



Full wwPDB EM Validation Report ⓘ

Mar 20, 2026 – 03:34 AM UTC

PDB ID : 7OSF / pdb_00007osf
EMDB ID : EMD-13049
Title : ABC Transporter complex NosDFYL, R-domain 1
Authors : Mueller, C.; Zhang, L.; Lu, W.; Du, J.; Einsle, O.
Deposited on : 2021-06-08
Resolution : 3.80 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:

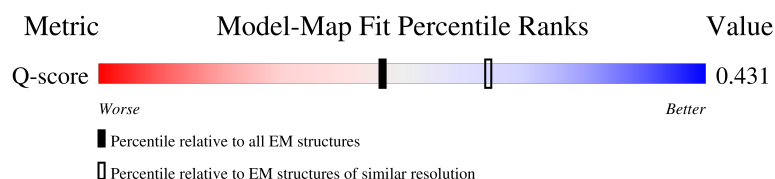
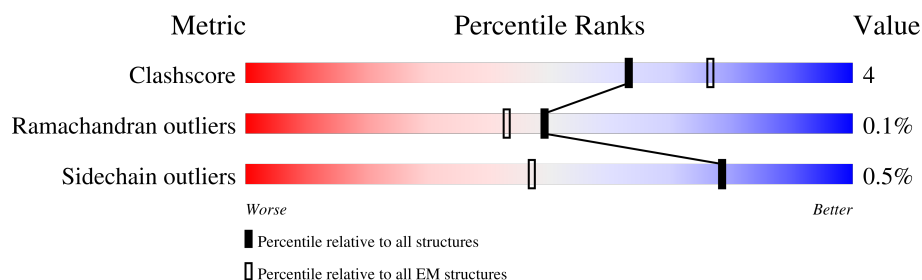
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh &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

1 Overall quality at a glance

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

The reported resolution of this entry is 3.80 Å.

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



Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	10198 (3.30 - 4.30)

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	436	
2	B	308	
2	C	308	
3	D	276	

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Mol	Chain	Length	Quality of chain
3	E	276	 84% 6% 10%
4	H	190	 66% 9% 25%

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 12826 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 Probable ABC transporter binding protein NosD.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	392	Total	C	N	O	S	0	0
			3077	1915	556	596	10		

- Molecule 2 is a protein called Probable ABC transporter ATP-binding protein NosF.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	308	Total	C	N	O	S	0	0
			2375	1483	443	442	7		
2	C	307	Total	C	N	O	S	0	0
			2367	1478	442	441	6		

- Molecule 3 is a protein called Probable ABC transporter permease protein NosY.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	D	267	Total	C	N	O	S	0	0
			2015	1355	319	334	7		
3	E	248	Total	C	N	O	S	0	0
			1896	1284	300	306	6		

- Molecule 4 is a protein called Copper-binding lipoprotein NosL.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	H	143	Total	C	N	O	S	0	0
			1092	685	189	209	9		

- Molecule 5 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

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

- Molecule 6 is COPPER (II) ION (CCD ID: CU) (formula: Cu) (labeled as "Ligand of

Interest" by depositor).

Mol	Chain	Residues	Atoms		AltConf
6	A	1	Total	Cu	0
			1	1	

- Molecule 7 is ZINC ION (CCD ID: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		AltConf
7	H	1	Total	Zn	0
			1	1	

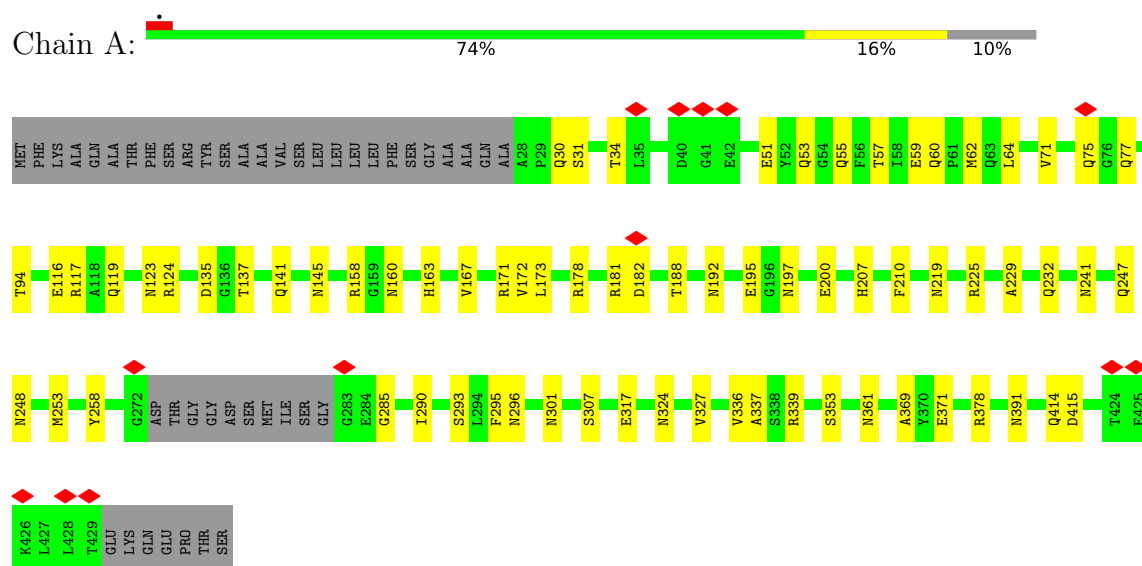
- Molecule 8 is water.

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

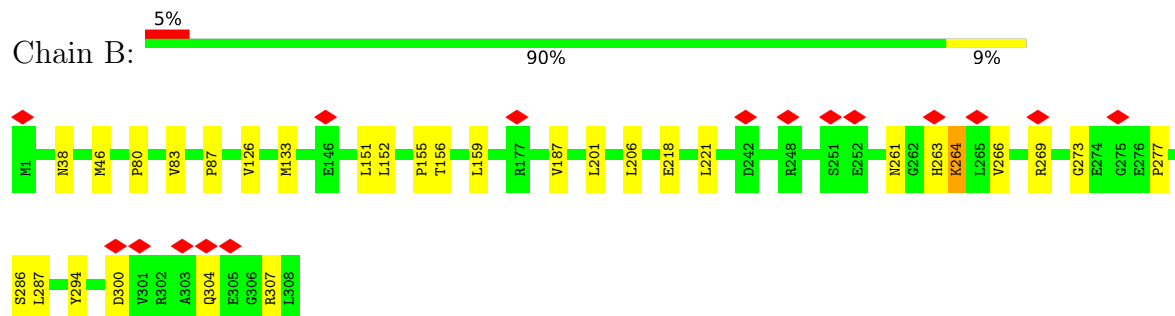
3 Residue-property plots

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

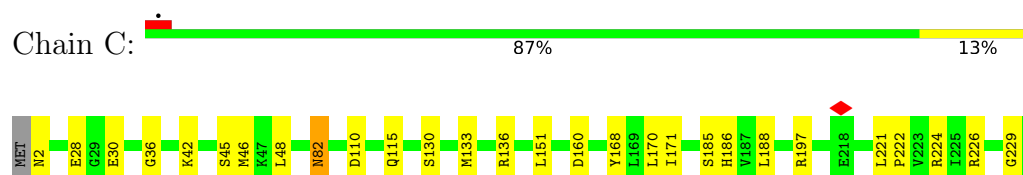
- Molecule 1: Probable ABC transporter binding protein NosD

