



Full wwPDB EM Validation Report ⓘ

Mar 8, 2026 – 12:06 PM UTC

PDB ID : 8WUT / pdb_00008wut
EMDB ID : EMD-37859
Title : SpCas9-MMLV RT-pegRNA-target DNA complex (initiation)
Authors : Shuto, Y.; Nakagawa, R.; Hoki, M.; Omura, S.N.; Hirano, H.; Itoh, Y.; Nureki, O.
Deposited on : 2023-10-21
Resolution : 3.00 Å(reported)

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

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

A user guide is available at

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

The types of validation reports are described at

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

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

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

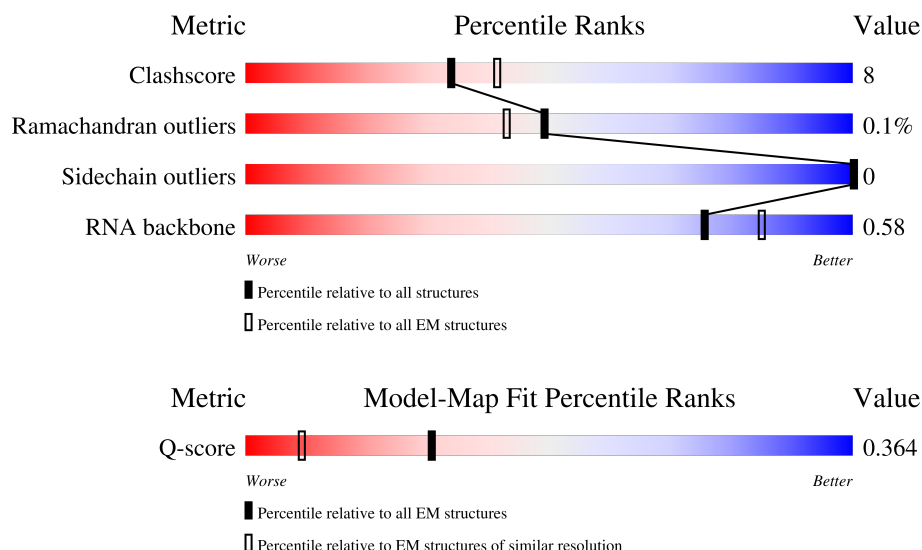
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.00 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
RNA backbone	8273	3508	-
Q-score	-	25397	14081 (2.50 - 3.50)

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	A	1367	
2	E	496	
3	C	51	

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Mol	Chain	Length	Quality of chain
4	D	17	65% 24% 12%
5	B	115	49% 41% 8%
6	F	35	17% 29% 34% 37%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 18316 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 CRISPR-associated endonuclease Cas9/Csn1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	1313	Total	C	N	O	S	0	0
			10751	6854	1868	2008	21		

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	10	ALA	ASP	engineered mutation	UNP Q99ZW2
A	840	ALA	HIS	engineered mutation	UNP Q99ZW2

- Molecule 2 is a protein called Gag-Pol polyprotein.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	E	453	Total	C	N	O	S	0	0
			3571	2313	619	625	14		

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	200	ASN	ASP	engineered mutation	UNP Q8UN00
E	249	ASN	ASP	engineered mutation	UNP Q8UN00
E	306	LYS	THR	engineered mutation	UNP Q8UN00
E	313	PHE	TRP	engineered mutation	UNP Q8UN00
E	330	PRO	THR	engineered mutation	UNP Q8UN00
E	335	ASN	ASP	engineered mutation	UNP Q8UN00

- Molecule 3 is a DNA chain called DNA (51-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	41	Total	C	N	O	P	0	0
			835	397	152	245	41		

- Molecule 4 is a DNA chain called DNA (5'-D(*TP*GP*AP*TP*GP*GP*CP*AP*GP*AP*GP*TP*AP*CP*TP*AP*G)-3').

Mol	Chain	Residues	Atoms					AltConf	Trace
4	D	15	Total	C	N	O	P	0	0
			313	148	62	88	15		

- Molecule 5 is a RNA chain called RNA (115-MER).

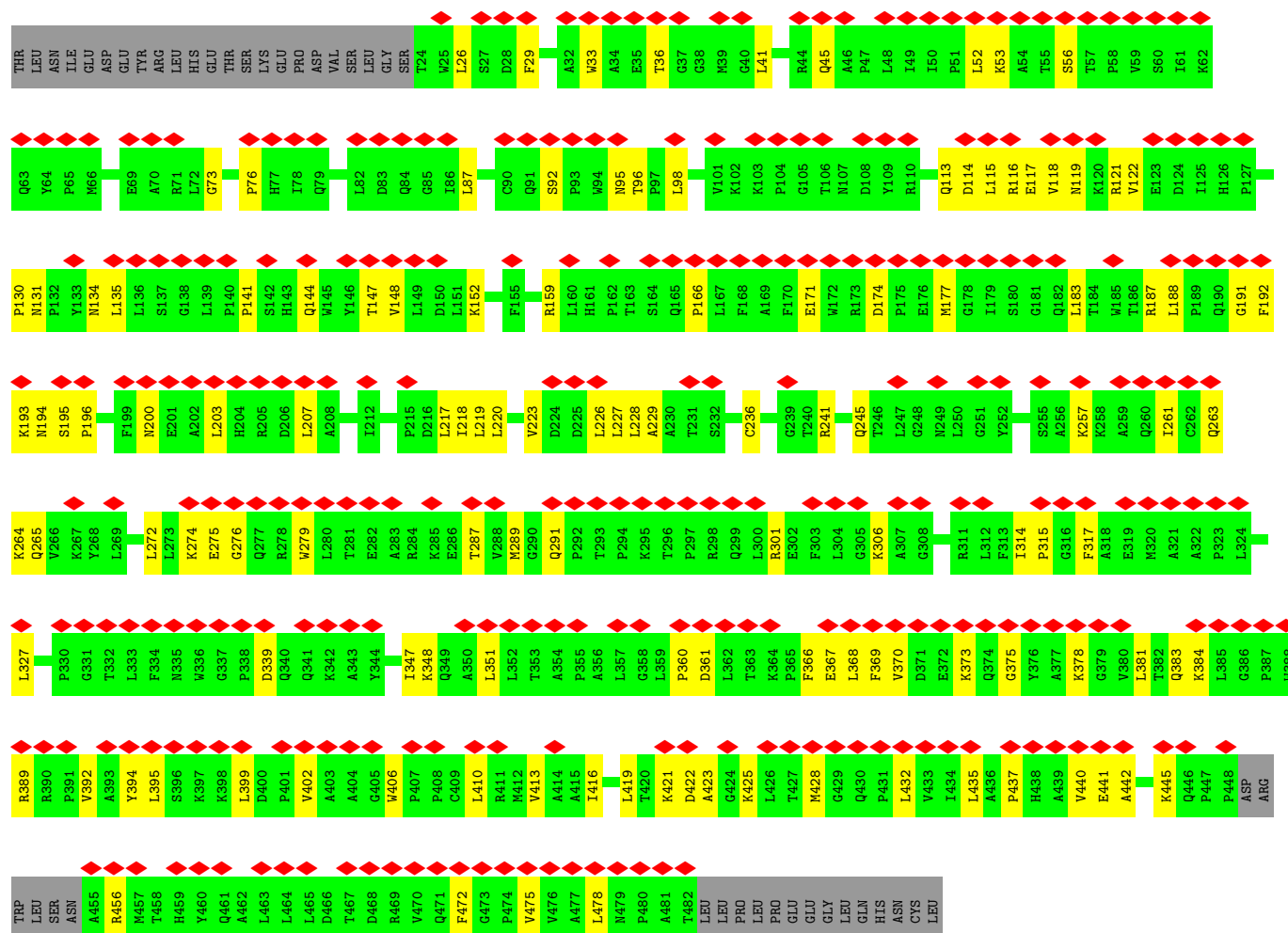
Mol	Chain	Residues	Atoms					AltConf	Trace
5	B	112	Total	C	N	O	P	0	0
			2393	1069	428	784	112		

- Molecule 6 is DNA/RNA hybrid called DNA/RNA (35-MER).

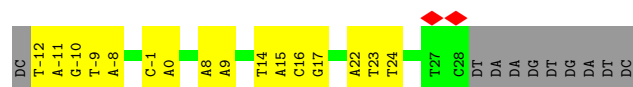
Mol	Chain	Residues	Atoms					AltConf	Trace
6	F	22	Total	C	N	O	P	0	0
			453	214	92	125	22		



• Molecule 2: Gag-Pol polyprotein

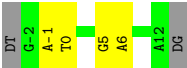


• Molecule 3: DNA (51-MER)



• Molecule 4: DNA (5'-D(*TP*GP*AP*TP*GP*GP*CP*AP*GP*AP*GP*TP*AP*CP*TP*AP*G)-3')

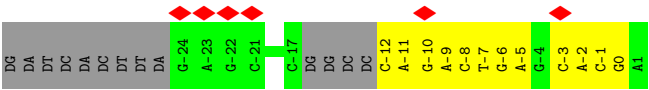
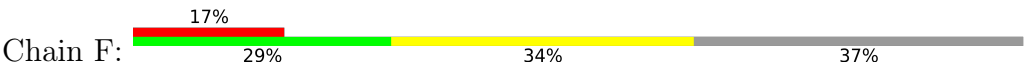




• Molecule 5: RNA (115-MER)



• Molecule 6: DNA/RNA (35-MER)



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	118125	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)	1600	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	2.032	Depositor
Minimum map value	-0.227	Depositor
Average map value	0.003	Depositor
Map value standard deviation	0.084	Depositor
Recommended contour level	0.393	Depositor
Map size ($\text{\AA}$)	223.0668, 223.0668, 223.0668	wwPDB
Map dimensions	172, 172, 172	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	1.2969, 1.2969, 1.2969	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 2DA

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	0.23	0/10940	0.30	0/14696
2	E	0.12	0/3671	0.29	0/5007
3	C	0.34	0/935	0.42	0/1439
4	D	0.31	0/352	0.36	0/542
5	B	0.30	0/2676	0.28	0/4167
6	F	0.18	0/486	0.37	0/745
All	All	0.23	0/19060	0.31	0/26596

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 [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	10751	0	10946	132	0
2	E	3571	0	3652	94	0
3	C	835	0	461	10	0
4	D	313	0	169	2	0
5	B	2393	0	1204	44	0
6	F	453	0	246	16	0
All	All	18316	0	16678	285	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 8.

