



wwPDB EM Validation Summary Report ⓘ

Sep 20, 2026 – 10:41 pm BST

PDB ID : 9SPE / pdb_00009spe
EMDB ID : EMD-55078
Title : Human tRNA ligase complex bound by PYROXD1
Authors : Pfeiderer, M.M.; Jinek, M.
Deposited on : 2025-09-16
Resolution : 3.25 Å(reported)

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

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

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

EMDB validation analysis : 0.0.1.dev133
Mogul : 1.8.4, CSD as541be (2020)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.50

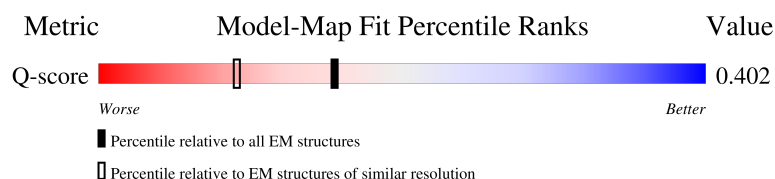
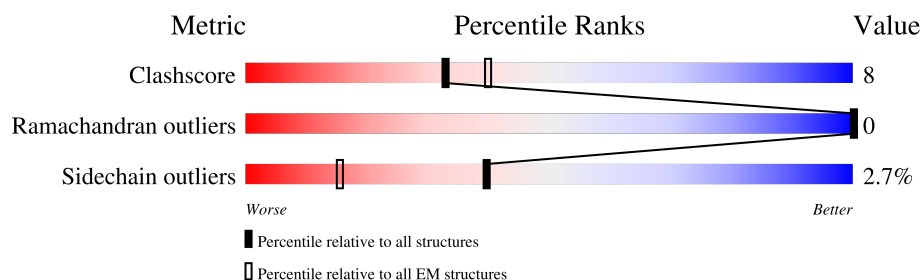
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.25 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.



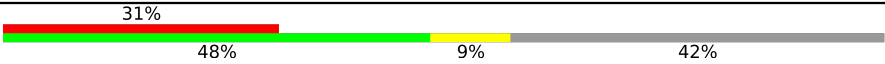
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	14599 (2.75 - 3.75)

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	259	16% 22% 75%
2	B	330	53% 62% 12% 25%
3	C	244	50% 62% 15% 22%
4	D	329	7% 90%

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Mol	Chain	Length	Quality of chain
5	P	731	
6	R	521	

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 11226 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 Ashwin.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	64	Total	C	N	O	S	0	0
			531	338	94	96	3		

There are 27 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-26	MET	-	initiating methionine	UNP Q9BVC5
A	-25	TRP	-	expression tag	UNP Q9BVC5
A	-24	SER	-	expression tag	UNP Q9BVC5
A	-23	HIS	-	expression tag	UNP Q9BVC5
A	-22	PRO	-	expression tag	UNP Q9BVC5
A	-21	GLN	-	expression tag	UNP Q9BVC5
A	-20	PHE	-	expression tag	UNP Q9BVC5
A	-19	GLU	-	expression tag	UNP Q9BVC5
A	-18	LYS	-	expression tag	UNP Q9BVC5
A	-17	GLY	-	expression tag	UNP Q9BVC5
A	-16	SER	-	expression tag	UNP Q9BVC5
A	-15	GLY	-	expression tag	UNP Q9BVC5
A	-14	TRP	-	expression tag	UNP Q9BVC5
A	-13	SER	-	expression tag	UNP Q9BVC5
A	-12	HIS	-	expression tag	UNP Q9BVC5
A	-11	PRO	-	expression tag	UNP Q9BVC5
A	-10	GLN	-	expression tag	UNP Q9BVC5
A	-9	PHE	-	expression tag	UNP Q9BVC5
A	-8	GLU	-	expression tag	UNP Q9BVC5
A	-7	LYS	-	expression tag	UNP Q9BVC5
A	-6	PRO	-	expression tag	UNP Q9BVC5
A	-5	PRO	-	expression tag	UNP Q9BVC5
A	-4	SER	-	expression tag	UNP Q9BVC5
A	-3	GLY	-	expression tag	UNP Q9BVC5
A	-2	ALA	-	expression tag	UNP Q9BVC5
A	-1	ASP	-	expression tag	UNP Q9BVC5
A	0	PRO	-	expression tag	UNP Q9BVC5

- Molecule 2 is a protein called Protein FAM98B.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	248	Total	C	N	O	S	0	0
			1948	1229	325	380	14		

- Molecule 3 is a protein called RNA transcription, translation and transport factor protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	191	Total	C	N	O	S	0	0
			1569	1003	281	280	5		

- Molecule 4 is a protein called ATP-dependent RNA helicase DDX1, Green fluorescent protein.

Mol	Chain	Residues	Atoms				AltConf	Trace
4	D	32	Total	C	N	O	0	0
			248	161	39	48		

There are 53 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	696	MET	-	initiating methionine	UNP Q92499
D	741	SER	-	linker	UNP Q92499
D	742	PHE	-	linker	UNP Q92499
D	743	GLY	-	linker	UNP Q92499
D	744	SER	-	linker	UNP Q92499
D	745	GLY	-	linker	UNP Q92499
D	746	ALA	-	linker	UNP Q92499
D	747	THR	-	linker	UNP Q92499
D	748	ASN	-	linker	UNP Q92499
D	749	PHE	-	linker	UNP Q92499
D	750	SER	-	linker	UNP Q92499
D	751	LEU	-	linker	UNP Q92499
D	752	LEU	-	linker	UNP Q92499
D	753	LYS	-	linker	UNP Q92499
D	754	GLN	-	linker	UNP Q92499
D	755	ALA	-	linker	UNP Q92499
D	756	GLY	-	linker	UNP Q92499
D	757	ASP	-	linker	UNP Q92499
D	758	VAL	-	linker	UNP Q92499
D	759	GLU	-	linker	UNP Q92499
D	760	GLU	-	linker	UNP Q92499
D	761	ASN	-	linker	UNP Q92499
D	762	PRO	-	linker	UNP Q92499

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Chain	Residue	Modelled	Actual	Comment	Reference
D	763	GLY	-	linker	UNP Q92499
D	764	PRO	-	linker	UNP Q92499
D	765	GLY	-	linker	UNP Q92499
D	766	SER	-	linker	UNP Q92499
D	767	GLY	-	linker	UNP Q92499
D	768	GLU	-	linker	UNP Q92499
D	769	GLY	-	linker	UNP Q92499
D	770	ARG	-	linker	UNP Q92499
D	771	GLY	-	linker	UNP Q92499
D	772	SER	-	linker	UNP Q92499
D	773	LEU	-	linker	UNP Q92499
D	774	LEU	-	linker	UNP Q92499
D	775	THR	-	linker	UNP Q92499
D	776	CYS	-	linker	UNP Q92499
D	777	GLY	-	linker	UNP Q92499
D	778	ASP	-	linker	UNP Q92499
D	779	VAL	-	linker	UNP Q92499
D	780	GLU	-	linker	UNP Q92499
D	781	GLU	-	linker	UNP Q92499
D	782	ASN	-	linker	UNP Q92499
D	783	PRO	-	linker	UNP Q92499
D	784	GLY	-	linker	UNP Q92499
D	785	PRO	-	linker	UNP Q92499
D	786	MET	-	linker	UNP Q92499
D	787	VAL	-	linker	UNP Q92499
D	850	LEU	PHE	engineered mutation	UNP P42212
D	851	THR	SER	engineered mutation	UNP P42212
D	943	HIS	GLN	engineered mutation	UNP P42212
D	992	LYS	ALA	engineered mutation	UNP P42212
D	1017	LEU	HIS	engineered mutation	UNP P42212

- Molecule 5 is a protein called Glutathione S-transferase class-mu 26 kDa isozyme, tRNA ligase complex-associated NAD(P)H dehydrogenase PYROXD1.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	P	422	Total	C	N	O	S	0	0
			3310	2114	544	628	24		

There are 13 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
P	-12	SER	-	linker	UNP P08515

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Chain	Residue	Modelled	Actual	Comment	Reference
P	-11	ASP	-	linker	UNP P08515
P	-10	LEU	-	linker	UNP P08515
P	-9	GLU	-	linker	UNP P08515
P	-8	VAL	-	linker	UNP P08515
P	-7	LEU	-	linker	UNP P08515
P	-6	PHE	-	linker	UNP P08515
P	-5	GLN	-	linker	UNP P08515
P	-4	GLY	-	linker	UNP P08515
P	-3	PRO	-	linker	UNP P08515
P	-2	LEU	-	linker	UNP P08515
P	-1	GLY	-	linker	UNP P08515
P	0	SER	-	linker	UNP P08515