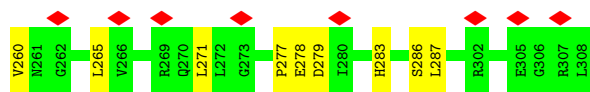


- Molecule 2: Probable ABC transporter ATP-binding protein NosF



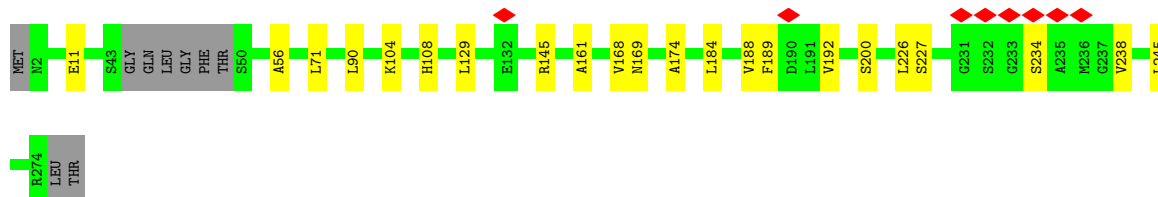
- Molecule 2: Probable ABC transporter ATP-binding protein NosF





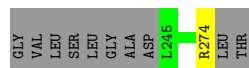
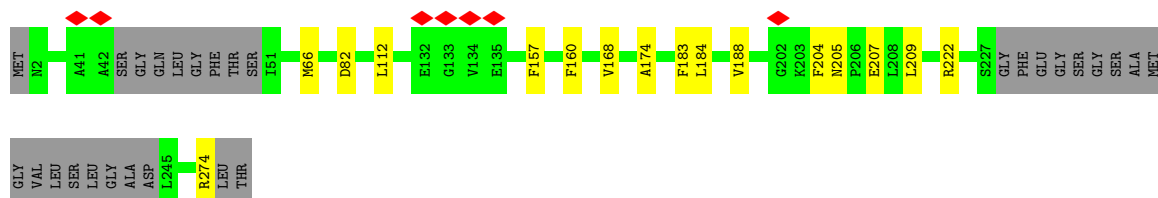
- Molecule 3: Probable ABC transporter permease protein NosY

Chain D: 89% 8%



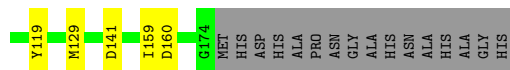
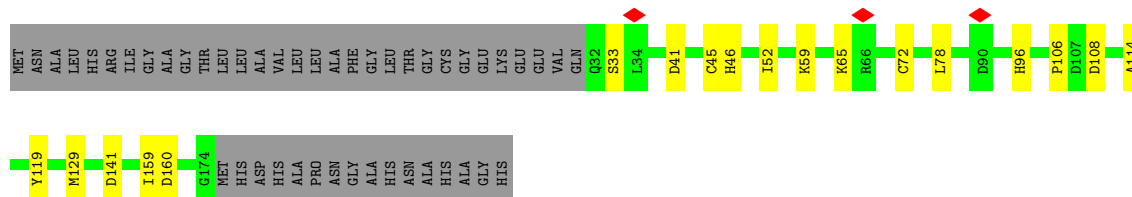
- Molecule 3: Probable ABC transporter permease protein NosY

Chain E: 84% 6% 10%



- Molecule 4: Copper-binding lipoprotein NosL

Chain H: 66% 9% 25%



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	37715	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)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.082	Depositor
Minimum map value	-0.026	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.004	Depositor
Recommended contour level	0.022	Depositor
Map size ($\text{\AA}$)	259.2, 259.2, 259.2	wwPDB
Map dimensions	200, 200, 200	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	1.296, 1.296, 1.296	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, CU, ZN

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	0.35	0/3140	0.66	0/4262
2	B	0.35	0/2410	0.70	0/3257
2	C	0.33	0/2402	0.65	2/3247 (0.1%)
3	D	0.32	0/2060	0.55	0/2809
3	E	0.31	0/1939	0.60	0/2646
4	H	0.30	0/1116	0.58	0/1505
All	All	0.33	0/13067	0.63	2/17726 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
2	B	0	2
4	H	0	1
All	All	0	3

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	260	VAL	CA-C-N	6.09	133.17	121.54
2	C	260	VAL	C-N-CA	6.09	133.17	121.54

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	263	HIS	Peptide
2	B	266	VAL	Peptide
4	H	108	ASP	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	3077	0	2962	42	0
2	B	2375	0	2429	16	0
2	C	2367	0	2417	23	0
3	D	2015	0	2136	13	0
3	E	1896	0	2026	8	0
4	H	1092	0	1047	11	0
5	A	1	0	0	0	0
6	A	1	0	0	0	0
7	H	1	0	0	0	0
8	A	1	0	0	0	0
All	All	12826	0	13017	108	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 4.

