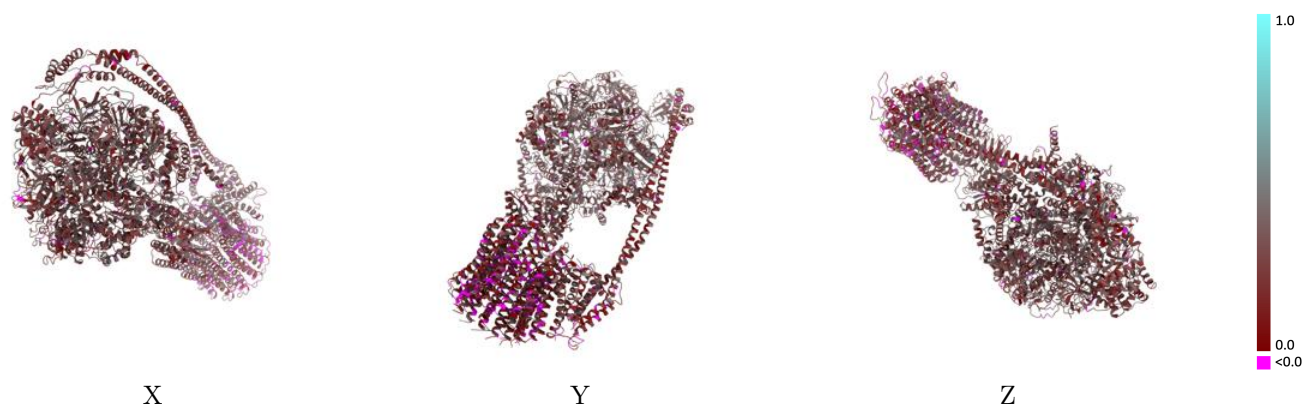
Y



Z

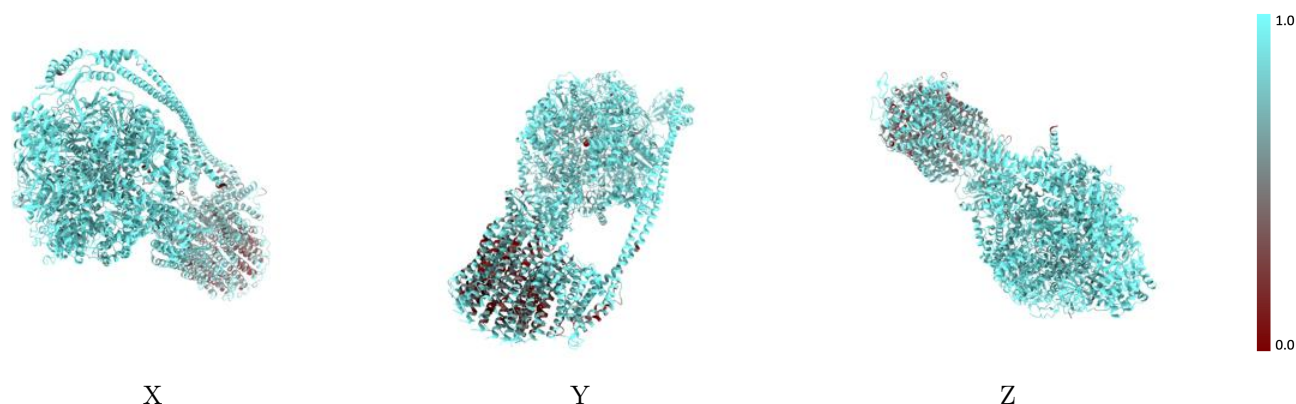
The images above show the 3D surface view of the map at the recommended contour level 0.035 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