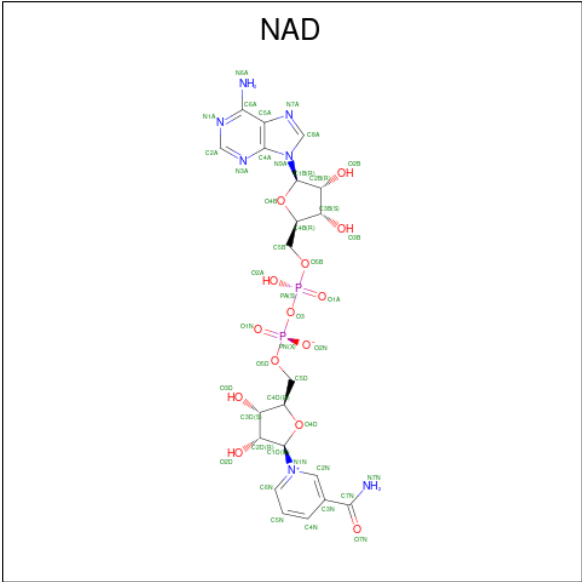
- Molecule 6 is a protein called RNA-splicing ligase RtcB homolog.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	R	460	Total	C	N	O	S	0	0
			3521	2210	628	657	26		

There are 16 discrepancies between the modelled and reference sequences:

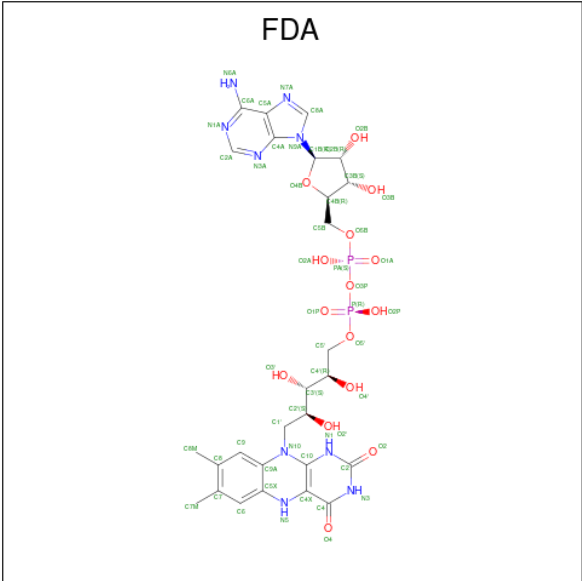
Chain	Residue	Modelled	Actual	Comment	Reference
R	-15	MET	-	initiating methionine	UNP Q9Y3I0
R	-14	HIS	-	expression tag	UNP Q9Y3I0
R	-13	HIS	-	expression tag	UNP Q9Y3I0
R	-12	HIS	-	expression tag	UNP Q9Y3I0
R	-11	HIS	-	expression tag	UNP Q9Y3I0
R	-10	HIS	-	expression tag	UNP Q9Y3I0
R	-9	HIS	-	expression tag	UNP Q9Y3I0
R	-8	HIS	-	expression tag	UNP Q9Y3I0
R	-7	HIS	-	expression tag	UNP Q9Y3I0
R	-6	PRO	-	expression tag	UNP Q9Y3I0
R	-5	PRO	-	expression tag	UNP Q9Y3I0
R	-4	SER	-	expression tag	UNP Q9Y3I0
R	-3	GLY	-	expression tag	UNP Q9Y3I0
R	-2	ALA	-	expression tag	UNP Q9Y3I0
R	-1	ASP	-	expression tag	UNP Q9Y3I0
R	0	PRO	-	expression tag	UNP Q9Y3I0

- Molecule 7 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (CCD ID: NAD) (formula: $C_{21}H_{27}N_7O_{14}P_2$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					AltConf
7	P	1	Total	C	N	O	P	0
			44	21	7	14	2	

- Molecule 8 is DIHYDROFLAVINE-ADENINE DINUCLEOTIDE (CCD ID: FDA) (formula: $C_{27}H_{35}N_9O_{15}P_2$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					AltConf
8	P	1	Total	C	N	O	P	0
			53	27	9	15	2	

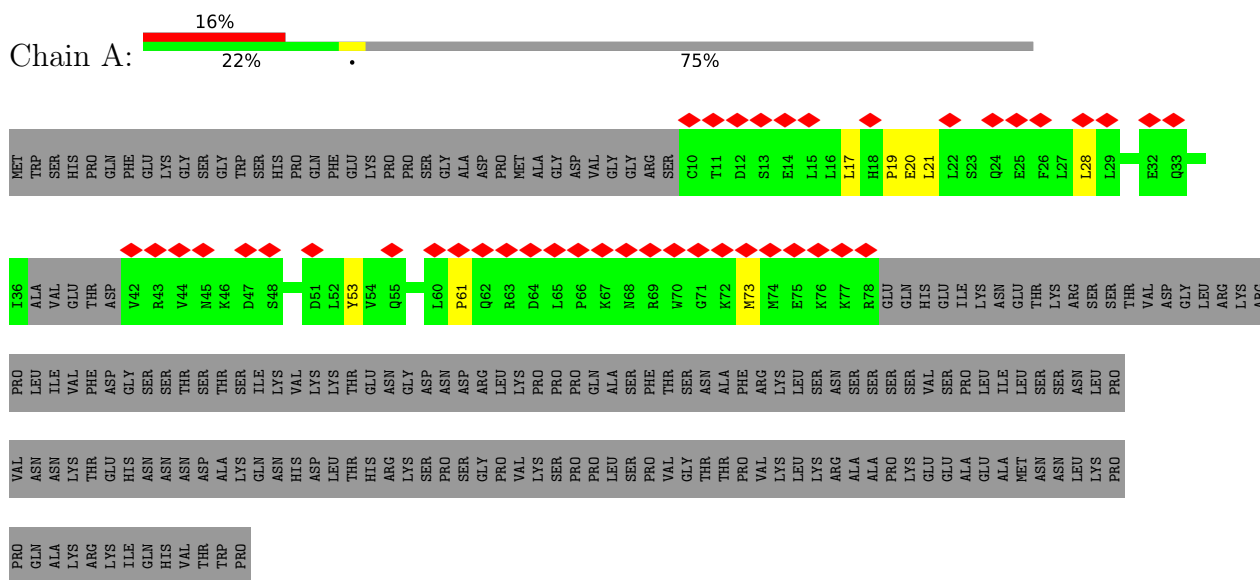
- Molecule 9 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		AltConf
9	R	2	Total	Mg	0
			2	2	

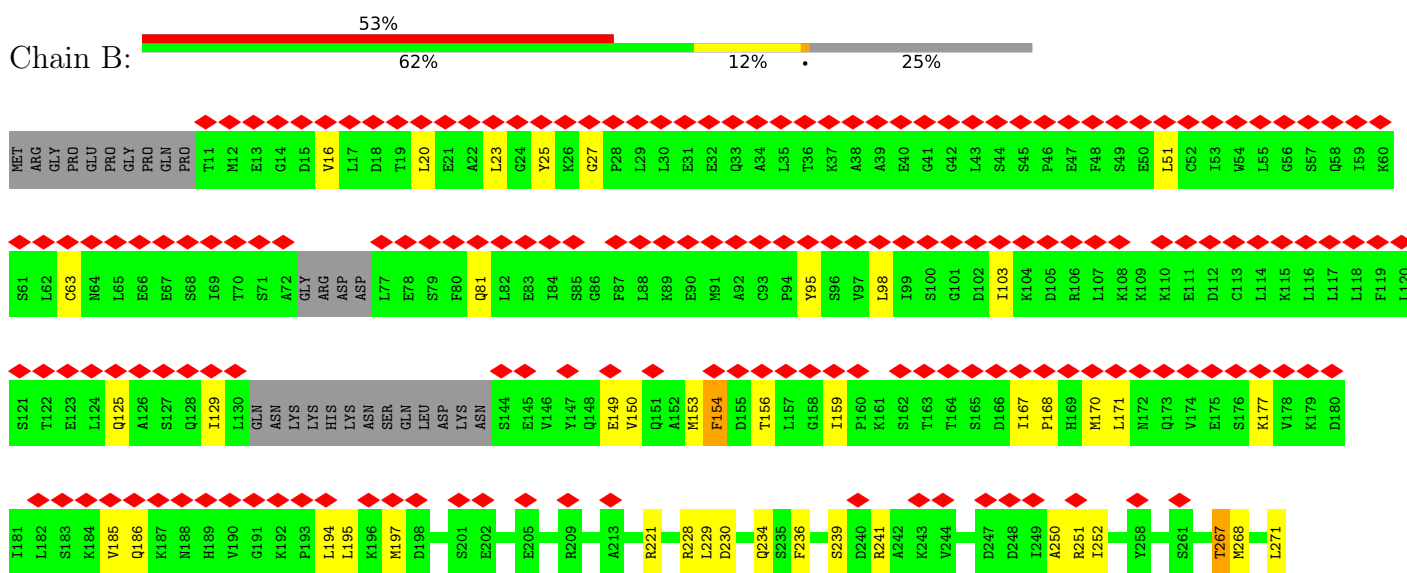
3 Residue-property plots

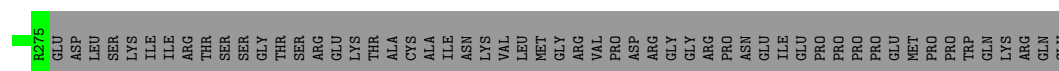
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion 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: Ashwin