All (108) 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:A:301:ASN:H	1:A:324:ASN:HD22	1.41	0.67
2:C:130:SER:H	2:C:133:MET:HE3	1.64	0.62
1:A:116:GLU:HG2	1:A:117:ARG:HG3	1.83	0.60
1:A:369:ALA:HB3	3:D:234:SER:HB2	1.84	0.60
1:A:119:GLN:HG3	1:A:141:GLN:HB2	1.85	0.57
2:B:264:LYS:NZ	2:C:279:ASP:OD2	2.38	0.57
1:A:327:VAL:HA	1:A:353:SER:HB3	1.86	0.57
1:A:160:ASN:ND2	1:A:182:ASP:OD1	2.39	0.56
1:A:75:GLN:HB2	1:A:77:GLN:HE22	1.72	0.55
1:A:141:GLN:HG2	1:A:171:ARG:HB2	1.88	0.55
2:C:229:GLY:HA3	2:C:278:GLU:HG2	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:286:SER:OG	2:C:287:LEU:N	2.41	0.54
4:H:78:LEU:HD13	4:H:159:ILE:HG21	1.89	0.54
2:C:30:GLU:OE2	2:C:197:ARG:NH1	2.39	0.54
1:A:160:ASN:OD1	1:A:163:HIS:ND1	2.39	0.54
1:A:337:ALA:O	1:A:378:ARG:NH2	2.41	0.54
2:C:115:GLN:HE21	2:C:170:LEU:HD11	1.70	0.54
2:C:160:ASP:N	2:C:160:ASP:OD1	2.40	0.53
1:A:158:ARG:HB2	1:A:181:ARG:HB2	1.89	0.53
1:A:30:GLN:HG3	1:A:57:THR:HB	1.91	0.53
2:B:277:PRO:HG2	2:C:265:LEU:HD11	1.91	0.53
1:A:301:ASN:OD1	1:A:324:ASN:ND2	2.42	0.53
3:D:11:GLU:OE2	3:D:104:LYS:NZ	2.42	0.52
2:B:269:ARG:O	2:B:273:GLY:N	2.42	0.52
3:E:205:ASN:ND2	3:E:207:GLU:OE1	2.42	0.52
1:A:317:GLU:OE1	1:A:339:ARG:NH1	2.42	0.52
1:A:371:GLU:HG3	1:A:414:GLN:HB3	1.93	0.51
1:A:178:ARG:HG2	1:A:200:GLU:HB2	1.92	0.51
4:H:141:ASP:N	4:H:141:ASP:OD1	2.43	0.50
1:A:197:ASN:OD1	1:A:219:ASN:ND2	2.44	0.49
3:E:66:MET:HE1	3:E:112:LEU:HB2	1.92	0.49
3:D:145:ARG:NH2	3:D:227:SER:OG	2.45	0.49
1:A:188:THR:HG22	1:A:210:PHE:HB2	1.95	0.49
2:B:152:LEU:HB3	2:B:155:PRO:HB3	1.94	0.49
1:A:192:ASN:N	1:A:192:ASN:OD1	2.44	0.49
4:H:114:ALA:O	4:H:119:TYR:OH	2.32	0.48
2:C:188:LEU:HD23	2:C:287:LEU:HG	1.95	0.48
3:D:11:GLU:OE1	3:D:108:HIS:NE2	2.47	0.48
1:A:225:ARG:HA	1:A:247:GLN:HB2	1.96	0.47
2:B:201:LEU:HB3	2:B:206:LEU:HD12	1.96	0.47
1:A:135:ASP:N	1:A:135:ASP:OD1	2.47	0.47
1:A:258:TYR:N	1:A:295:PHE:O	2.47	0.46
3:D:184:LEU:HA	3:D:188:VAL:HB	1.96	0.46
2:C:226:ARG:HA	2:C:256:GLU:HA	1.97	0.46
1:A:31:SER:O	1:A:34:THR:OG1	2.33	0.46
1:A:173:LEU:HG	1:A:195:GLU:HB2	1.98	0.46
2:B:221:LEU:O	2:B:261:ASN:ND2	2.48	0.46
3:D:168:VAL:HG21	3:D:174:ALA:HB2	1.97	0.46
1:A:55:GLN:NE2	1:A:77:GLN:O	2.42	0.46
4:H:45:CYS:HB2	4:H:72:CYS:HB3	1.96	0.45
4:H:160:ASP:N	4:H:160:ASP:OD1	2.49	0.45
1:A:123:ASN:OD1	1:A:145:ASN:ND2	2.49	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:245:LEU:HD23	3:D:245:LEU:HA	1.87	0.45
2:B:286:SER:OG	2:B:287:LEU:N	2.49	0.45
2:C:46:MET:HG3	2:C:151:LEU:HD21	1.99	0.45
4:H:45:CYS:SG	4:H:46:HIS:N	2.90	0.45
1:A:137:THR:H	1:A:167:VAL:HG12	1.81	0.45
2:B:87:PRO:HA	2:B:126:VAL:HB	1.99	0.44
4:H:33:SER:HB3	4:H:65:LYS:HB3	1.99	0.44
2:B:155:PRO:HG2	2:B:156:THR:HG23	2.00	0.44
3:D:189:PHE:HA	3:D:192:VAL:HG12	1.99	0.44
1:A:253:MET:HG3	1:A:290:ILE:HA	1.99	0.44
3:D:169:ASN:OD1	3:D:169:ASN:N	2.51	0.44
2:B:133:MET:HE3	2:B:133:MET:HB2	1.93	0.43
2:C:2:ASN:N	2:C:28:GLU:OE2	2.52	0.43
2:B:46:MET:HE1	2:B:151:LEU:HB3	2.00	0.43
2:C:231:SER:OG	2:C:232:GLU:N	2.52	0.43
1:A:59:GLU:HG2	1:A:60:GLN:HG2	2.01	0.43
2:C:82:ASN:N	2:C:82:ASN:OD1	2.49	0.43
3:D:90:LEU:HD23	3:D:90:LEU:HA	1.88	0.42
3:D:200:SER:O	3:D:200:SER:OG	2.33	0.42
3:E:204:PHE:HE2	3:E:209:LEU:HB2	1.83	0.42
1:A:415:ASP:N	1:A:415:ASP:OD1	2.51	0.42
2:C:224:ARG:HB2	2:C:283:HIS:HB2	2.01	0.42
3:E:157:PHE:HA	3:E:160:PHE:HB2	2.01	0.42
1:A:248:ASN:HA	1:A:285:GLY:HA2	2.01	0.42
1:A:94:THR:HG23	1:A:124:ARG:HB2	2.01	0.42
2:B:156:THR:HA	2:B:159:LEU:HD12	2.02	0.42
2:C:271:LEU:HD12	2:C:277:PRO:HG3	2.01	0.42
3:E:168:VAL:HG21	3:E:174:ALA:HB2	2.02	0.42
2:B:80:PRO:HG2	2:B:83:VAL:HB	2.02	0.42
3:D:71:LEU:HD23	3:D:161:ALA:HB2	2.02	0.42
2:C:45:SER:HA	2:C:48:LEU:HD12	2.02	0.41
2:B:38:ASN:ND2	2:B:300:ASP:OD1	2.53	0.41
4:H:96:HIS:CD2	4:H:106:PRO:HB2	2.55	0.41
2:B:304:GLN:HB2	2:B:307:ARG:HG2	2.02	0.41
1:A:285:GLY:O	1:A:307:SER:OG	2.33	0.41
2:B:218:GLU:OE1	2:B:294:TYR:OH	2.33	0.41
3:E:184:LEU:HA	3:E:188:VAL:HB	2.03	0.41
1:A:62:MET:HG3	1:A:64:LEU:HD23	2.02	0.41
1:A:293:SER:OG	1:A:296:ASN:ND2	2.54	0.41
2:C:168:TYR:HA	2:C:171:ILE:HB	2.03	0.41
4:H:41:ASP:N	4:H:41:ASP:OD1	2.47	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:207:HIS:HA	1:A:229:ALA:HB3	2.03	0.41
1:A:219:ASN:H	1:A:241:ASN:HD22	1.69	0.41
1:A:391:ASN:HB3	3:E:222:ARG:HD2	2.03	0.41
2:C:221:LEU:HA	2:C:222:PRO:HD3	1.90	0.41
2:C:133:MET:HA	2:C:136:ARG:HD2	2.03	0.40
3:E:82:ASP:OD2	3:E:274:ARG:NH2	2.54	0.40
1:A:232:GLN:NE2	4:H:129:MET:O	2.54	0.40
2:C:185:SER:OG	2:C:186:HIS:N	2.53	0.40
2:C:36:GLY:H	2:C:42:LYS:HD2	1.86	0.40
3:D:56:ALA:HA	3:D:226:LEU:HD11	2.03	0.40
4:H:52:ILE:HG23	4:H:59:LYS:HG2	2.03	0.40
1:A:51:GLU:HG2	1:A:71:VAL:HB	2.04	0.40
1:A:53:GLN:HE21	1:A:75:GLN:HG2	1.86	0.40
1:A:361:ASN:OD1	1:A:361:ASN:N	2.54	0.40
2:C:110:ASP:OD1	2:C:110:ASP:N	2.41	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	388/436 (89%)	357 (92%)	31 (8%)	0	100	100
2	B	306/308 (99%)	270 (88%)	35 (11%)	1 (0%)	36	67
2	C	305/308 (99%)	277 (91%)	28 (9%)	0	100	100
3	D	263/276 (95%)	256 (97%)	7 (3%)	0	100	100
3	E	242/276 (88%)	237 (98%)	5 (2%)	0	100	100
4	H	141/190 (74%)	133 (94%)	8 (6%)	0	100	100
All	All	1645/1794 (92%)	1530 (93%)	114 (7%)	1 (0%)	49	79