All (285) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:370:VAL:HG13	2:E:419:LEU:HD12	1.62	0.79
2:E:272:LEU:HB3	2:E:279:TRP:HB2	1.65	0.78
1:A:212:LEU:HD13	1:A:300:ILE:HD11	1.65	0.77
2:E:218:ILE:HB	2:E:229:ALA:HB3	1.67	0.76
3:C:16:DC:H2''	3:C:17:DG:H5'	1.68	0.76
2:E:117:GLU:HB3	2:E:121:ARG:HH22	1.51	0.75
1:A:44:LYS:NZ	5:B:92:G:O6	2.22	0.72
2:E:301:ARG:HD3	6:F:-3:DC:H5'	1.72	0.71
2:E:301:ARG:HB3	6:F:-3:DC:H4'	1.71	0.71
3:C:15:DA:H2''	3:C:16:DC:H5''	1.72	0.71
2:E:147:THR:OG1	2:E:236:CYS:SG	2.50	0.69
2:E:203:LEU:HD21	2:E:226:LEU:HD21	1.72	0.69
2:E:366:PHE:HB2	2:E:432:LEU:HB2	1.75	0.69
3:C:-1:DC:H2''	3:C:0:DA:C8	2.29	0.67
1:A:171:GLU:HG2	1:A:269:ASP:HB3	1.76	0.67
1:A:802:GLU:HG2	1:A:805:GLN:H	1.60	0.67
2:E:413:VAL:HG22	2:E:440:VAL:HG21	1.77	0.67
1:A:1151:LYS:NZ	3:C:-8:DA:OP1	2.28	0.66
2:E:141:PRO:O	2:E:144:GLN:NE2	2.29	0.64
5:B:109:C:H2'	5:B:110:A:H8	1.63	0.63
6:F:-12:DC:H2'	6:F:-11:DA:C8	2.33	0.63
2:E:317:PHE:HE1	2:E:347:ILE:HG12	1.64	0.63
1:A:9:LEU:HD12	1:A:761:ILE:HG22	1.81	0.62
1:A:49:GLY:HA2	1:A:1092:VAL:HG13	1.81	0.62
2:E:131:ASN:ND2	2:E:134:ASN:OD1	2.27	0.62
1:A:499:ASP:HB2	1:A:663:SER:HB3	1.80	0.62
2:E:193:LYS:HA	5:B:102:U:O2'	2.00	0.62
1:A:554:LYS:NZ	1:A:608:ASP:OD1	2.30	0.61
5:B:3:C:O2'	5:B:4:C:O5'	2.19	0.61
5:B:110:A:H2'	5:B:111:G:C8	2.36	0.61
2:E:130:PRO:HD3	2:E:200:ASN:HD21	1.66	0.60
5:B:52:A:OP2	5:B:62:G:N2	2.32	0.60
6:F:-12:DC:H2'	6:F:-11:DA:H8	1.65	0.60
5:B:110:A:H2'	5:B:111:G:H8	1.67	0.60
1:A:995:THR:HG22	1:A:999:LYS:HD2	1.82	0.60
1:A:979:ASN:ND2	1:A:1226:LEU:O	2.34	0.60
6:F:-3:DC:H2'	6:F:-2:DA:C8	2.36	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:606:PHE:O	1:A:612:ASN:ND2	2.35	0.59
6:F:-11:DA:H2'	6:F:-10:DG:C8	2.36	0.59
1:A:95:ASP:OD1	1:A:95:ASP:N	2.35	0.59
2:E:402:VAL:HG11	6:F:-12:DC:H5''	1.84	0.59
2:E:29:PHE:HE1	2:E:241:ARG:HH11	1.50	0.59
2:E:367:GLU:OE1	2:E:389:ARG:NH1	2.36	0.58
1:A:1193:ASP:OD1	1:A:1193:ASP:N	2.37	0.58
2:E:92:SER:HB2	2:E:166:PRO:HB3	1.85	0.57
2:E:369:PHE:HB3	2:E:478:LEU:HD22	1.85	0.57
2:E:122:VAL:HG21	2:E:192:PHE:HZ	1.69	0.57
2:E:265:GLN:HE21	2:E:272:LEU:HD21	1.70	0.57
1:A:278:LEU:O	1:A:282:ILE:HG22	2.05	0.57
2:E:148:VAL:HG22	2:E:227:LEU:HD13	1.87	0.56
1:A:1050:ILE:HG22	1:A:1051:THR:HG23	1.87	0.56
6:F:-9:DA:H2'	6:F:-8:DC:C6	2.41	0.56
2:E:265:GLN:HG3	2:E:272:LEU:HD11	1.88	0.56
1:A:512:SER:OG	1:A:617:GLU:OE1	2.22	0.56
1:A:791:LEU:HD13	1:A:889:ALA:HB2	1.88	0.55
3:C:8:DA:H2'	3:C:9:DA:C8	2.41	0.55
5:B:111:G:H2'	5:B:112:U:C6	2.42	0.55
5:B:108:U:H2'	5:B:109:C:C6	2.41	0.55
1:A:8:GLY:HA3	1:A:991:ALA:HB2	1.89	0.55
1:A:801:VAL:HG23	1:A:805:GLN:HB3	1.88	0.55
1:A:870:VAL:HG21	1:A:902:LYS:HB3	1.89	0.55
1:A:525:THR:O	1:A:690:ASN:ND2	2.39	0.55
1:A:686:ASP:HB3	1:A:690:ASN:HA	1.89	0.54
2:E:442:ALA:HA	2:E:445:LYS:HD2	1.89	0.54
2:E:122:VAL:HG21	2:E:192:PHE:CZ	2.42	0.54
2:E:456:ARG:HB2	6:F:-11:DA:H4'	1.88	0.54
5:B:105:U:H2'	5:B:106:G:H8	1.73	0.54
6:F:-11:DA:H2'	6:F:-10:DG:H8	1.73	0.53
1:A:232:GLU:HG3	1:A:238:PHE:HB3	1.90	0.53
2:E:207:LEU:HD13	2:E:219:LEU:HD21	1.89	0.53
1:A:827:GLU:O	1:A:859:ARG:NH2	2.42	0.53
2:E:315:PRO:HB3	2:E:394:TYR:CD1	2.44	0.53
5:B:108:U:H2'	5:B:109:C:H6	1.73	0.53
1:A:213:SER:O	1:A:215:ARG:NH2	2.37	0.53
2:E:289:MET:O	2:E:348:LYS:NZ	2.42	0.53
1:A:536:LYS:NZ	1:A:537:PRO:O	2.36	0.53
2:E:192:PHE:HB3	2:E:195:SER:OG	2.09	0.53
5:B:107:C:H2'	5:B:108:U:C6	2.44	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:565:LYS:HA	1:A:569:PHE:HB2	1.91	0.52
2:E:217:LEU:HB3	2:E:228:LEU:HD11	1.89	0.52
2:E:314:ILE:HB	2:E:317:PHE:HB2	1.92	0.52
5:B:109:C:H2'	5:B:110:A:C8	2.43	0.52
2:E:26:LEU:HA	2:E:33:TRP:CD1	2.44	0.52
2:E:87:LEU:HD13	2:E:183:LEU:HD13	1.92	0.52
2:E:147:THR:HG23	2:E:261:ILE:HA	1.92	0.52
1:A:521:TYR:CE2	1:A:684:LYS:HG2	2.45	0.52
2:E:361:ASP:O	2:E:383:GLN:NE2	2.38	0.52
2:E:95:ASN:ND2	2:E:171:GLU:OE1	2.43	0.51
2:E:406:TRP:HB3	2:E:410:LEU:HD21	1.92	0.51
2:E:131:ASN:O	2:E:135:LEU:HD12	2.11	0.51
5:B:112:U:H2'	5:B:113:C:C6	2.45	0.51
1:A:473:ILE:HG12	1:A:481:VAL:HG11	1.92	0.51
1:A:847:LEU:HD13	1:A:919:ARG:HH21	1.76	0.51
2:E:29:PHE:HE2	2:E:245:GLN:HB2	1.75	0.51
6:F:-8:DC:H2'	6:F:-7:DT:C6	2.45	0.51
1:A:539:PHE:HB3	1:A:690:ASN:ND2	2.26	0.51
3:C:14:DT:H2''	3:C:15:DA:H5'	1.94	0.50
5:B:12:U:H2'	5:B:13:U:C6	2.46	0.50
1:A:508:LEU:HD21	1:A:664:ARG:HB2	1.93	0.50
2:E:395:LEU:HD13	2:E:422:ASP:HB2	1.92	0.50
2:E:115:LEU:O	2:E:119:ASN:HB2	2.10	0.50
5:B:111:G:H2'	5:B:112:U:H6	1.76	0.50
1:A:1110:ILE:HG12	1:A:1122:ARG:HD3	1.92	0.50
2:E:144:GLN:HE22	2:E:276:GLY:HA3	1.77	0.50
5:B:22:U:H2'	5:B:23:U:C6	2.46	0.50
2:E:366:PHE:HB2	2:E:432:LEU:CB	2.41	0.50
5:B:5:G:H2'	5:B:6:U:H6	1.76	0.50
5:B:74:A:H2'	5:B:75:A:O4'	2.11	0.50
1:A:161:MET:HE1	1:A:422:ILE:HD12	1.94	0.49
1:A:671:ARG:HG2	1:A:678:THR:HG22	1.93	0.49
3:C:-10:DG:H2''	3:C:-9:DT:C6	2.47	0.49
5:B:103:C:H2'	5:B:104:G:C8	2.47	0.49
2:E:366:PHE:HB3	2:E:381:LEU:HD11	1.94	0.49
1:A:1272:GLN:NE2	5:B:89:G:O6	2.43	0.49
1:A:105:PHE:HE2	5:B:23:U:O2	1.95	0.49
1:A:82:LEU:HD22	1:A:162:ILE:HD12	1.95	0.49
1:A:103:GLU:HB2	1:A:106:LEU:HD12	1.95	0.49
2:E:26:LEU:HD21	2:E:36:THR:HG21	1.96	0.48
1:A:1266:LEU:HD22	1:A:1309:ILE:HD12	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:95:ASN:HD21	2:E:171:GLU:HB2	1.78	0.48
2:E:114:ASP:OD1	2:E:116:ARG:NE	2.47	0.48
1:A:1004:LEU:HD23	1:A:1008:PHE:HE2	1.78	0.48
5:B:84:A:H2'	5:B:85:C:C6	2.49	0.48
1:A:75:ARG:NH2	5:B:50:U:O4	2.42	0.48
2:E:192:PHE:CE2	2:E:194:ASN:HB2	2.48	0.48
1:A:1156:LYS:HE2	1:A:1158:LYS:HG2	1.95	0.47
5:B:75:A:H2'	5:B:76:A:H8	1.79	0.47
1:A:628:ASP:HB3	1:A:631:MET:HG3	1.97	0.47
1:A:1006:SER:OG	1:A:1007:GLU:OE1	2.30	0.47
1:A:631:MET:O	1:A:634:GLU:HG3	2.14	0.47
1:A:65:LYS:O	1:A:69:ARG:HG3	2.15	0.47
2:E:257:LYS:HD2	2:E:257:LYS:HA	1.71	0.47
5:B:5:G:H2'	5:B:6:U:C6	2.49	0.47
5:B:37:U:H2'	5:B:38:A:C8	2.50	0.47
5:B:54:G:H2'	5:B:55:C:C6	2.50	0.47
1:A:860:SER:OG	1:A:863:ASN:OD1	2.31	0.47
2:E:441:GLU:HG3	2:E:472:PHE:CE1	2.49	0.47
1:A:1325:LYS:HE2	1:A:1325:LYS:HB2	1.73	0.47
1:A:864:ARG:NH2	1:A:871:PRO:HD3	2.30	0.47
1:A:908:LEU:O	1:A:913:LYS:NZ	2.47	0.47
2:E:423:ALA:C	2:E:425:LYS:H	2.23	0.46
1:A:832:ARG:NH2	1:A:835:ASP:OD2	2.49	0.46
5:B:8:G:H2'	5:B:9:U:C6	2.50	0.46
5:B:75:A:H2'	5:B:76:A:C8	2.51	0.46
1:A:264:LEU:HD11	1:A:278:LEU:HD23	1.96	0.46
1:A:870:VAL:HG22	1:A:871:PRO:HD2	1.98	0.46