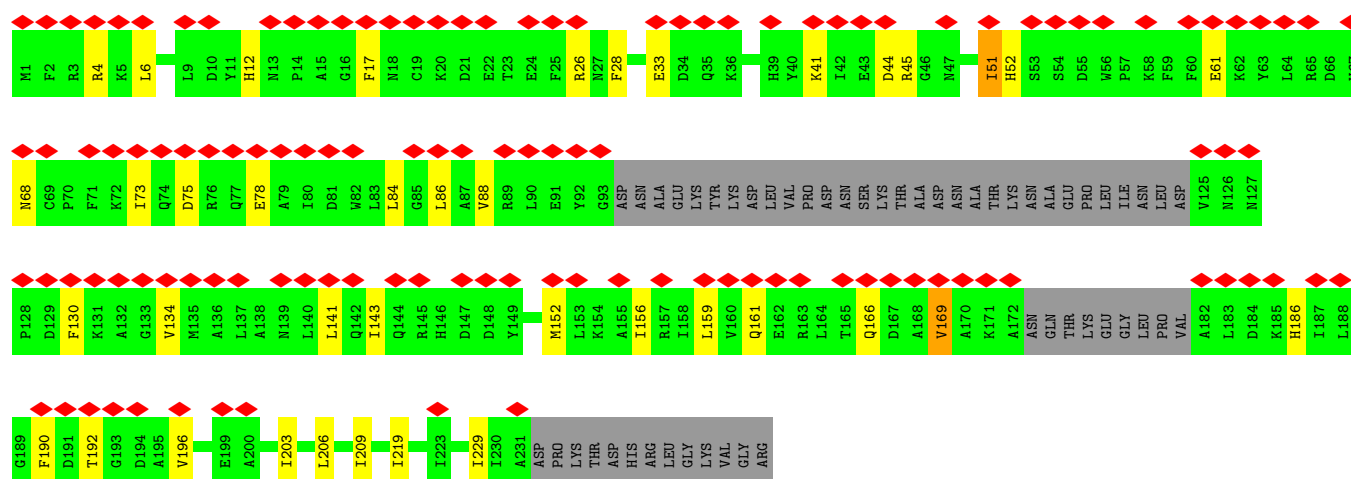


• Molecule 2: Protein FAM98B

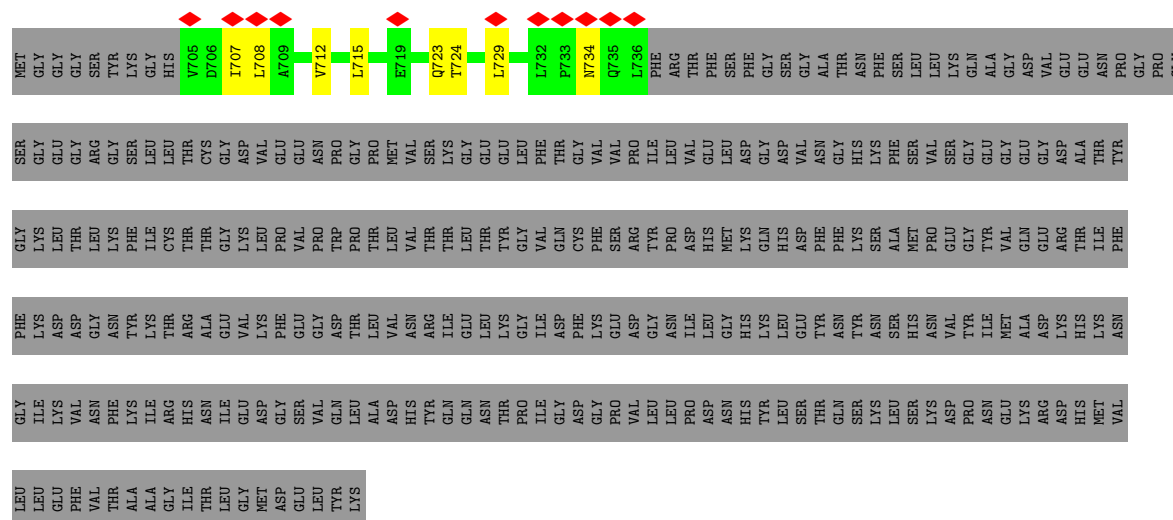




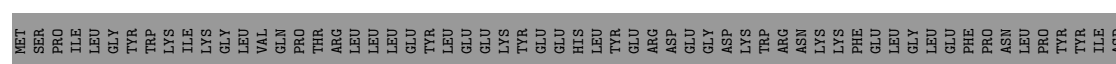
• Molecule 3: RNA transcription, translation and transport factor protein

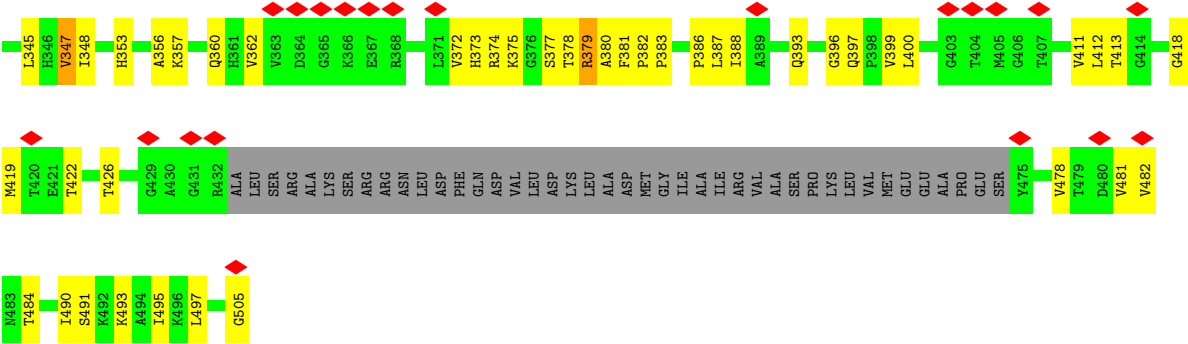


• Molecule 4: ATP-dependent RNA helicase DDX1, Green fluorescent protein



• Molecule 5: Glutathione S-transferase class-mu 26 kDa isozyme, tRNA ligase complex-associated NAD(P)H dehydrogenase PYROXD1





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	269780	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	60	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2600	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	0.520	Depositor
Minimum map value	-0.372	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.015	Depositor
Recommended contour level	0.1	Depositor
Map size (Å)	233.99998, 233.99998, 233.99998	wwPDB
Map dimensions	360, 360, 360	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.65, 0.65, 0.65	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: MG, NAD, FDA

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	0.12	0/539	0.31	0/725
2	B	0.13	0/1972	0.30	0/2651
3	C	0.14	0/1596	0.33	0/2152
4	D	0.16	0/252	0.37	0/344
5	P	0.14	0/3377	0.34	0/4553
6	R	0.18	0/3589	0.39	0/4845
All	All	0.15	0/11325	0.35	0/15270

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	531	0	553	6	0
2	B	1948	0	1997	32	0
3	C	1569	0	1585	29	0
4	D	248	0	251	5	0
5	P	3310	0	3261	50	0
6	R	3521	0	3490	89	0
7	P	44	0	26	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
8	P	53	0	33	1	0
9	R	2	0	0	0	0
All	All	11226	0	11196	189	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.

The worst 5 of 189 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)
6:R:374:ARG:HB2	6:R:377:SER:HB2	1.61	0.82
6:R:129:THR:HG22	6:R:131:LEU:H	1.50	0.76
3:C:61:GLU:HB2	3:C:73:ILE:HD11	1.66	0.75
1:A:17:LEU:HD13	2:B:250:ALA:HB2	1.74	0.69
5:P:420:VAL:HG21	5:P:468:GLU:HG2	1.76	0.68

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	60/259 (23%)	58 (97%)	2 (3%)	0	100	100
2	B	242/330 (73%)	237 (98%)	5 (2%)	0	100	100
3	C	185/244 (76%)	181 (98%)	4 (2%)	0	100	100
4	D	30/329 (9%)	29 (97%)	1 (3%)	0	100	100
5	P	414/731 (57%)	401 (97%)	13 (3%)	0	100	100
6	R	456/521 (88%)	443 (97%)	13 (3%)	0	100	100
All	All	1387/2414 (58%)	1349 (97%)	38 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	62/235 (26%)	61 (98%)	1 (2%)	55	70
2	B	221/294 (75%)	219 (99%)	2 (1%)	70	76
3	C	165/211 (78%)	159 (96%)	6 (4%)	31	56
4	D	27/279 (10%)	23 (85%)	4 (15%)	3	13
5	P	357/629 (57%)	351 (98%)	6 (2%)	53	69
6	R	370/422 (88%)	357 (96%)	13 (4%)	32	56
All	All	1202/2070 (58%)	1170 (97%)	32 (3%)	40	61