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	264	LYS

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	332/365 (91%)	330 (99%)	2 (1%)	78	79
2	B	251/251 (100%)	250 (100%)	1 (0%)	84	83
2	C	250/251 (100%)	249 (100%)	1 (0%)	84	83
3	D	205/212 (97%)	203 (99%)	2 (1%)	68	74
3	E	193/212 (91%)	192 (100%)	1 (0%)	81	81
4	H	112/144 (78%)	112 (100%)	0	100	100
All	All	1343/1435 (94%)	1336 (100%)	7 (0%)	78	81

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	172	VAL
1	A	336	VAL
2	B	187	VAL
2	C	82	ASN
3	D	129	LEU
3	D	238	VAL
3	E	183	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (29) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	53	GLN
1	A	77	GLN
1	A	123	ASN
1	A	141	GLN
1	A	145	ASN
1	A	175	ASN

Continued on next page...

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Mol	Chain	Res	Type
1	A	176	HIS
1	A	179	ASN
1	A	190	ASN
1	A	212	ASN
1	A	241	ASN
1	A	248	ASN
1	A	301	ASN
1	A	324	ASN
2	B	2	ASN
2	B	23	ASN
2	B	88	GLN
2	B	97	HIS
2	B	176	GLN
2	B	245	HIS
2	B	283	HIS
2	C	97	HIS
2	C	115	GLN
2	C	284	GLN
3	D	3	GLN
3	D	205	ASN
3	D	225	ASN
3	E	225	ASN
4	H	161	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 3 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

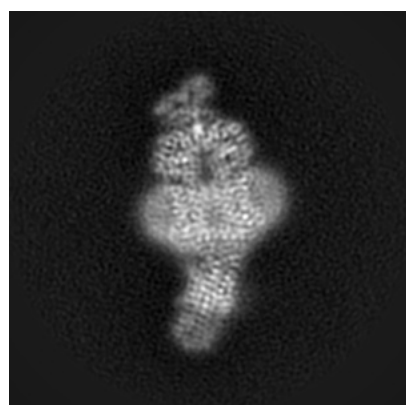
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-13049. These allow visual inspection of the internal detail of the map and identification of artifacts.

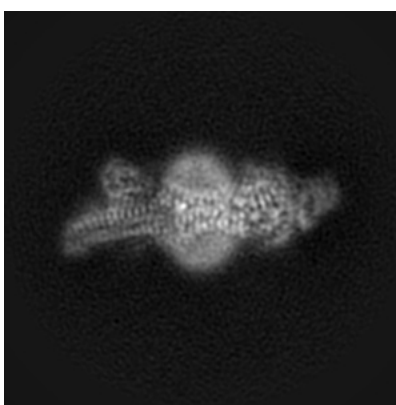
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

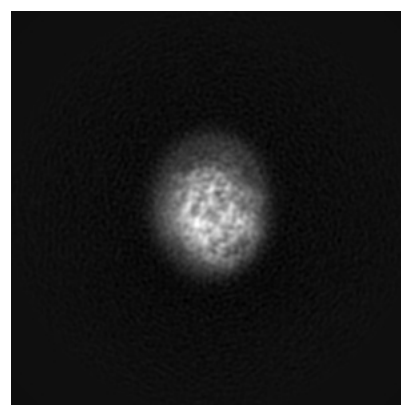
6.1.1 Primary map



X



Y

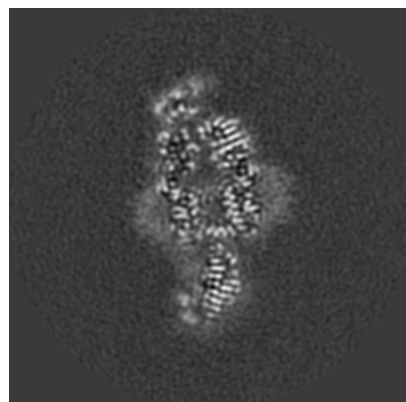


Z

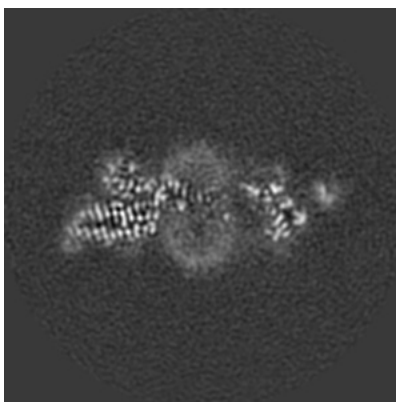
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

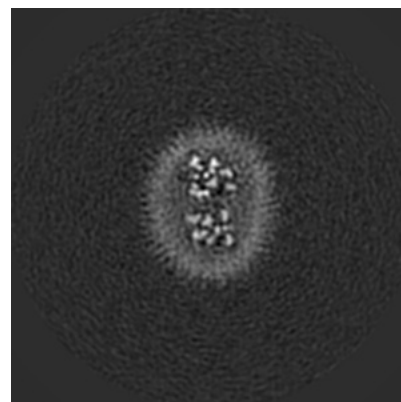
6.2.1 Primary map



X Index: 100



Y Index: 100

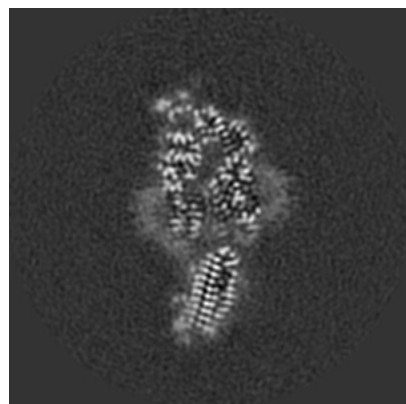


Z Index: 100

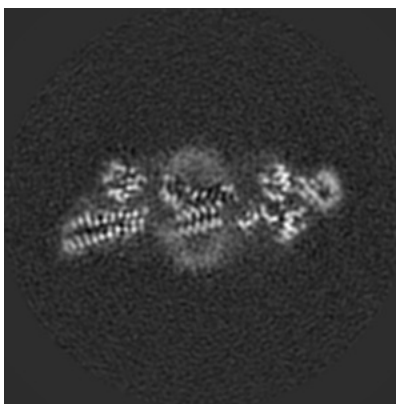
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