5:B:83:C:H2'	5:B:84:A:H8	1.80	0.46
2:E:45:GLN:O	2:E:159:ARG:NH2	2.49	0.46
1:A:78:ARG:NH1	1:A:162:ILE:O	2.48	0.46
1:A:673:LYS:HB2	1:A:703:THR:HG21	1.98	0.46
2:E:265:GLN:NE2	2:E:272:LEU:HD21	2.31	0.46
2:E:339:ASP:OD1	2:E:339:ASP:N	2.46	0.46
1:A:1215:ALA:HB2	1:A:1221:GLN:HG3	1.98	0.46
5:B:38:A:H2'	5:B:39:G:C8	2.51	0.46
1:A:223:GLU:OE2	1:A:234:LYS:NZ	2.37	0.46
1:A:58:THR:HG22	1:A:731:PRO:HG2	1.99	0.45
1:A:890:LYS:HA	1:A:890:LYS:HD3	1.72	0.45
1:A:1212:ARG:HB3	1:A:1220:LEU:HD22	1.98	0.45
5:B:46:A:H2'	5:B:47:A:C8	2.51	0.45
1:A:273:ASP:O	1:A:653:ARG:NH2	2.50	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:152:LYS:HD2	2:E:152:LYS:HA	1.71	0.45
1:A:268:LYS:HD3	1:A:268:LYS:HA	1.81	0.45
1:A:1124:LYS:HB3	1:A:1124:LYS:HE3	1.62	0.45
2:E:373:LYS:HD3	2:E:378:LYS:HZ3	1.81	0.45
2:E:375:GLY:HA2	2:E:399:LEU:HD12	1.97	0.45
5:B:37:U:H2'	5:B:38:A:H8	1.82	0.45
5:B:76:A:C6	5:B:77:A:H1'	2.52	0.45
1:A:784:ILE:HG21	1:A:806:LEU:HD11	1.98	0.45
2:E:119:ASN:ND2	2:E:191:GLY:O	2.50	0.45
1:A:1050:ILE:HD11	1:A:1060:ARG:HE	1.82	0.45
6:F:-6:DG:H2'	6:F:-5:DA:C8	2.52	0.45
1:A:802:GLU:CD	1:A:805:GLN:HB2	2.42	0.45
2:E:217:LEU:HD13	2:E:228:LEU:HD21	1.98	0.45
2:E:96:THR:HG21	2:E:115:LEU:HD23	1.97	0.44
1:A:270:THR:OG1	1:A:629:ARG:NH1	2.51	0.44
1:A:788:ILE:HG13	1:A:796:LEU:HD12	1.99	0.44
1:A:569:PHE:HA	1:A:573:GLU:HB2	1.99	0.44
2:E:317:PHE:CE1	2:E:347:ILE:HG12	2.47	0.44
2:E:360:PRO:HB3	2:E:392:VAL:HG13	1.99	0.44
1:A:691:ARG:HG3	1:A:696:LEU:HD11	1.98	0.44
1:A:836:TYR:CD1	1:A:859:ARG:HA	2.53	0.44
1:A:545:LYS:NZ	1:A:683:LEU:O	2.50	0.44
4:D:5:DG:H2''	4:D:6:DA:C8	2.53	0.44
1:A:226:ILE:HD13	1:A:226:ILE:HA	1.90	0.44
2:E:41:LEU:HD11	2:E:159:ARG:HD3	1.99	0.44
2:E:121:ARG:NH1	2:E:121:ARG:HB2	2.33	0.44
2:E:134:ASN:CG	2:E:428:MET:HE1	2.43	0.44
2:E:218:ILE:O	2:E:229:ALA:N	2.38	0.44
2:E:41:LEU:HD12	2:E:187:ARG:NH2	2.33	0.43
2:E:306:LYS:N	6:F:-2:DA:H4'	2.33	0.43
1:A:222:LEU:HB2	1:A:246:LEU:HD21	2.00	0.43
2:E:116:ARG:HG3	5:B:101:A:H4'	1.99	0.43
2:E:188:LEU:HD21	2:E:195:SER:HA	2.00	0.43
1:A:1050:ILE:HB	1:A:1058:ARG:HB3	1.99	0.43
2:E:368:LEU:O	2:E:435:LEU:HD23	2.18	0.43
1:A:220:ARG:HD2	1:A:220:ARG:HA	1.84	0.43
1:A:273:ASP:HB3	1:A:653:ARG:NH1	2.34	0.43
1:A:999:LYS:HB3	1:A:1073:VAL:HG12	1.99	0.43
1:A:1085:LYS:HB3	1:A:1085:LYS:HE3	1.75	0.43
2:E:327:LEU:HD23	2:E:327:LEU:HA	1.82	0.43
3:C:-12:DT:H2''	3:C:-11:DA:C8	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:500:LYS:HE2	1:A:500:LYS:HB3	1.81	0.43
1:A:1003:LYS:HB2	1:A:1003:LYS:HE2	1.74	0.43
1:A:512:SER:O	1:A:516:GLU:HG2	2.18	0.43
1:A:1064:GLU:HG2	1:A:1076:LYS:NZ	2.33	0.43
2:E:98:LEU:HD12	2:E:113:GLN:HA	1.99	0.43
2:E:220:LEU:O	2:E:227:LEU:N	2.44	0.43
5:B:112:U:H2'	5:B:113:C:H6	1.84	0.43
1:A:106:LEU:HD23	1:A:106:LEU:HA	1.87	0.43
1:A:1047:LYS:HE2	1:A:1047:LYS:HB3	1.90	0.43
1:A:1156:LYS:HB2	1:A:1156:LYS:HE3	1.82	0.43
1:A:1357:GLU:OE1	1:A:1359:ARG:NH1	2.52	0.43
5:B:22:U:H2'	5:B:23:U:H6	1.84	0.43
1:A:557:ARG:HA	1:A:595:HIS:CD2	2.53	0.42
1:A:691:ARG:HG3	1:A:696:LEU:CD1	2.49	0.42
1:A:779:GLU:O	1:A:782:LYS:HG2	2.19	0.42
3:C:22:DA:H2''	3:C:23:DT:H72	2.01	0.42
1:A:434:LYS:HE3	1:A:434:LYS:HB2	1.86	0.42
1:A:464:TRP:NE1	1:A:491:PHE:HB2	2.34	0.42
1:A:641:HIS:CE1	1:A:642:LEU:HG	2.55	0.42
1:A:1047:LYS:H	1:A:1047:LYS:HG2	1.69	0.42
1:A:1205:GLU:HB2	1:A:1348:ILE:HD11	2.01	0.42
2:E:115:LEU:HD22	2:E:118:VAL:HB	1.99	0.42
4:D:-1:DA:H2''	4:D:0:DT:C6	2.54	0.42
5:B:19:G:H2'	5:B:20:A:O4'	2.19	0.42
1:A:1000:LYS:HB2	1:A:1073:VAL:HG21	2.01	0.42
2:E:261:ILE:N	2:E:263:GLN:OE1	2.53	0.42
1:A:225:LEU:O	1:A:229:LEU:HB2	2.19	0.42
1:A:465:MET:HE3	1:A:465:MET:HB3	1.98	0.42
2:E:113:GLN:CG	2:E:115:LEU:HG	2.49	0.42
2:E:370:VAL:HG11	2:E:416:ILE:HA	2.01	0.42
5:B:103:C:H2'	5:B:104:G:H8	1.85	0.42
1:A:781:MET:HE3	1:A:781:MET:HB3	1.77	0.42
1:A:814:TYR:HE1	1:A:828:LEU:HB2	1.84	0.42
5:B:104:G:C2	6:F:0:DG:C2	3.08	0.42
2:E:264:LYS:HE3	2:E:264:LYS:HB3	1.92	0.42
5:B:102:U:H2'	5:B:103:C:C6	2.55	0.42
1:A:963:VAL:O	1:A:967:ARG:HG3	2.20	0.42
1:A:861:ASP:OD1	1:A:861:ASP:N	2.49	0.42
1:A:424:ARG:HD2	1:A:424:ARG:HA	1.81	0.41
1:A:528:LYS:HD3	1:A:539:PHE:CZ	2.55	0.41
1:A:1169:MET:HB2	1:A:1169:MET:HE2	1.80	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:52:LEU:HB3	2:E:56:SER:HB2	2.02	0.41
2:E:53:LYS:NZ	2:E:122:VAL:O	2.53	0.41
2:E:174:ASP:HB3	2:E:177:MET:HG2	2.02	0.41
2:E:384:LYS:HB2	2:E:384:LYS:HE2	1.66	0.41
1:A:314:LYS:HA	1:A:314:LYS:HD2	1.83	0.41
1:A:1060:ARG:HG3	1:A:1061:PRO:HD2	2.03	0.41
1:A:1185:LYS:HA	1:A:1185:LYS:HD2	1.75	0.41
1:A:134:THR:OG1	1:A:137:HIS:ND1	2.50	0.41
1:A:241:LEU:HD23	1:A:241:LEU:HA	1.87	0.41
1:A:683:LEU:HD12	1:A:683:LEU:HA	1.89	0.41
5:B:46:A:H2'	5:B:47:A:H8	1.84	0.41
1:A:739:GLN:O	1:A:743:VAL:HG13	2.21	0.41
1:A:226:ILE:HG23	1:A:233:LYS:HA	2.01	0.41
1:A:309:ASN:HB3	1:A:311:GLU:OE1	2.21	0.41
1:A:699:ASP:OD2	1:A:702:LEU:HB2	2.21	0.41
1:A:1064:GLU:HB2	1:A:1074:TRP:HB3	2.01	0.41
5:B:84:A:H2'	5:B:85:C:H6	1.85	0.41
1:A:282:ILE:HG23	1:A:283:GLY:H	1.86	0.41
1:A:784:ILE:HD12	1:A:815:TYR:HB3	2.01	0.41
2:E:274:LYS:HG3	2:E:275:GLU:OE2	2.21	0.41
1:A:37:ASN:ND2	1:A:1364:GLN:OE1	2.53	0.41
1:A:131:LYS:HD2	1:A:132:TYR:CZ	2.56	0.41
1:A:633:GLU:HB2	1:A:652:LYS:HE2	2.03	0.41
2:E:351:LEU:HD23	2:E:351:LEU:HA	1.81	0.41
2:E:437:PRO:HA	2:E:475:VAL:HG23	2.03	0.41
6:F:-1:DC:H2''	6:F:0:DG:H8	1.85	0.41
6:F:-1:DC:H2''	6:F:0:DG:O5'	2.20	0.41
1:A:433:LEU:HD23	1:A:433:LEU:HA	1.78	0.41
2:E:73:GLY:O	2:E:76:PRO:HD2	2.21	0.41
1:A:187:GLN:O	1:A:191:THR:HG23	2.21	0.40
1:A:560:THR:HG23	1:A:563:GLN:H	1.86	0.40
2:E:421:LYS:HB2	2:E:421:LYS:HE2	1.97	0.40
1:A:760:VAL:HG22	1:A:956:ILE:HD12	2.02	0.40
2:E:26:LEU:HA	2:E:33:TRP:NE1	2.36	0.40
2:E:287:THR:O	2:E:291:GLN:HG2	2.21	0.40
1:A:60:GLU:O	1:A:64:LEU:HG	2.22	0.40
1:A:275:LEU:HD11	1:A:290:PHE:CD2	2.57	0.40
1:A:814:TYR:CE1	1:A:828:LEU:HB2	2.56	0.40
1:A:269:ASP:OD1	1:A:269:ASP:N	2.53	0.40
1:A:1158:LYS:HB3	1:A:1158:LYS:HE2	1.77	0.40
2:E:87:LEU:HD23	2:E:87:LEU:HA	1.88	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:195:SER:HB2	2:E:196:PRO:HD3	2.04	0.40
3:C:23:DT:H2''	3:C:24:DT:H72	2.03	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	1303/1367 (95%)	1277 (98%)	26 (2%)	0	100	100
2	E	449/496 (90%)	435 (97%)	13 (3%)	1 (0%)	43	76
All	All	1752/1863 (94%)	1712 (98%)	39 (2%)	1 (0%)	49	80