5 of 32 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
6	R	379	ARG
6	R	413	THR
4	D	734	ASN
4	D	724	THR
6	R	478	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 12 such sidechains are listed below:

Mol	Chain	Res	Type
6	R	232	GLN
6	R	250	HIS
6	R	428	HIS
6	R	289	ASN
3	C	186	HIS

5.3.3 RNA ⓘ

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 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
8	FDA	P	602	-	55,58,58	0.72	1 (1%)	73,89,89	0.82	3 (4%)
7	NAD	P	601	-	45,48,48	0.72	1 (2%)	63,73,73	0.84	2 (3%)

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
8	FDA	P	602	-	-	12/34/50/50	0/6/6/6
7	NAD	P	601	-	-	5/30/62/62	0/5/5/5

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	P	601	NAD	O4D-C1D	-2.20	1.38	1.41
8	P	602	FDA	C4X-N5	2.10	1.40	1.35

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	P	602	FDA	P-O3P-PA	-2.76	123.35	132.83
7	P	601	NAD	O7N-C7N-N7N	2.32	125.87	122.58
7	P	601	NAD	O2A-PA-O1A	2.14	122.80	112.24
8	P	602	FDA	O2'-C2'-C3'	2.13	114.28	109.10
8	P	602	FDA	O3'-C3'-C4'	2.02	113.68	108.81

There are no chirality outliers.

5 of 17 torsion outliers are listed below:

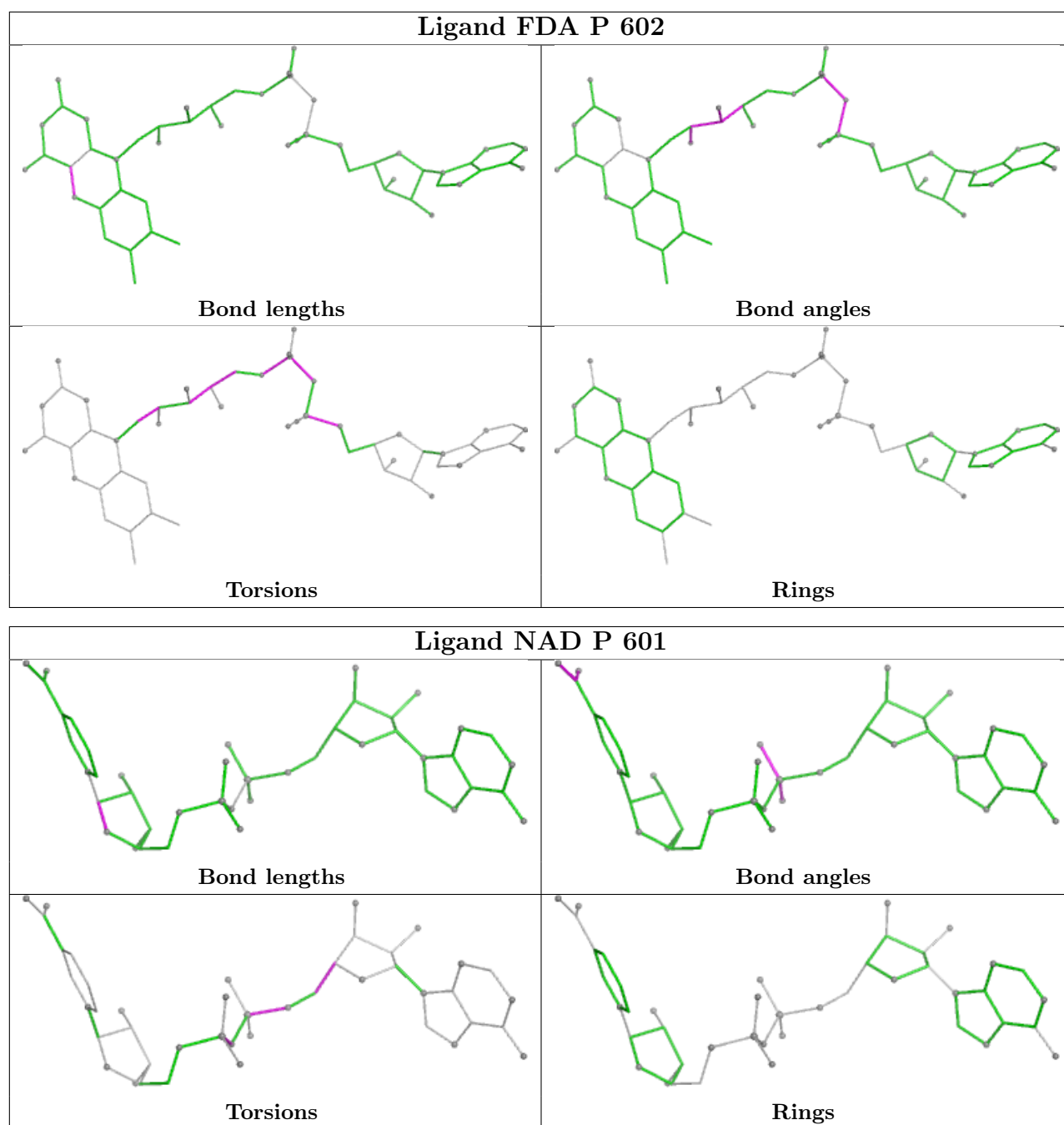
Mol	Chain	Res	Type	Atoms
8	P	602	FDA	C5B-O5B-PA-O3P
8	P	602	FDA	N10-C1'-C2'-O2'
8	P	602	FDA	C3'-C4'-C5'-O5'
8	P	602	FDA	O4'-C4'-C5'-O5'
8	P	602	FDA	C5'-O5'-P-O3P

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	P	602	FDA	1	0
7	P	601	NAD	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

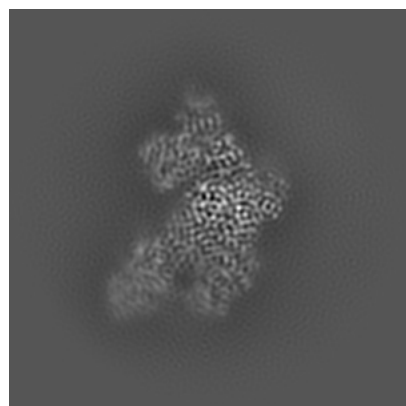
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-55078. These allow visual inspection of the internal detail of the map and identification of artifacts.

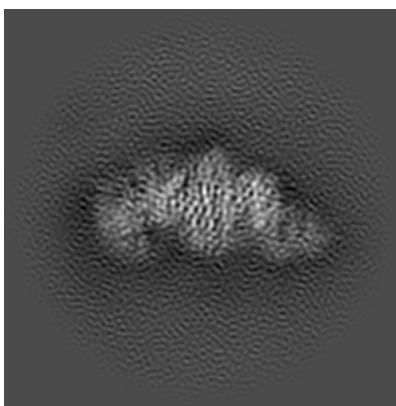
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