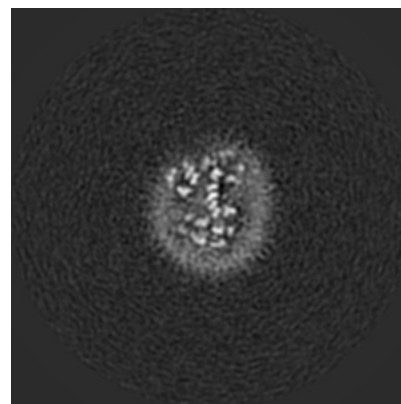
6.3.1 Primary map



X Index: 96



Y Index: 94



Z Index: 89

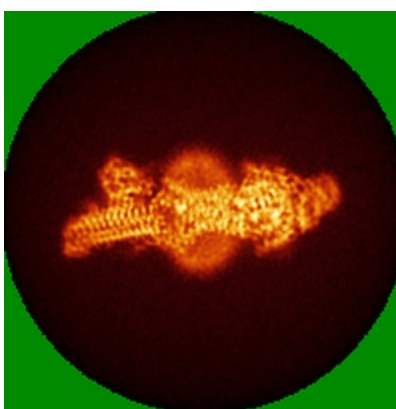
The images above show the largest variance slices of the map in three orthogonal directions.

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

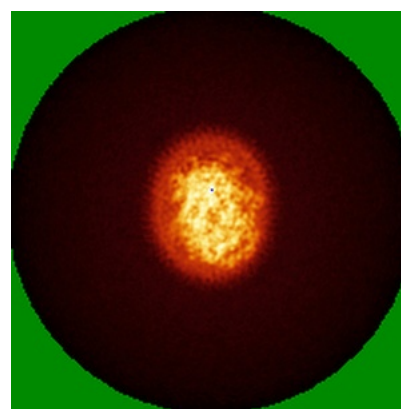
6.4.1 Primary map



X



Y

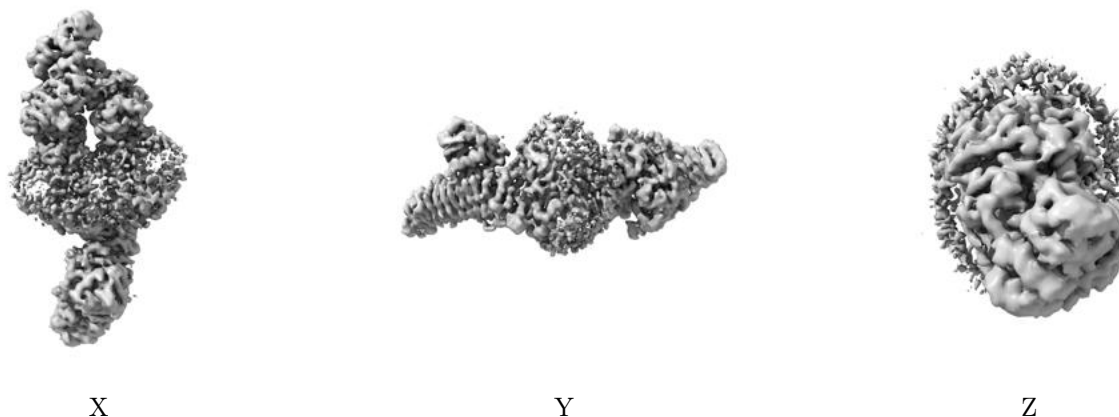


Z

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



The images above show the 3D surface view of the map at the recommended contour level 0.022. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

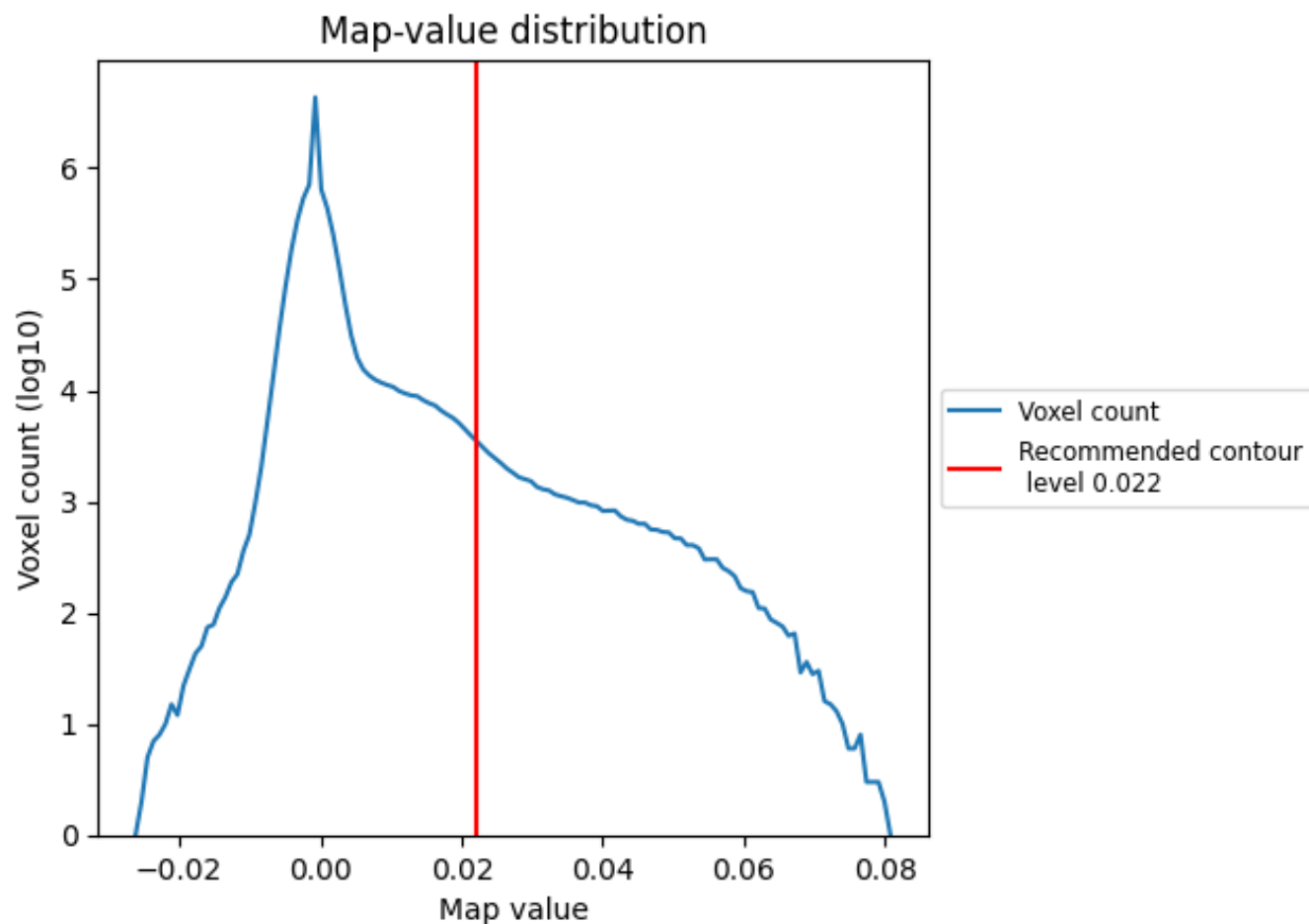
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