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	E	223	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	1180/1224 (96%)	1180 (100%)	0	100	100
2	E	383/425 (90%)	383 (100%)	0	100	100
All	All	1563/1649 (95%)	1563 (100%)	0	100	100

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

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	46	ASN
1	A	175	ASN
1	A	415	HIS
1	A	436	ASN
1	A	674	GLN
1	A	690	ASN
1	A	698	HIS
1	A	920	GLN
1	A	985	HIS
1	A	1101	GLN
1	A	1177	ASN
1	A	1264	HIS
1	A	1350	GLN
2	E	45	GLN
2	E	200	ASN
2	E	204	HIS
2	E	221	GLN
2	E	249	ASN
2	E	277	GLN
2	E	291	GLN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
5	B	110/115 (95%)	21 (19%)	0
6	F	0/35	-	-
All	All	110/150 (73%)	21 (19%)	0

All (21) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
5	B	2	G
5	B	3	C
5	B	4	C
5	B	5	G
5	B	20	A
5	B	28	A
5	B	35	A

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Mol	Chain	Res	Type
5	B	36	A
5	B	51	A
5	B	53	G
5	B	56	U
5	B	59	U
5	B	68	A
5	B	77	A
5	B	81	G
5	B	87	G
5	B	89	G
5	B	91	C
5	B	102	U
5	B	103	C
5	B	104	G

There are no RNA pucker outliers to report.

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

1 non-standard protein/DNA/RNA residue is 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
6	2DA	F	1	6,5	19,22,23	0.31	0	25,31,34	0.32	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	2DA	F	1	6,5	-	2/7/18/19	0/3/3/3

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	F	1	2DA	C4'-C5'-O5'-P
6	F	1	2DA	O4'-C1'-N9-C8

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides 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](#)

There are no chain breaks in this entry.

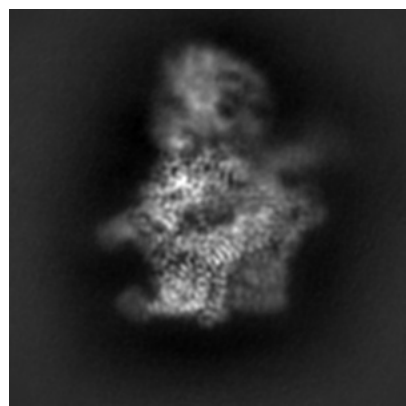
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-37859. These allow visual inspection of the internal detail of the map and identification of artifacts.