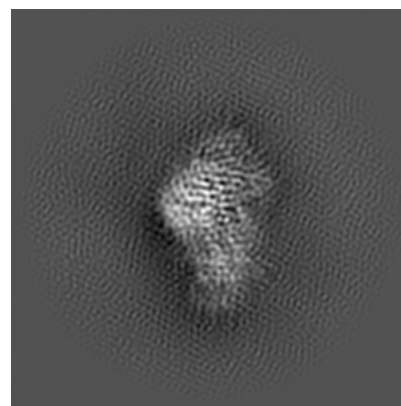
6.1.1 Primary map



X

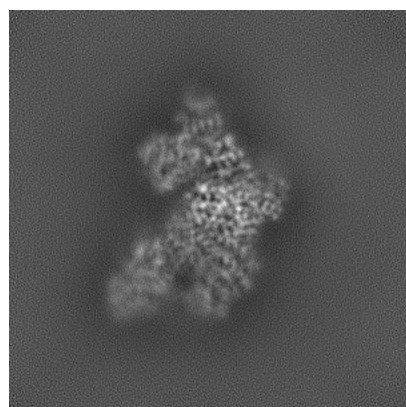


Y

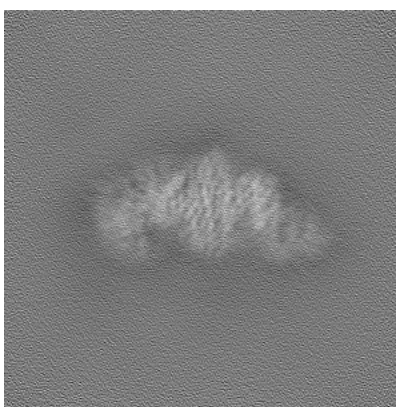


Z

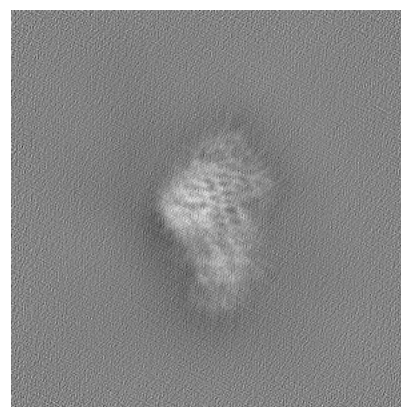
6.1.2 Raw map



X



Y

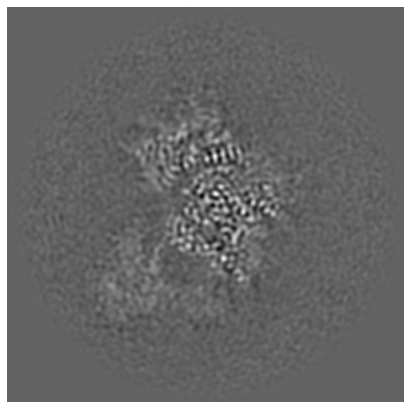


Z

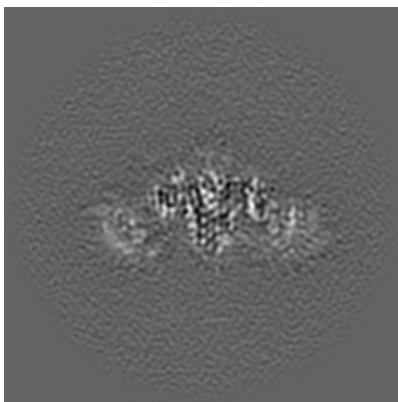
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

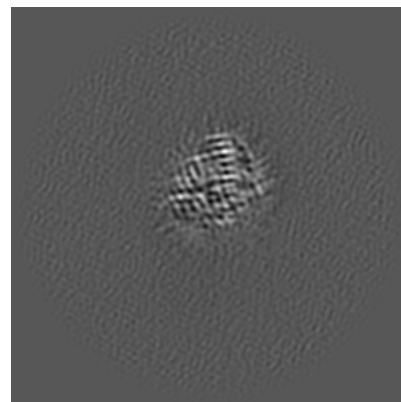
6.2.1 Primary map



X Index: 180

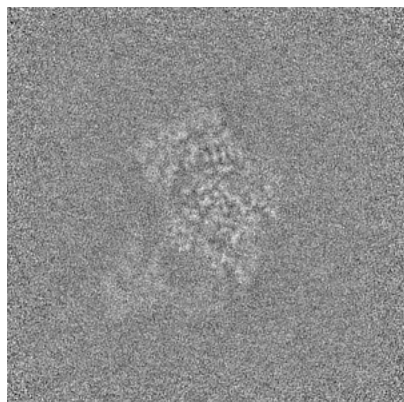


Y Index: 180

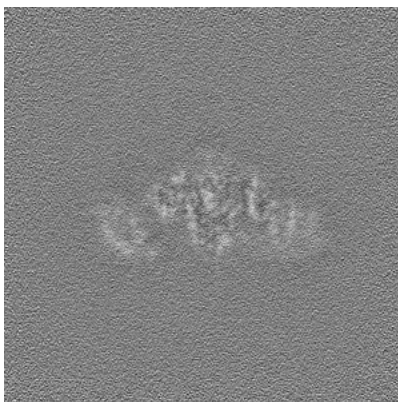


Z Index: 180

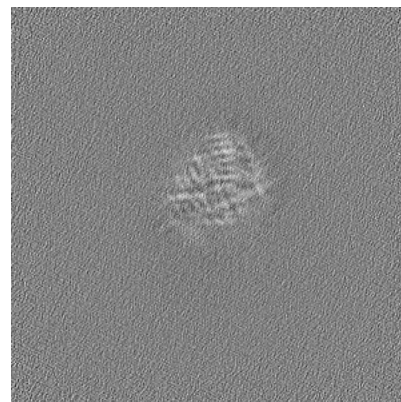
6.2.2 Raw map



X Index: 180



Y Index: 180

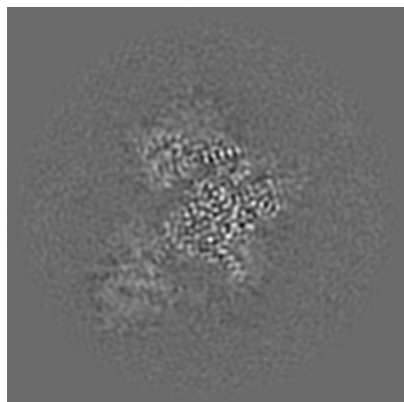


Z Index: 180

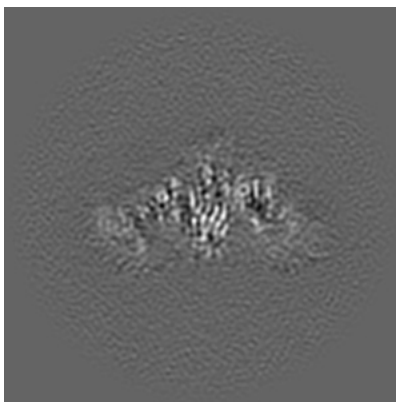
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

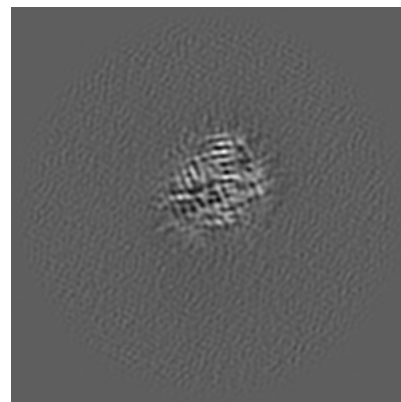
6.3.1 Primary map



X Index: 185

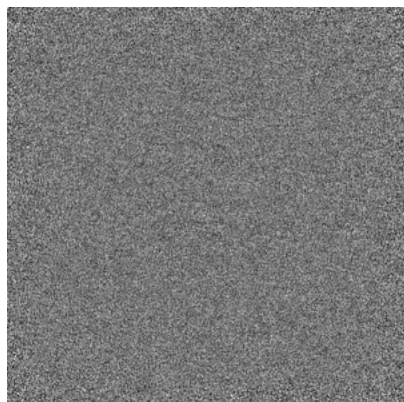


Y Index: 190

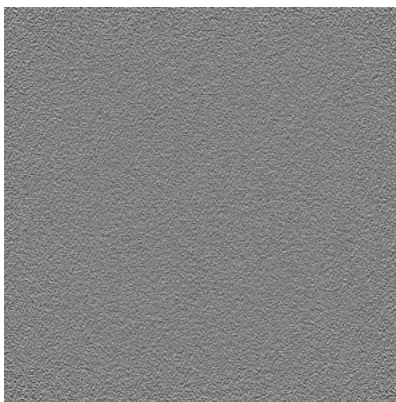


Z Index: 181

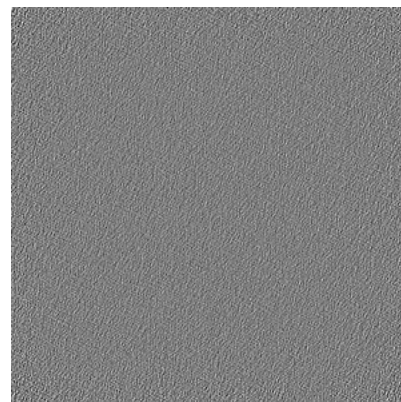
6.3.2 Raw map



X Index: 0



Y Index: 0

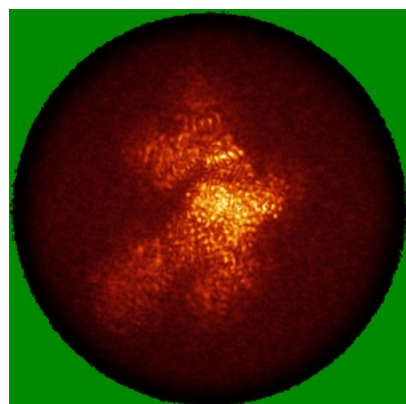


Z Index: 359

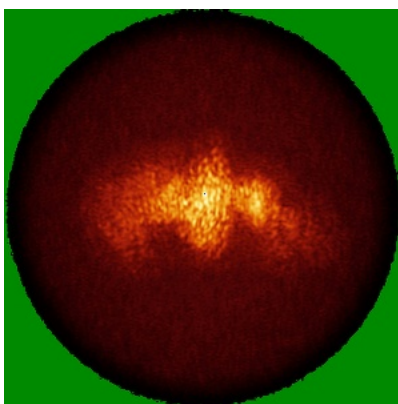
The images above show the largest variance slices of the map in three orthogonal directions.