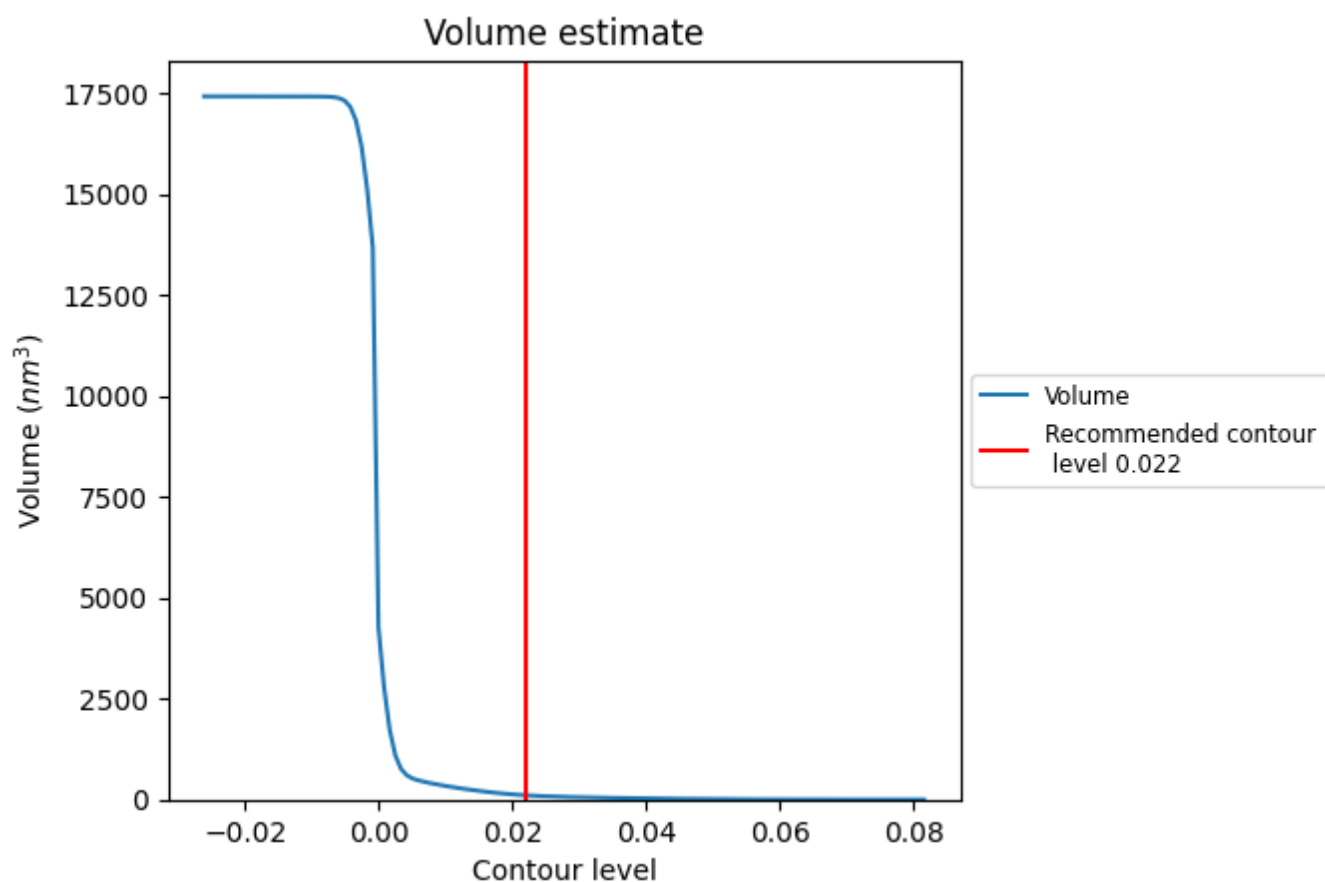
This section contains the results of statistical analysis of the map.

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.

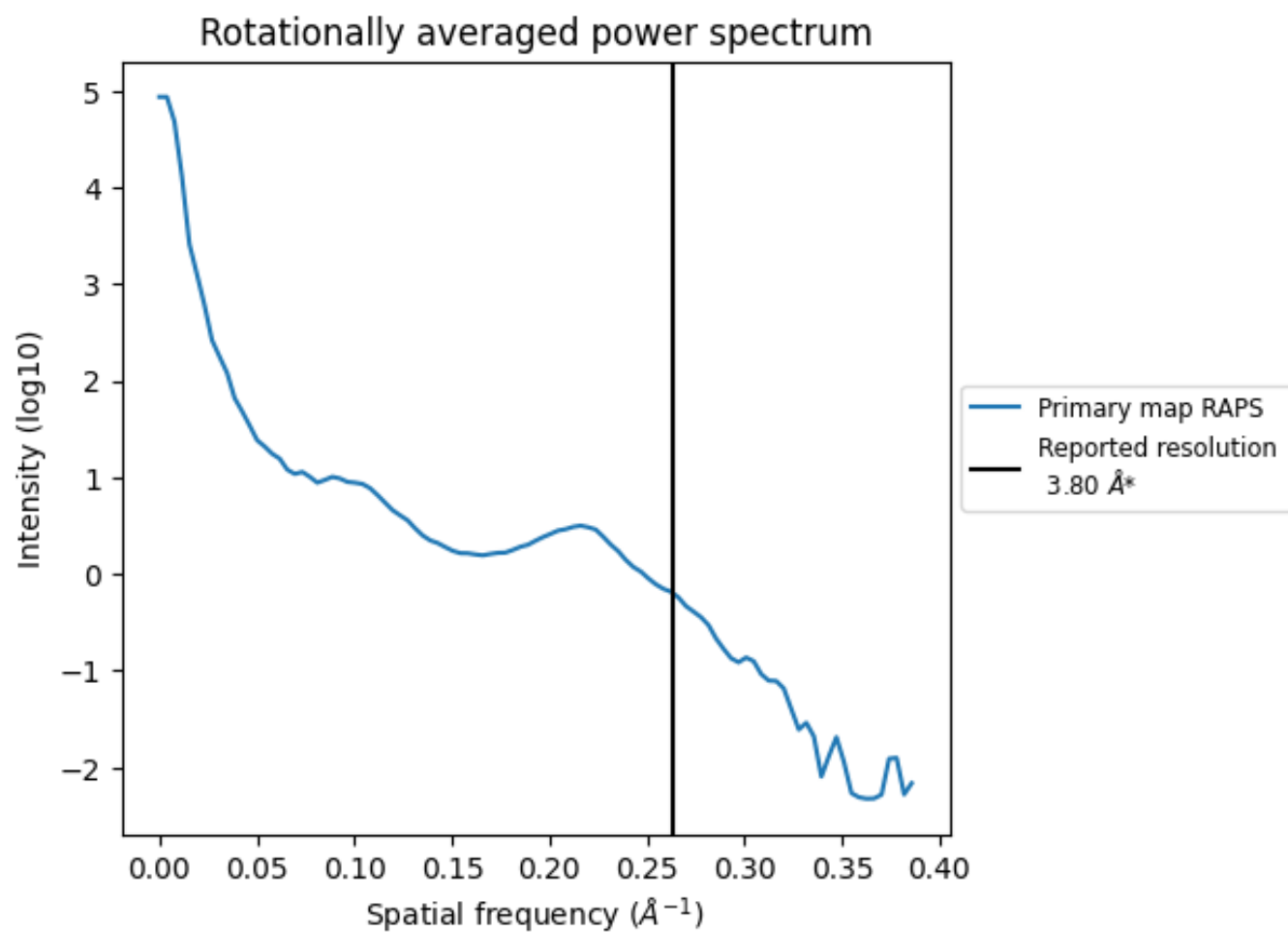
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 105 nm³; this corresponds to an approximate mass of 95 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 ⓘ