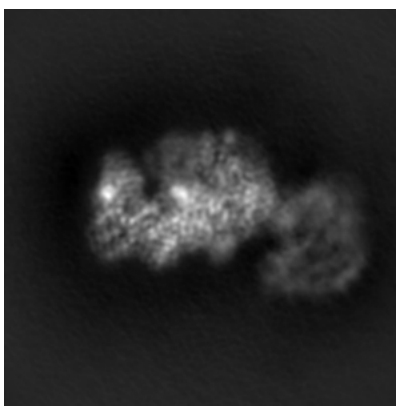
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

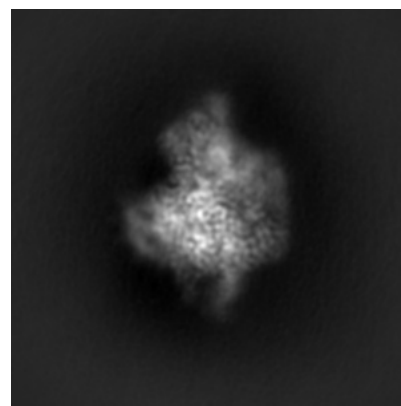
6.1.1 Primary map



X

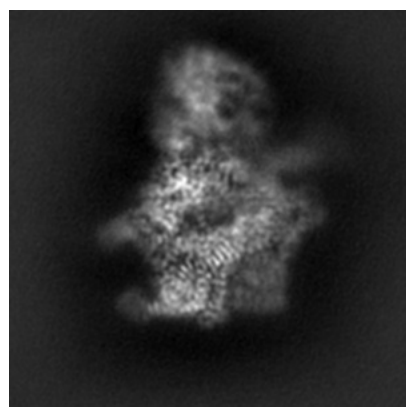


Y

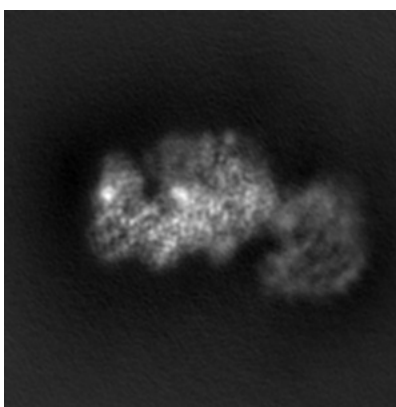


Z

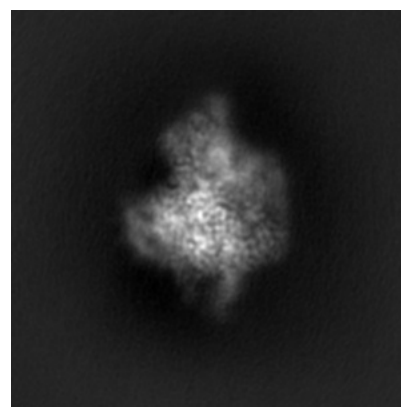
6.1.2 Raw map



X



Y

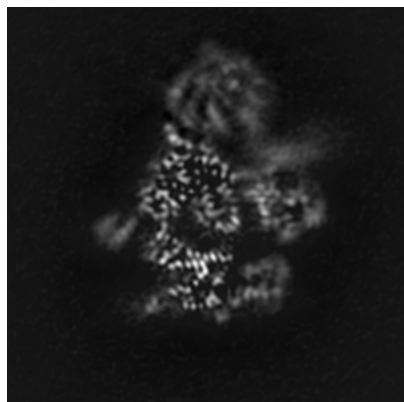


Z

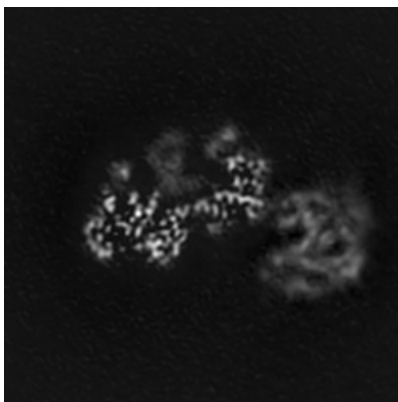
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

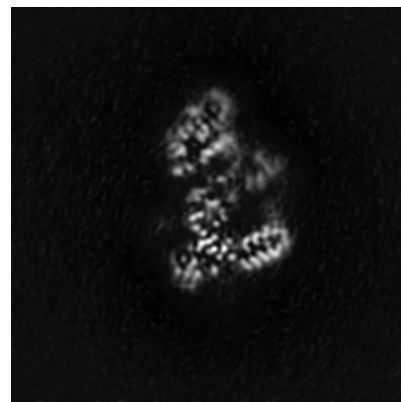
6.2.1 Primary map



X Index: 86

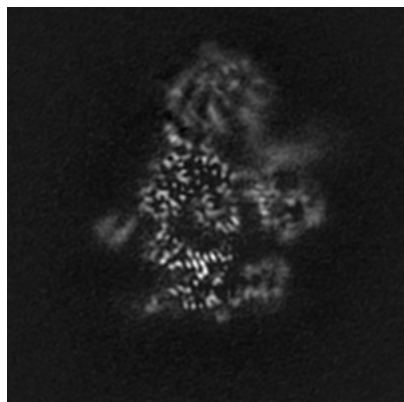


Y Index: 86

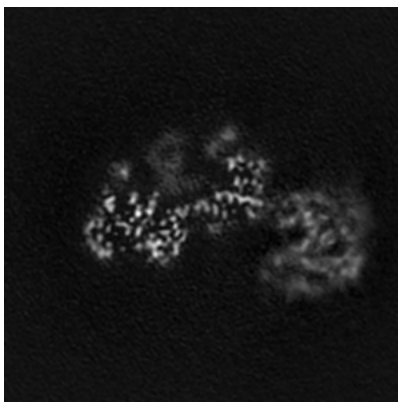


Z Index: 86

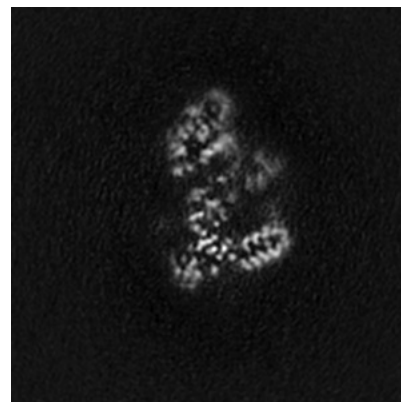
6.2.2 Raw map



X Index: 86



Y Index: 86

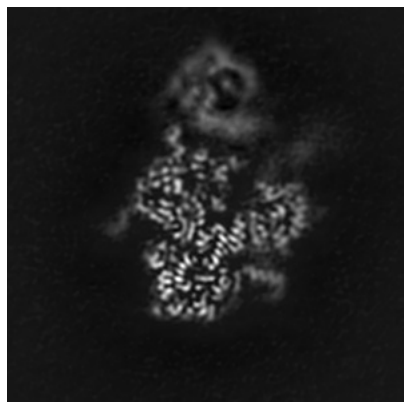


Z Index: 86

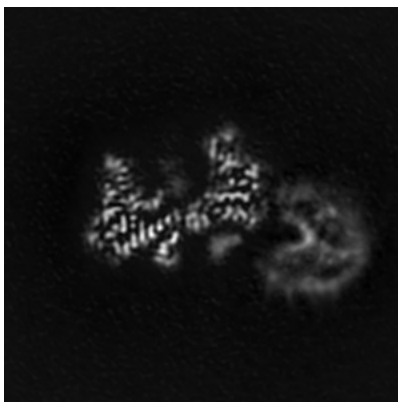
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

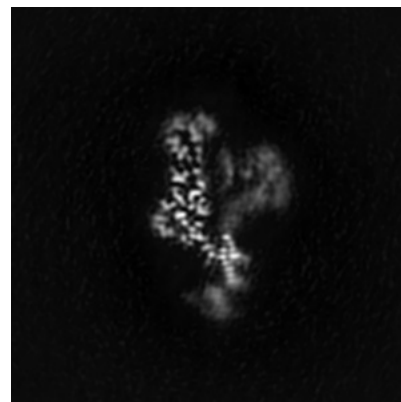
6.3.1 Primary map



X Index: 80

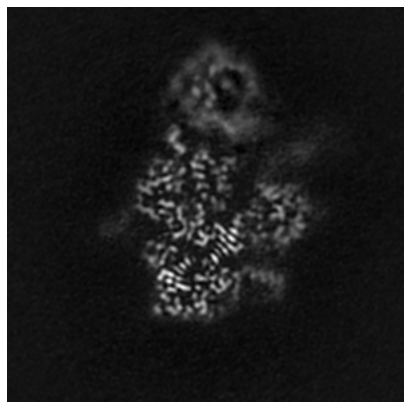


Y Index: 82

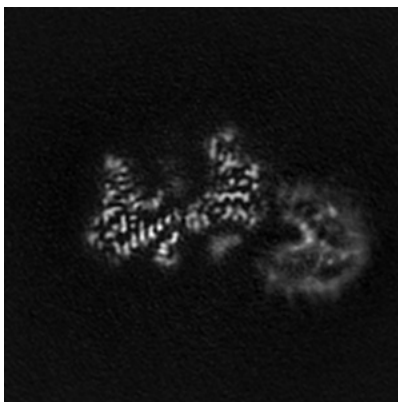


Z Index: 73

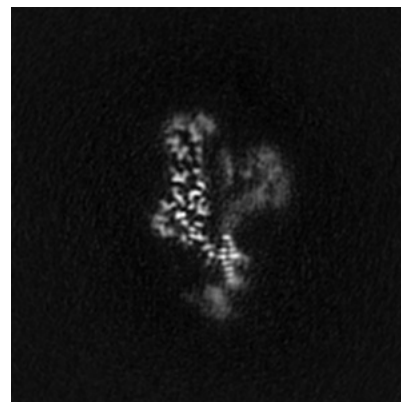
6.3.2 Raw map



X Index: 81



Y Index: 82

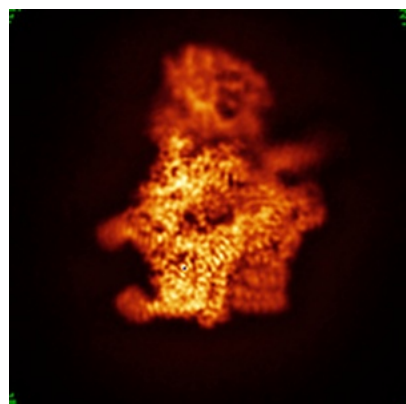


Z Index: 73

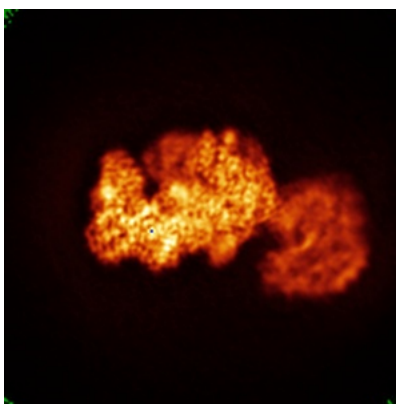
The images above show the largest variance slices of the map in three orthogonal directions.