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

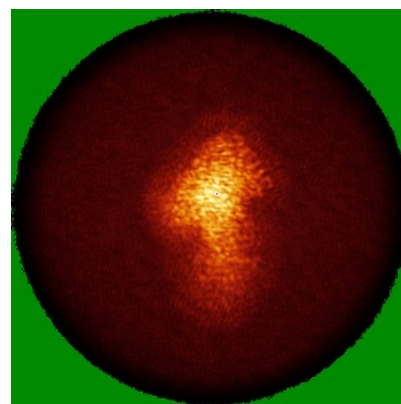
6.4.1 Primary map



X

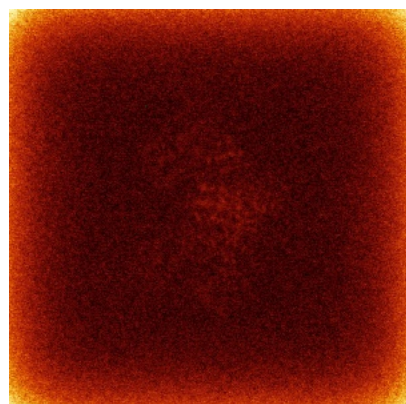


Y

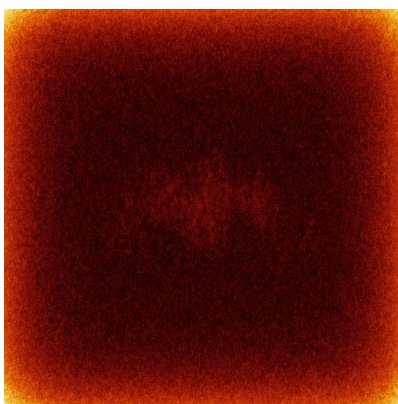


Z

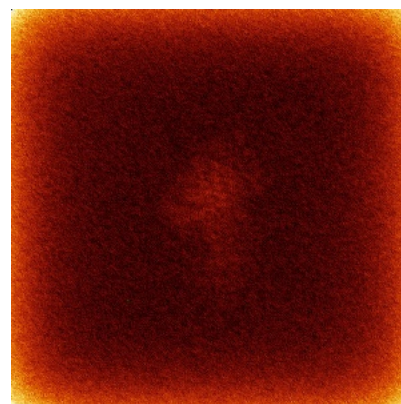
6.4.2 Raw map



X



Y



Z

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

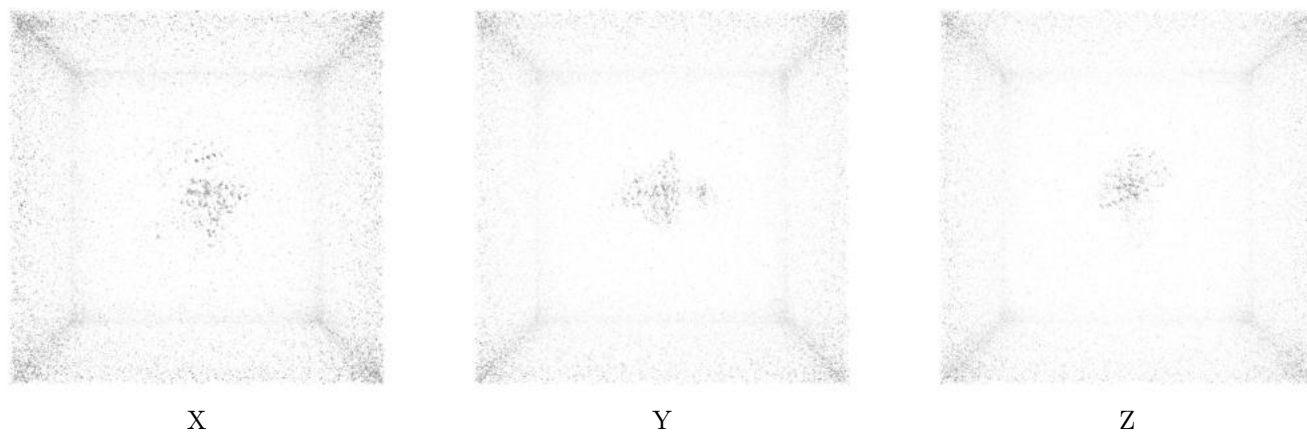
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



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

6.5.2 Raw map



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

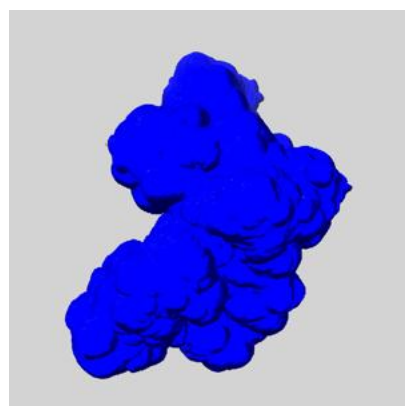
6.6 Mask visualisation [i](#)

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

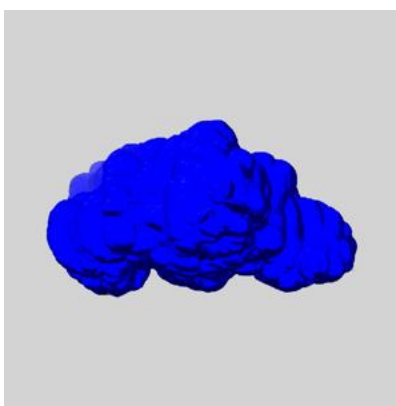
A mask typically either:

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

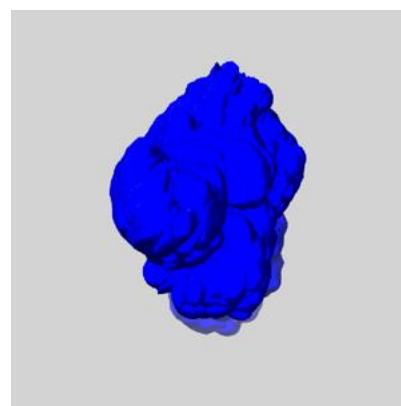
6.6.1 emd_55078_msk_1.map [i](#)