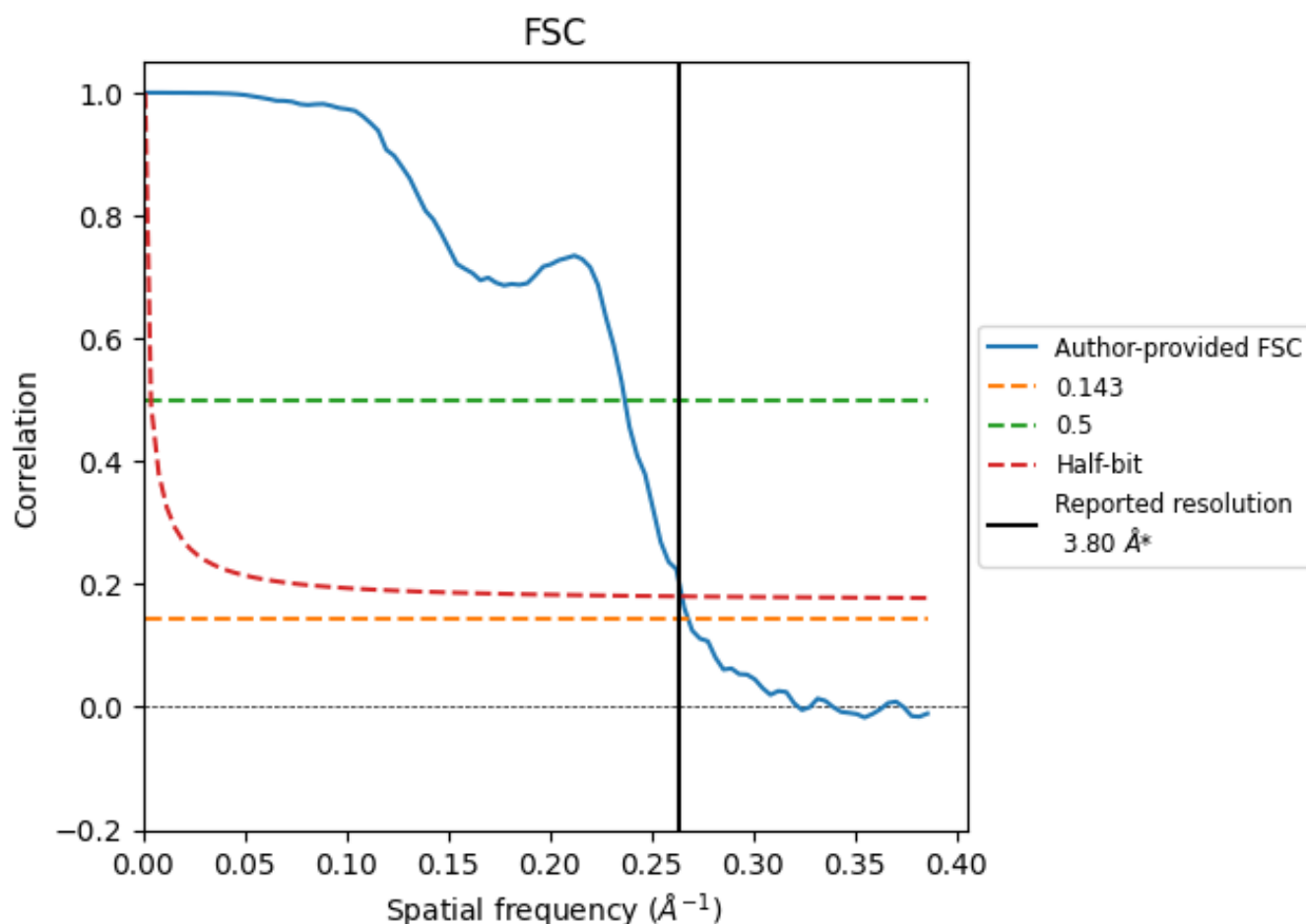


*Reported resolution corresponds to spatial frequency of 0.263 Å⁻¹

8 Fourier-Shell correlation [i](#)

Fourier-Shell Correlation (FSC) is the most commonly used method to estimate the resolution of single-particle and subtomogram-averaged maps. The shape of the curve depends on the imposed symmetry, mask and whether or not the two 3D reconstructions used were processed from a common reference. The reported resolution is shown as a black line. A curve is displayed for the half-bit criterion in addition to lines showing the 0.143 gold standard cut-off and 0.5 cut-off.

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.263 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

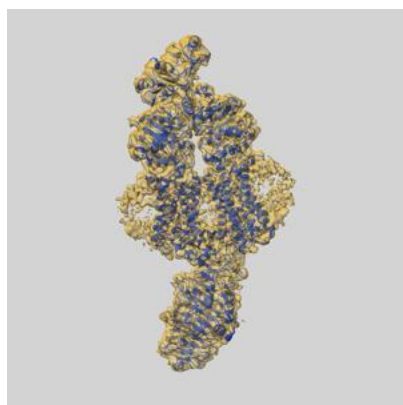
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.80	-	-
Author-provided FSC curve	3.73	4.22	3.77
Unmasked-calculated*	-	-	-

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

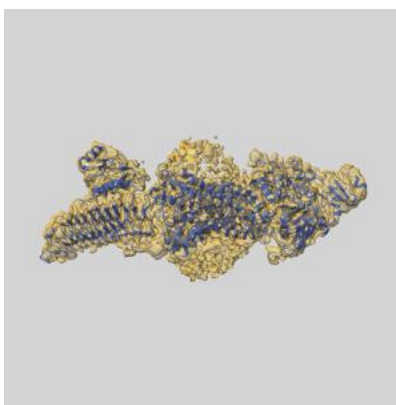
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-13049 and PDB model 7OSF. Per-residue inclusion information can be found in section [3](#) on page [6](#).