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

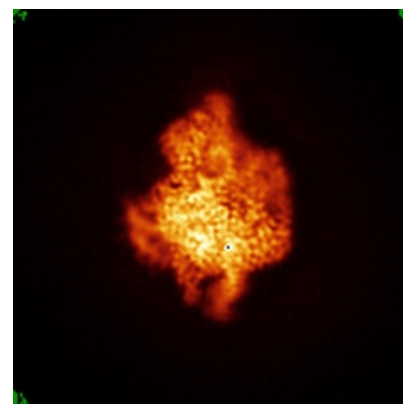
6.4.1 Primary map



X

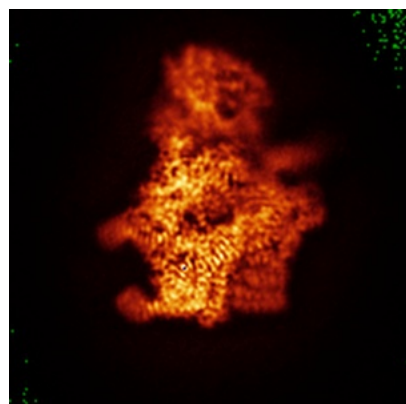


Y

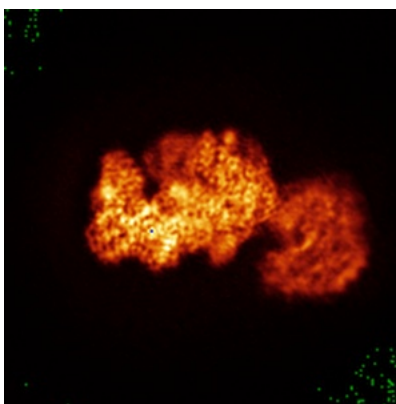


Z

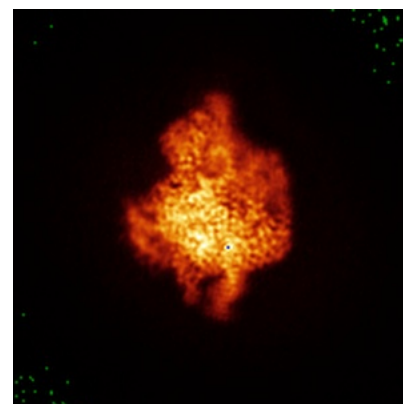
6.4.2 Raw map



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](#)

This section was not generated.

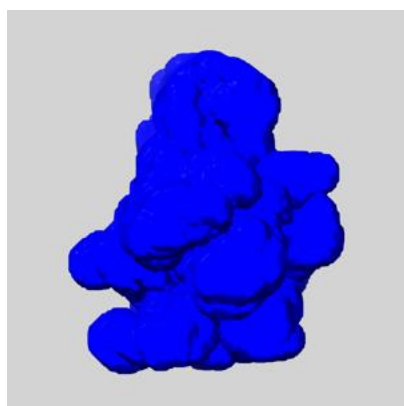
6.6 Mask visualisation [i](#)

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

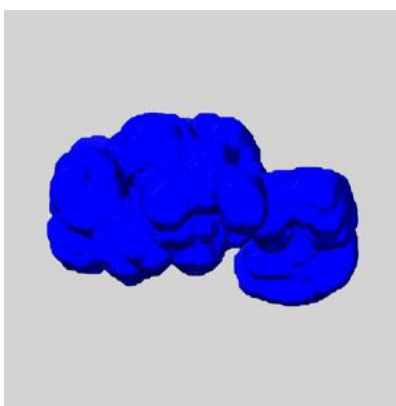
A mask typically either:

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

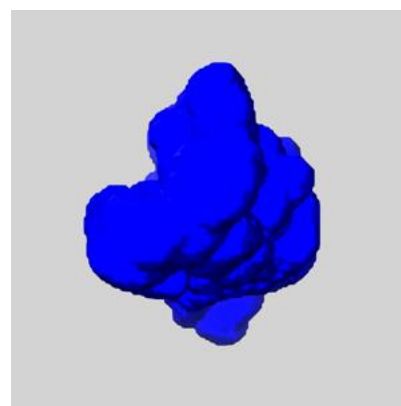
6.6.1 emd_37859_msk_1.map [i](#)