X



Y

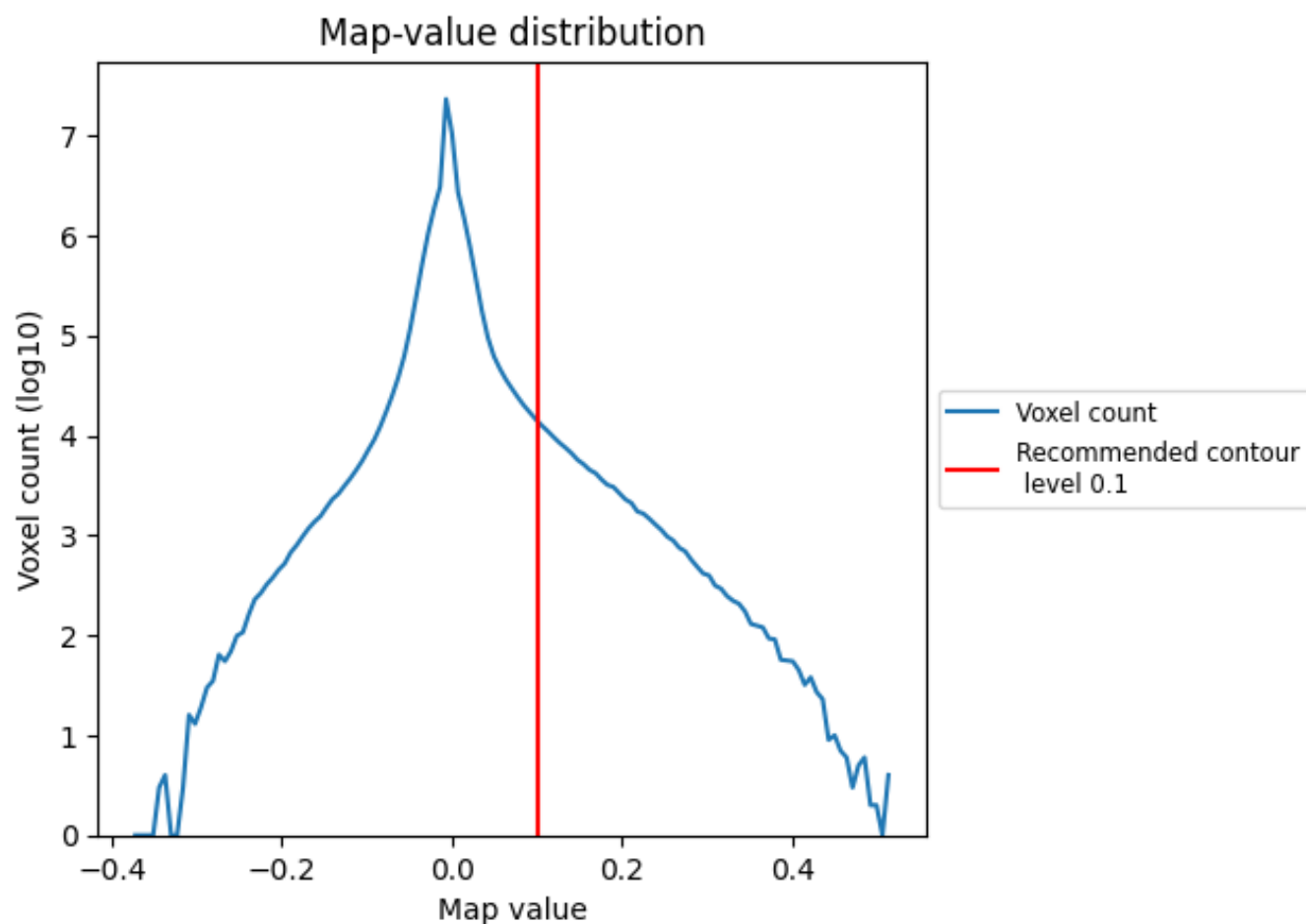


Z

7 Map analysis [i](#)

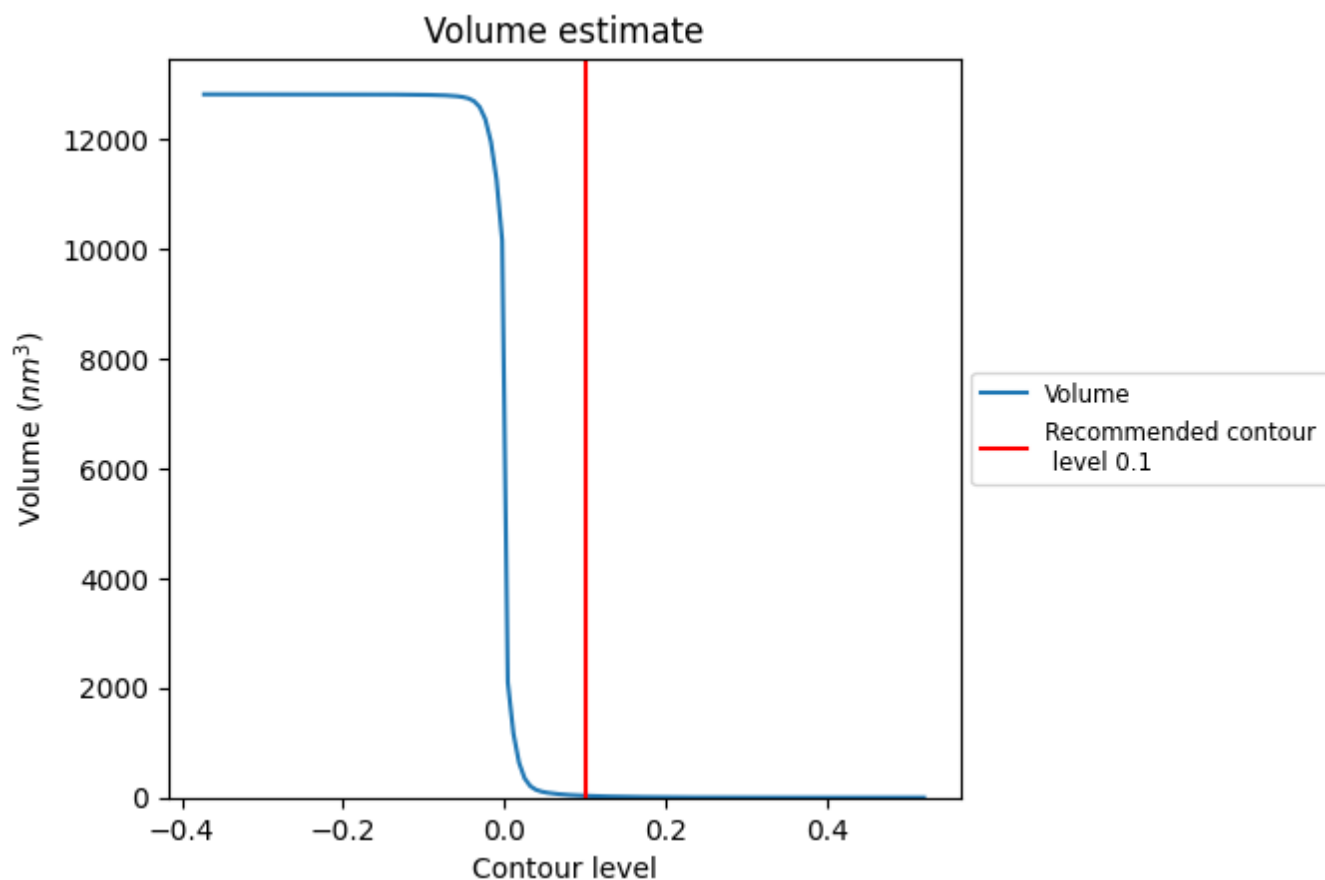
This section contains the results of statistical analysis of the map.

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.

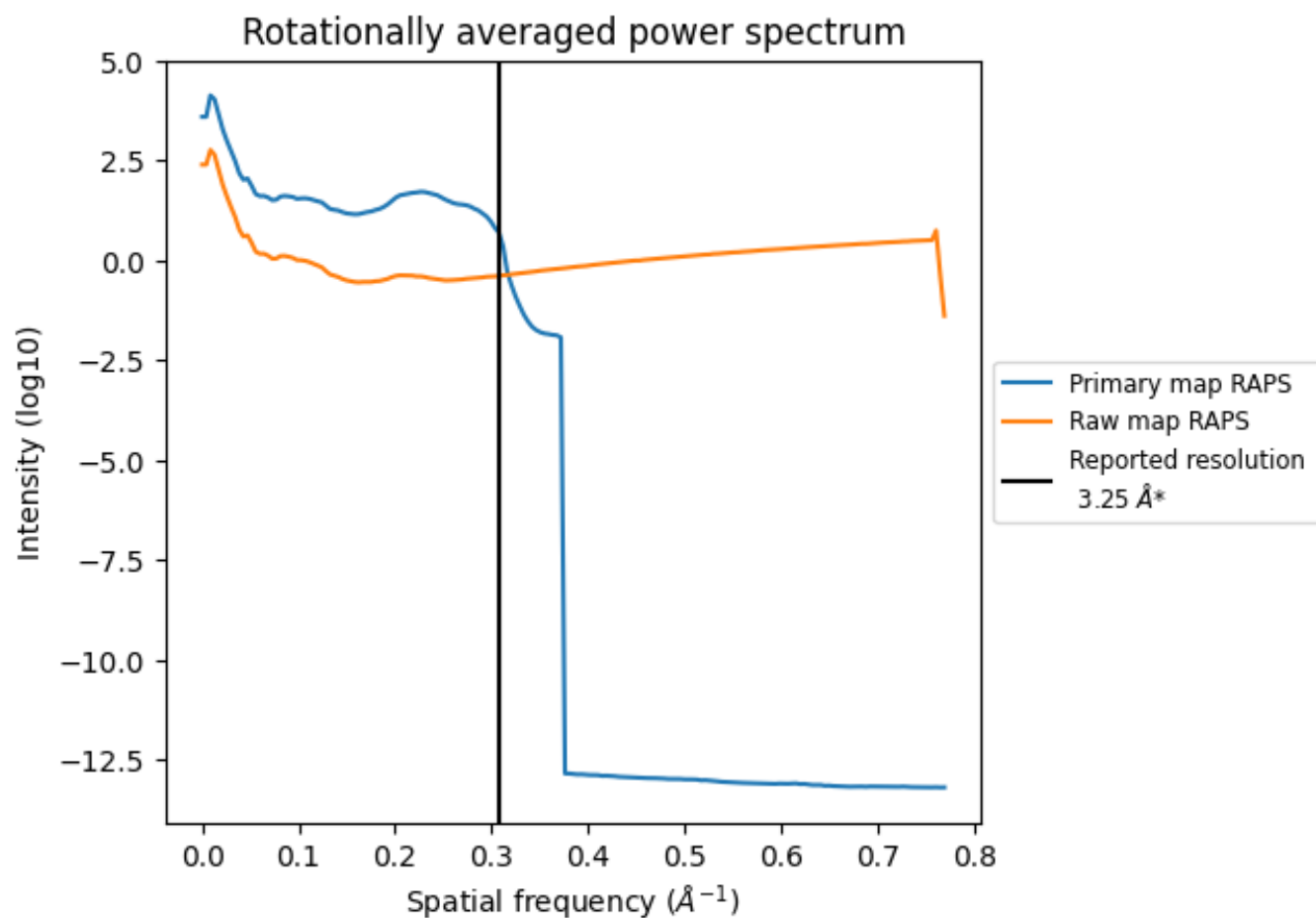
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 33 nm^3 ; this corresponds to an approximate mass of 29 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 ⓘ