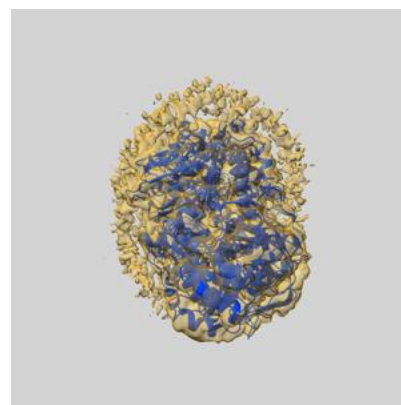
9.1 Map-model overlay [i](#)



X



Y



Z

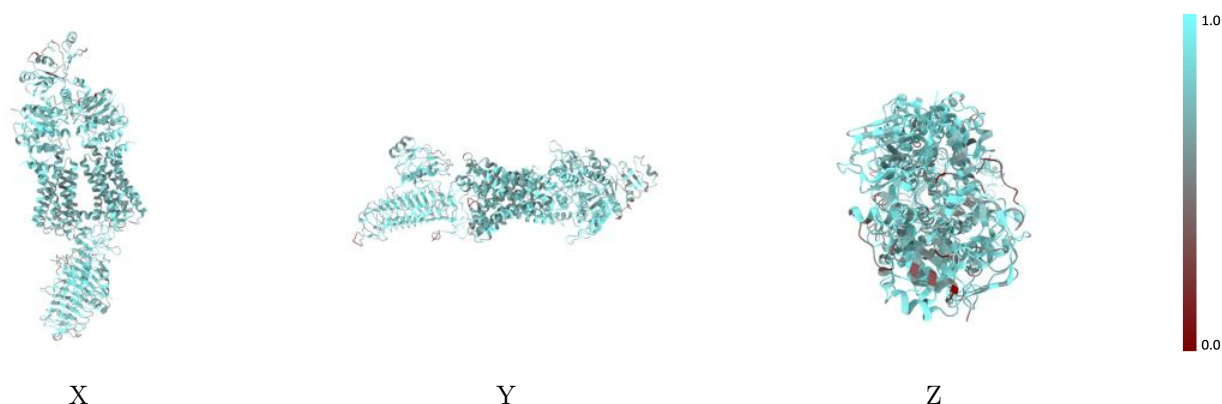
The images above show the 3D surface view of the map at the recommended contour level 0.022 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



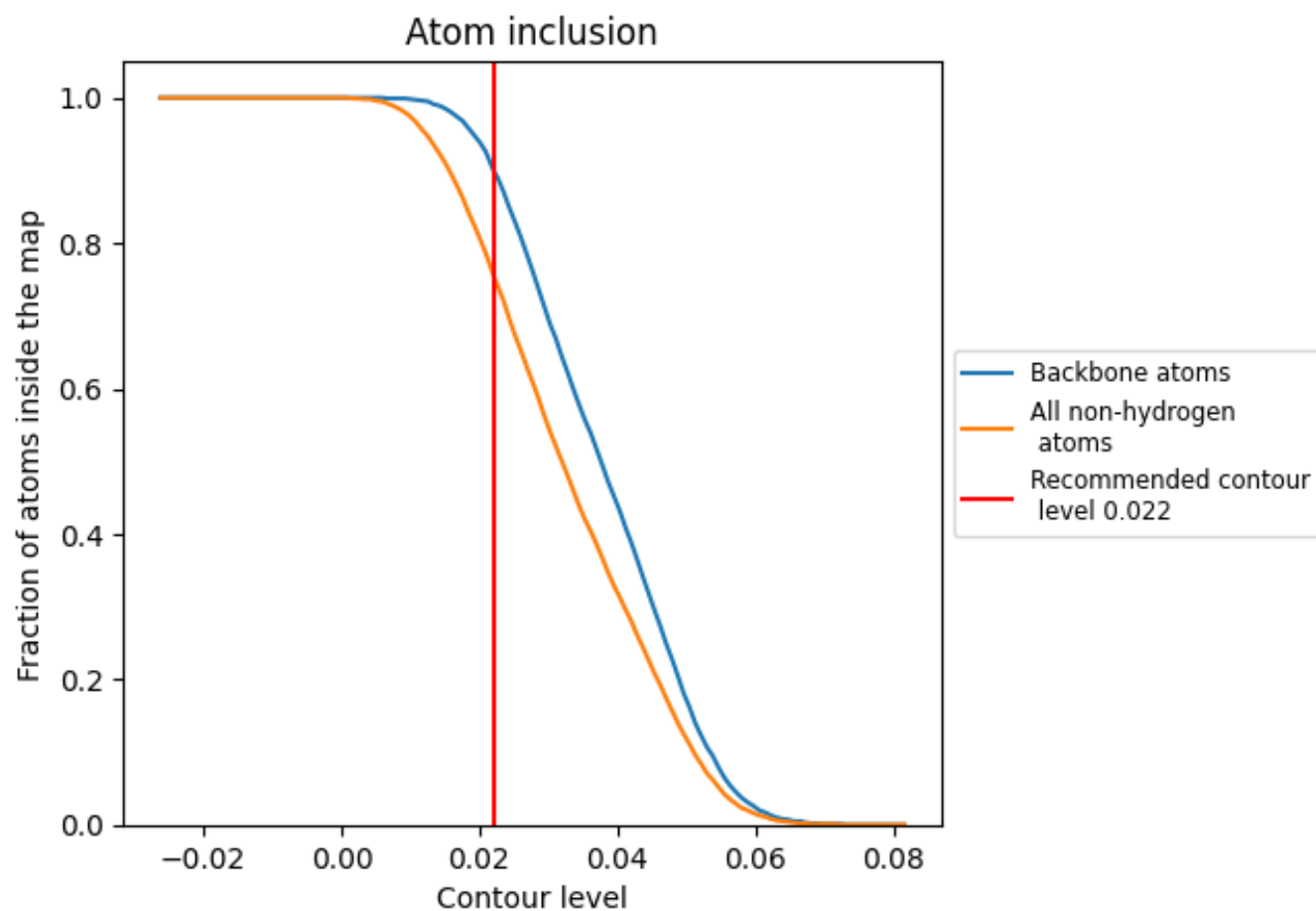
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.022).

9.4 Atom inclusion [i](#)



At the recommended contour level, 90% of all backbone atoms, 76% 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 (0.022) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7570	0.4310
A	0.7880	0.4350
B	0.7580	0.4120
C	0.7790	0.4280
D	0.7540	0.4560
E	0.7220	0.4290
H	0.7710	0.4210

1.0

0.0

<0.0