X



Y

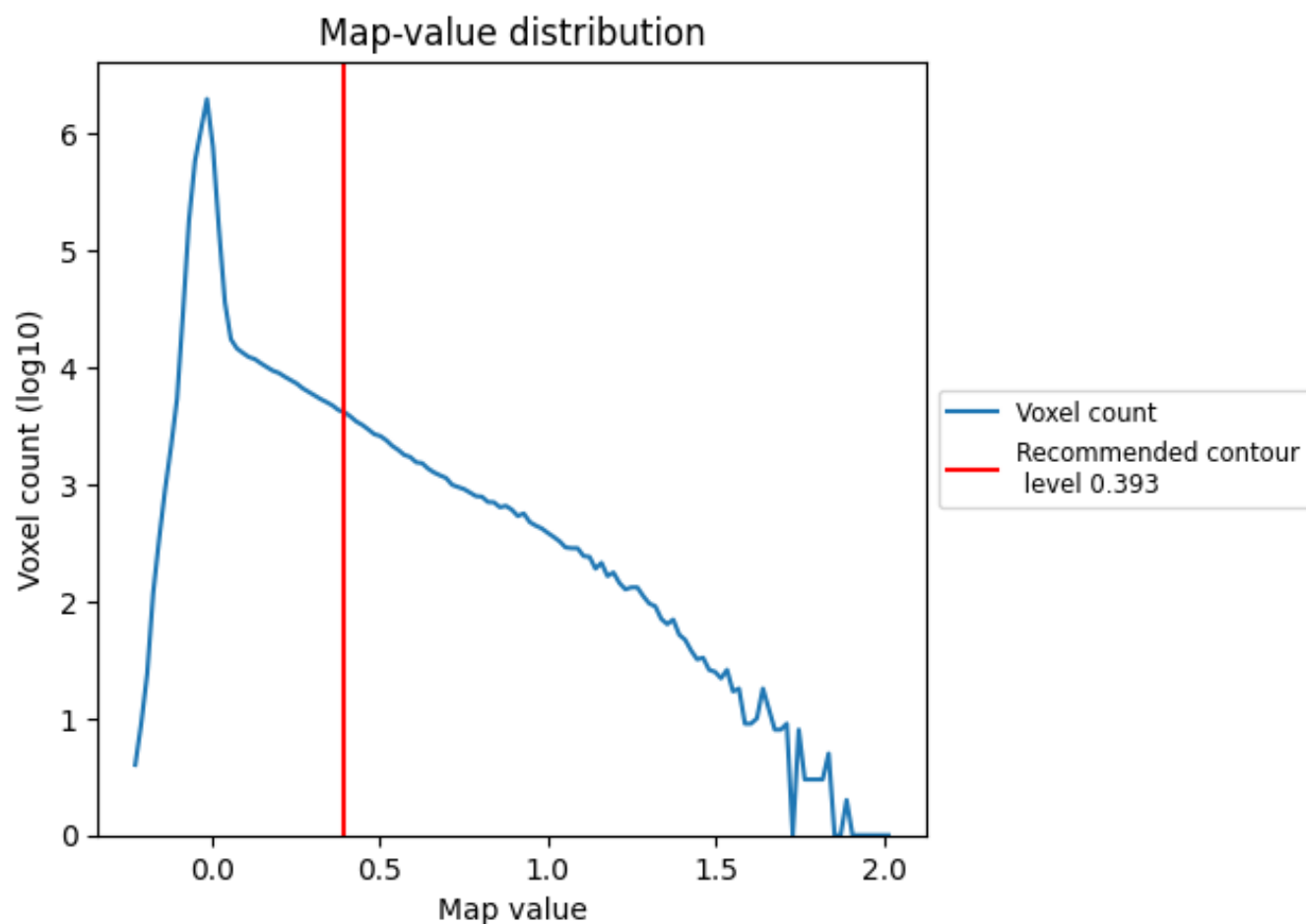


Z

7 Map analysis [i](#)

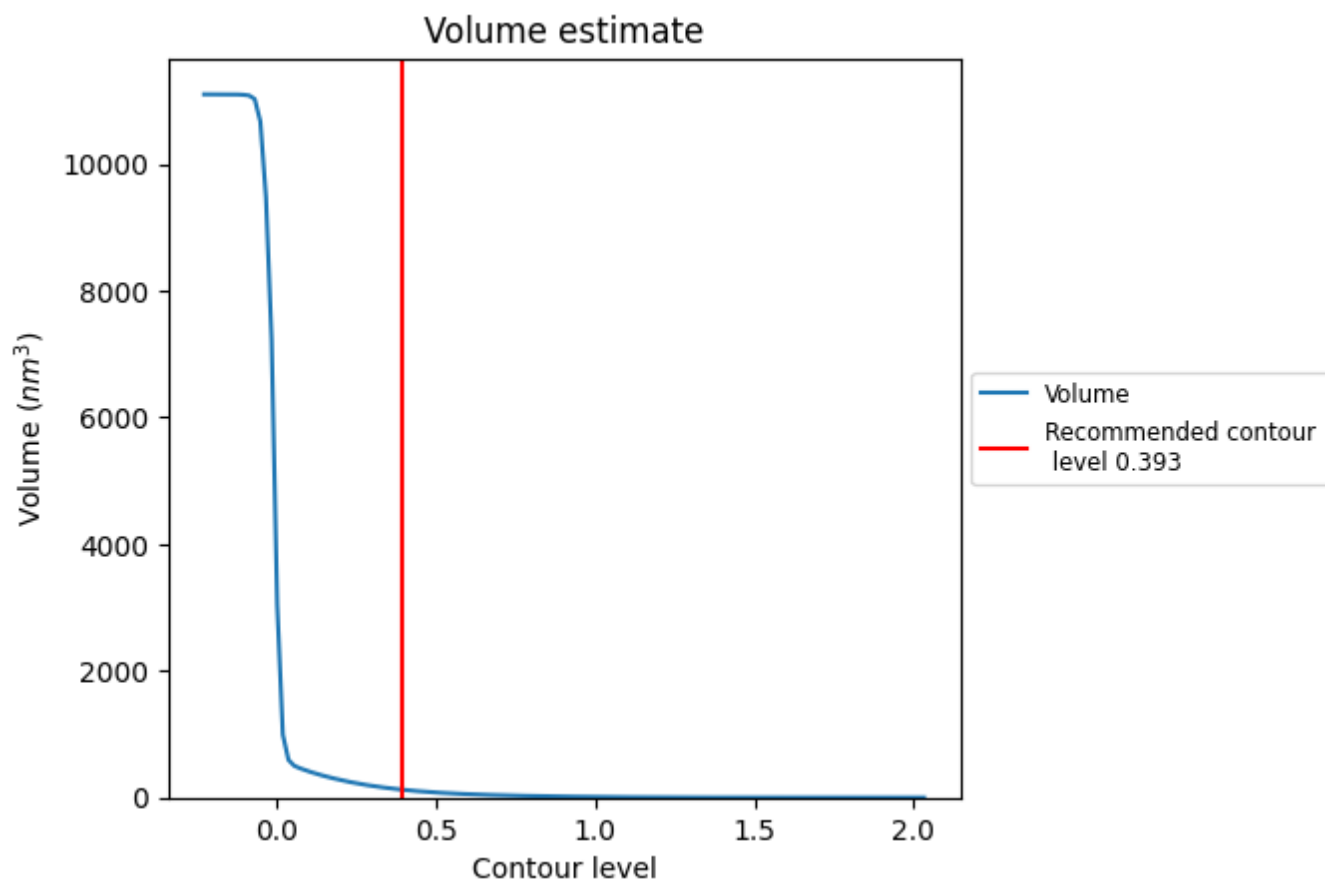
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



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

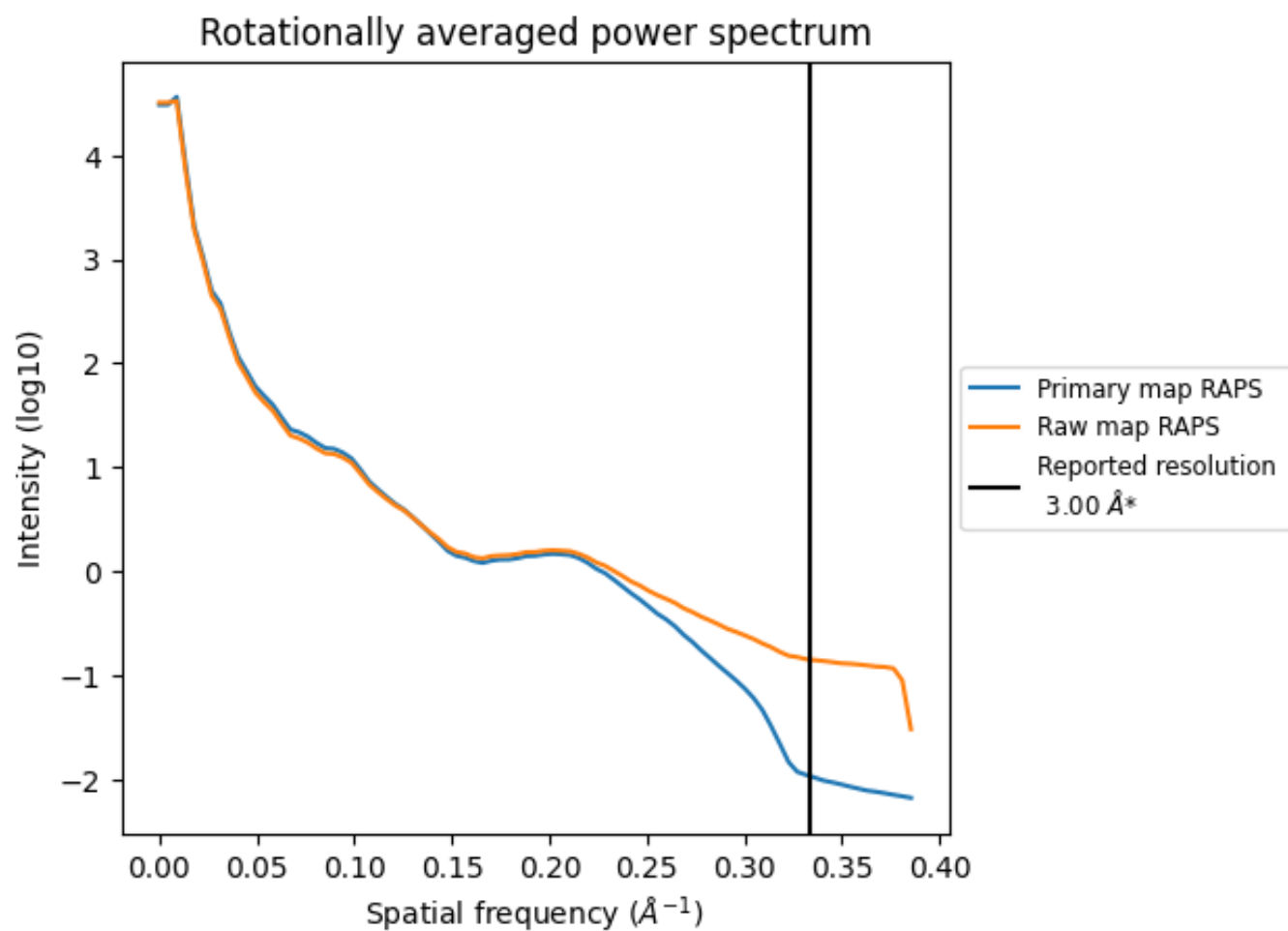
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 125 nm³; this corresponds to an approximate mass of 113 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 ⓘ