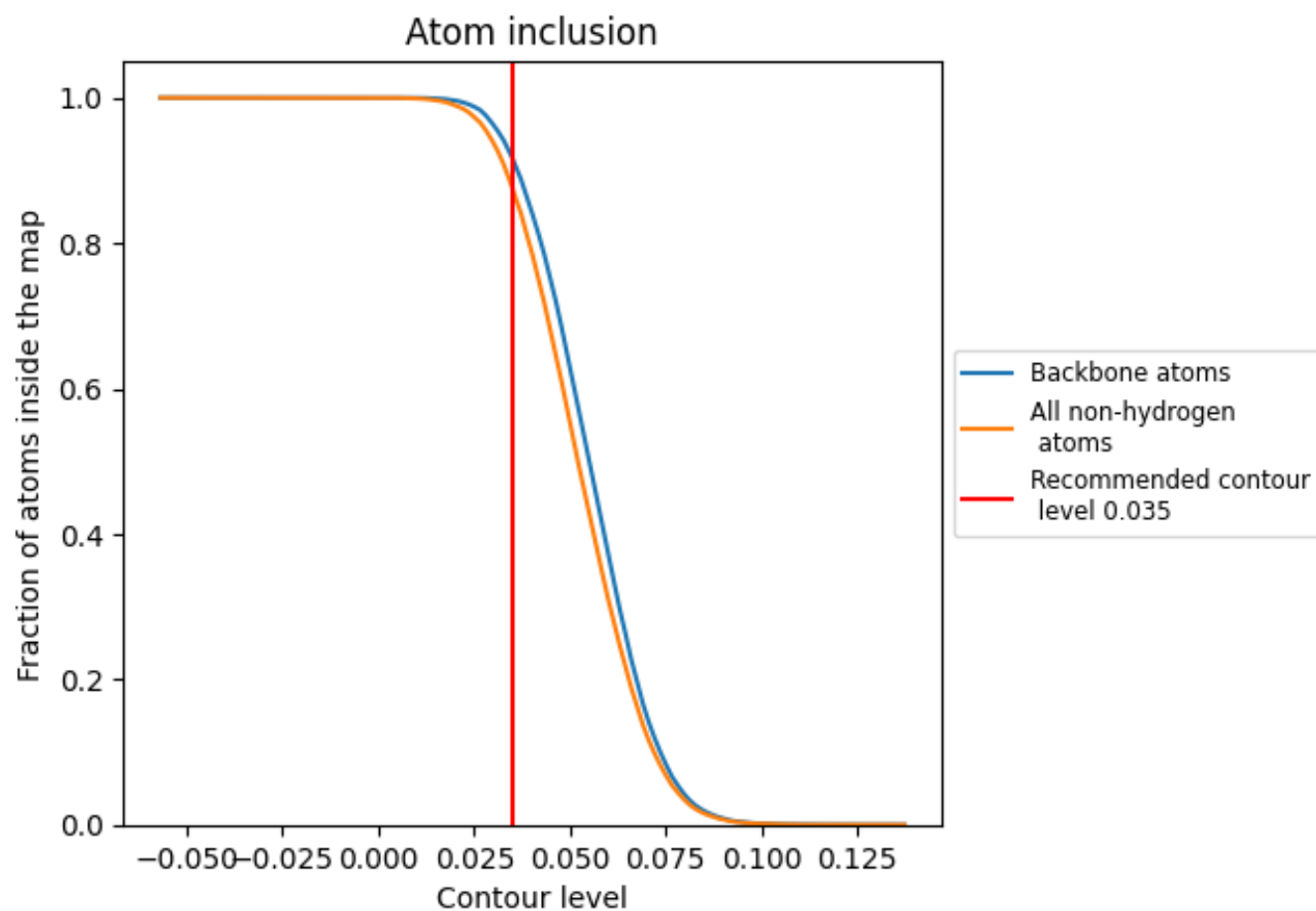
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.035).























































## 9.4 Atom inclusion [i](#)



At the recommended contour level, 92% of all backbone atoms, 88% 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 (0.035) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8761	 0.2620
A	 0.9479	 0.3030
B	 0.9480	 0.2930
C	 0.9432	 0.3040
D	 0.9395	 0.2920
E	 0.9373	 0.3080
F	 0.9302	 0.3010
I	 0.9083	 0.2250
J	 0.9078	 0.2120
M	 0.6510	 0.1710
N	 0.6458	 0.1670
O	 0.7266	 0.1740
P	 0.7083	 0.1590
Q	 0.6406	 0.1600
R	 0.5885	 0.1400
S	 0.6042	 0.1380
T	 0.7344	 0.1850
U	 0.7266	 0.1800
V	 0.7604	 0.1780
W	 0.7135	 0.1720
X	 0.6562	 0.1410
Y	 0.7292	 0.1670
Z	 0.6901	 0.1510
a	 0.8323	 0.2210
d	 0.9715	 0.2800
e	 0.8383	 0.2660
g	 0.9000	 0.2960