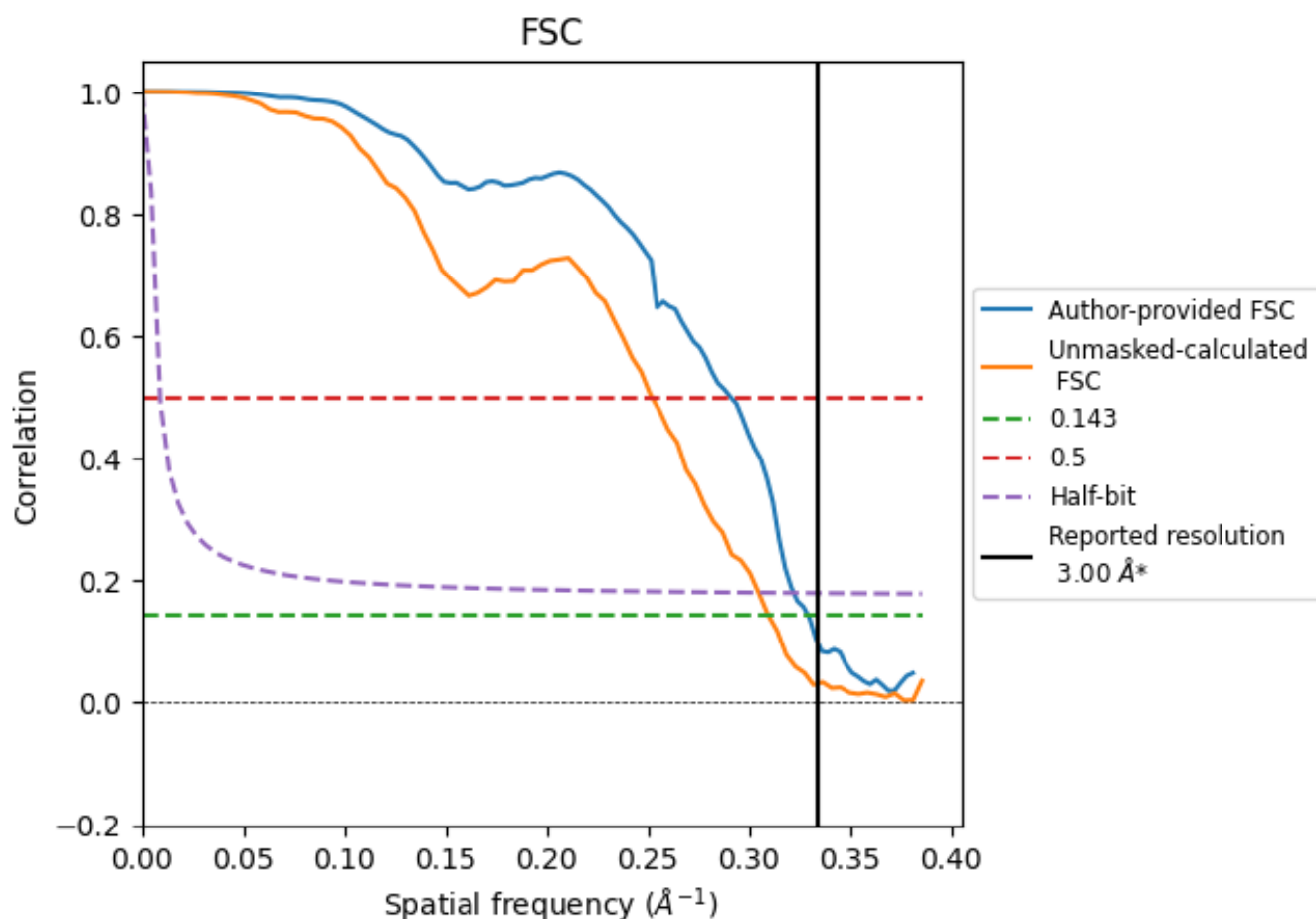


*Reported resolution corresponds to spatial frequency of 0.333 $\text{\AA}^{-1}$

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 0.333 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.00	-	-
Author-provided FSC curve	3.04	3.44	3.10
Unmasked-calculated*	3.23	3.97	3.28

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

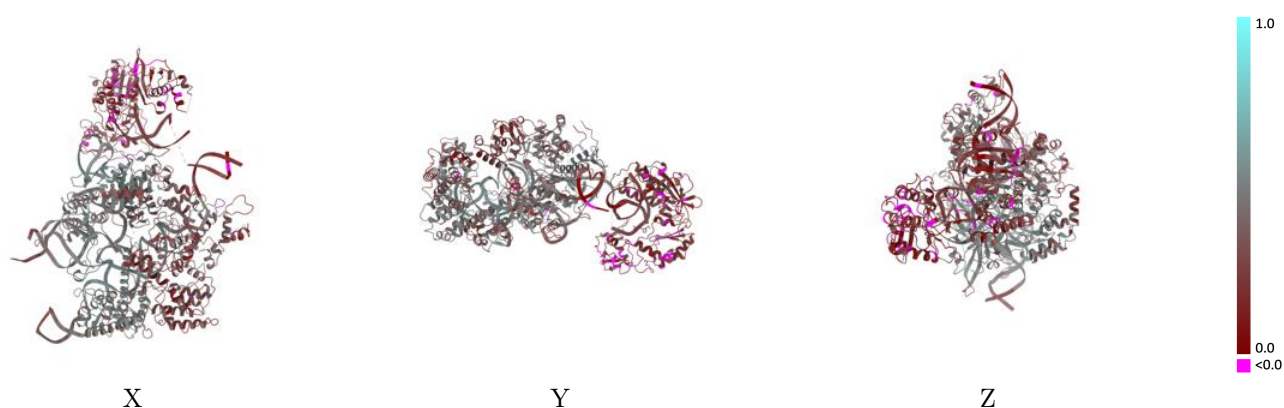
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-37859 and PDB model 8WUT. Per-residue inclusion information can be found in section 3 on page 6.

9.1 Map-model overlay [i](#)

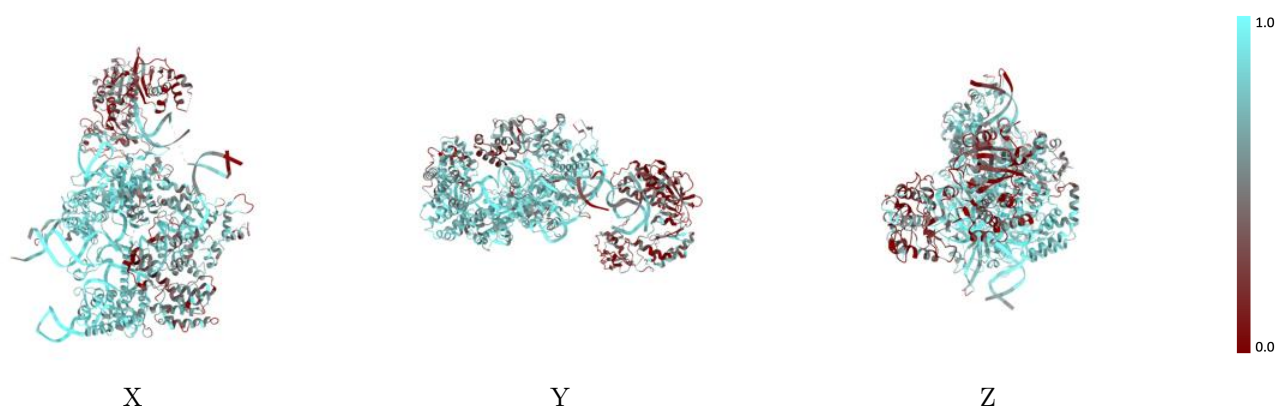
This section was not generated.

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



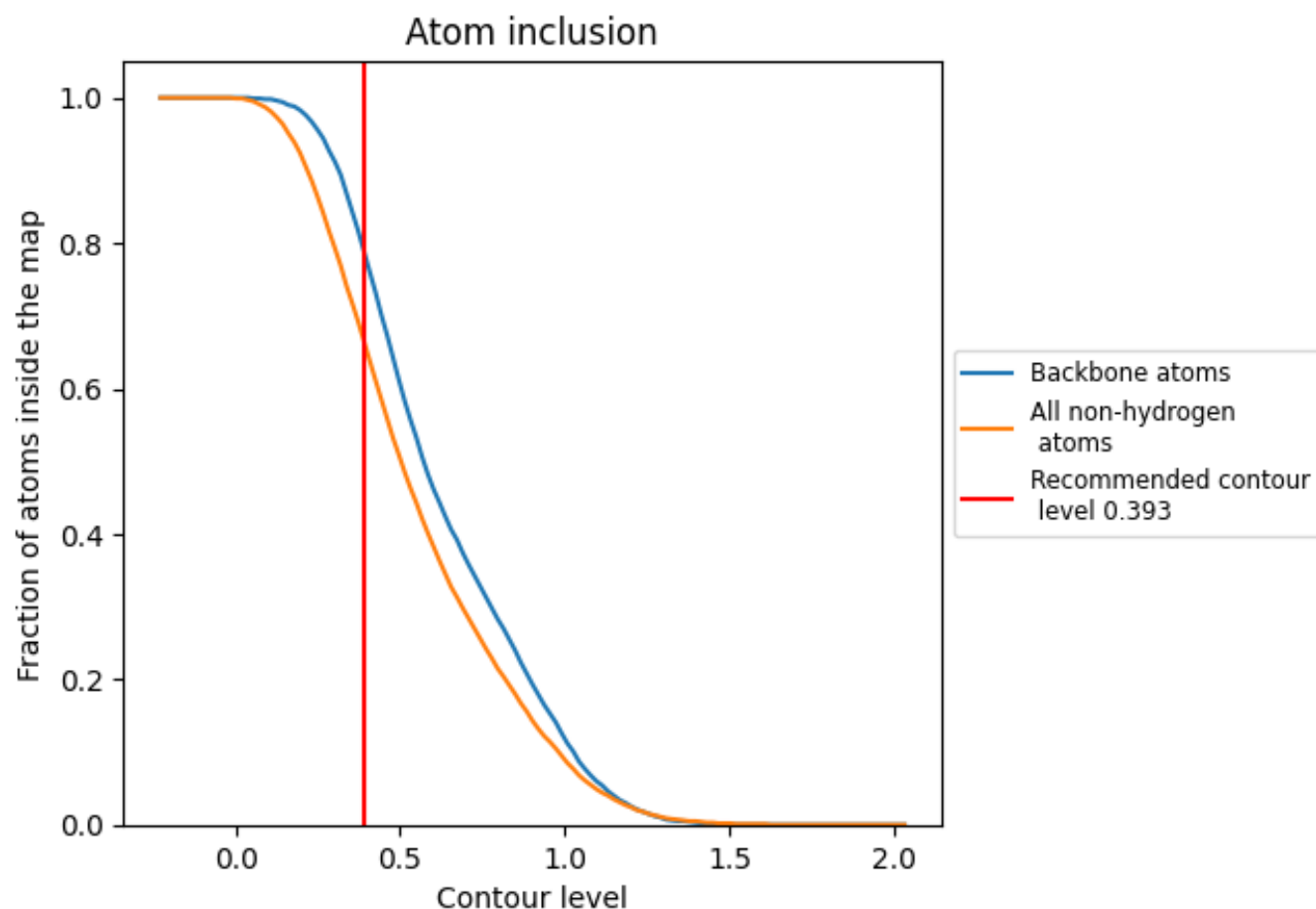
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.393).

9.4 Atom inclusion [i](#)



At the recommended contour level, 78% of all backbone atoms, 66% 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 (0.393) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.6610	0.3640
A	0.7080	0.4120
B	0.9330	0.4490
C	0.8350	0.4180
D	0.8910	0.4440
E	0.2790	0.1620
F	0.5960	0.2070

1.0

0.0

<0.0