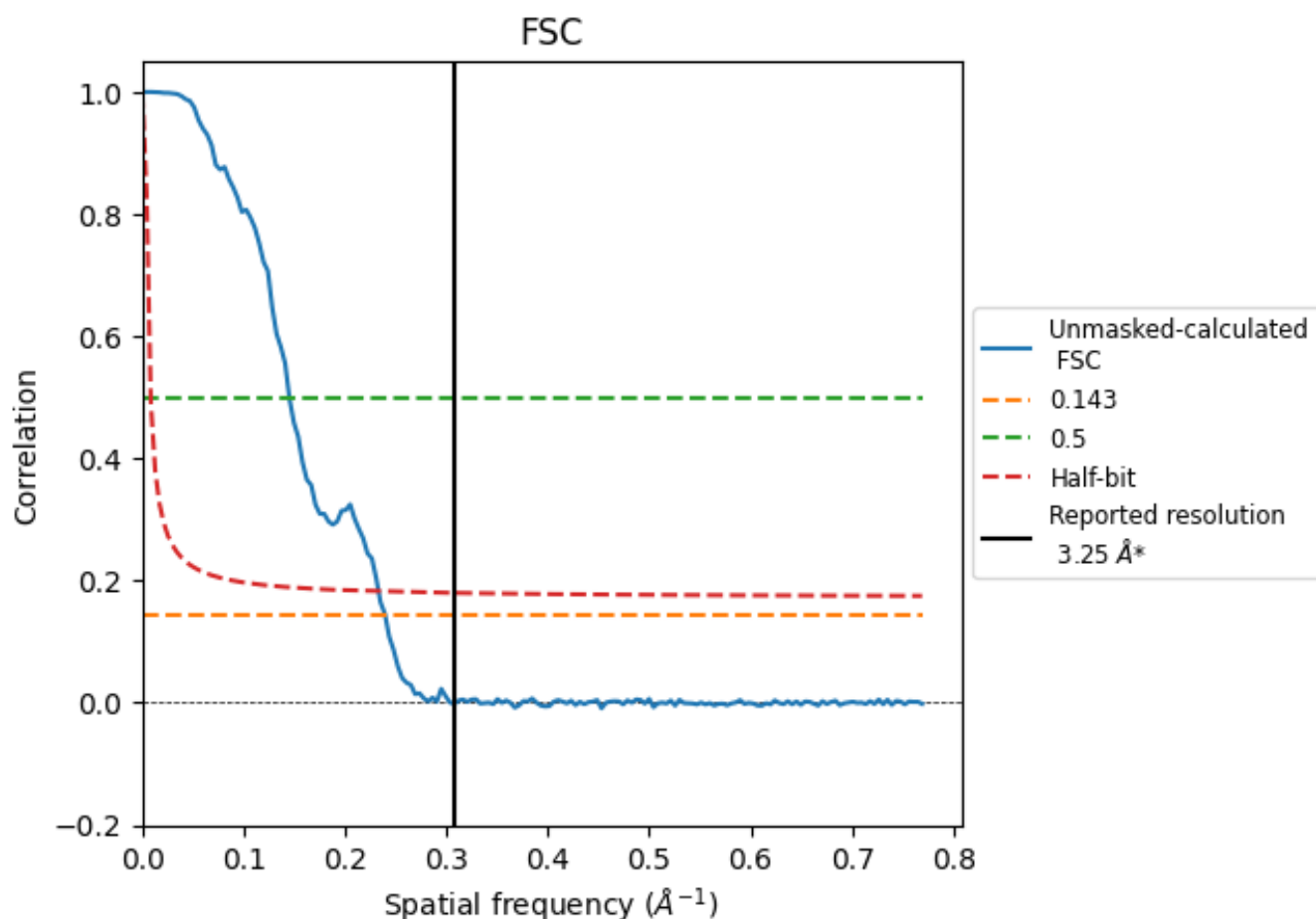


*Reported resolution corresponds to spatial frequency of 0.308 Å⁻¹

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 0.308 Å⁻¹

8.2 Resolution estimates [i](#)

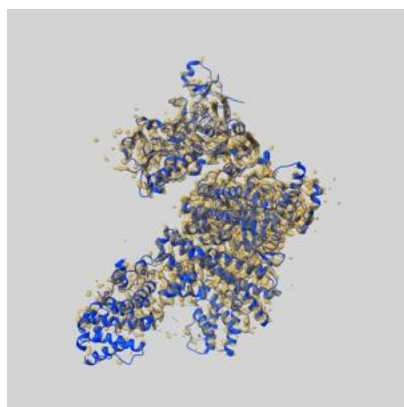
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.25	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	4.17	6.89	4.29

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 4.17 differs from the reported value 3.25 by more than 10 %

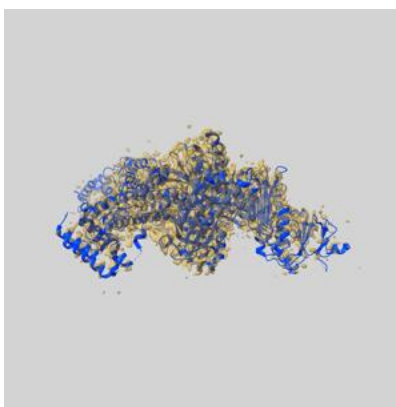
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-55078 and PDB model 9SPE. Per-residue inclusion information can be found in section [3](#) on page [10](#).

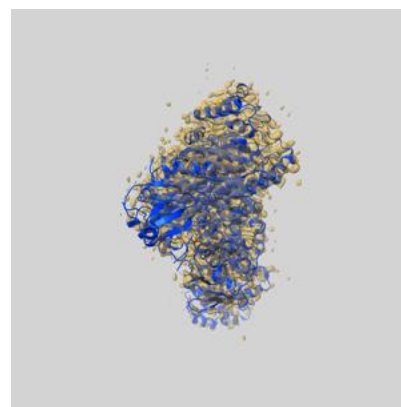
9.1 Map-model overlay [i](#)



X



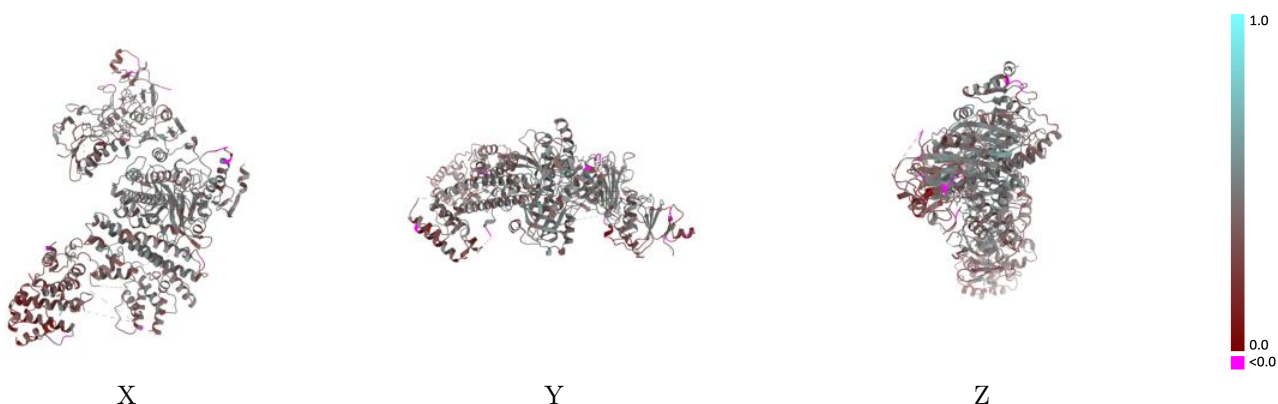
Y



Z

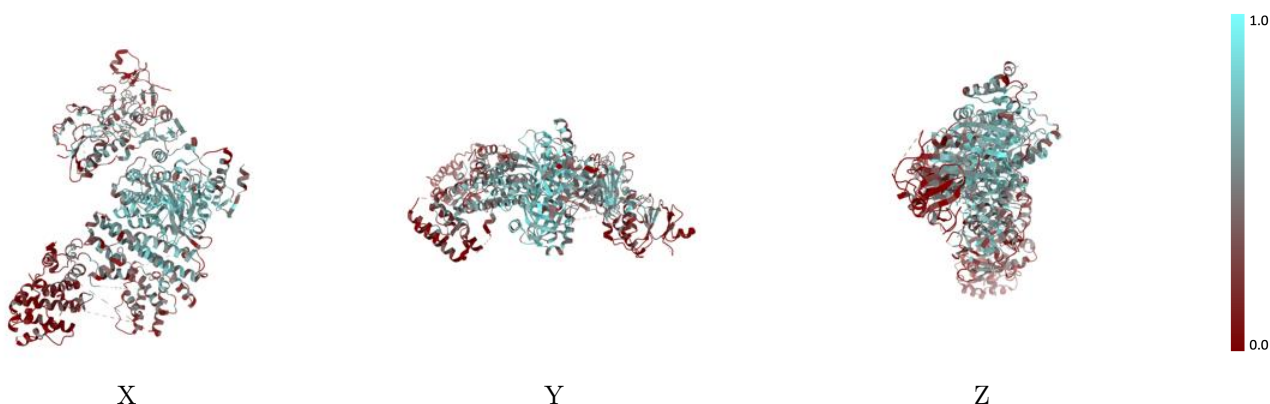
The images above show the 3D surface view of the map at the recommended contour level 0.1 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



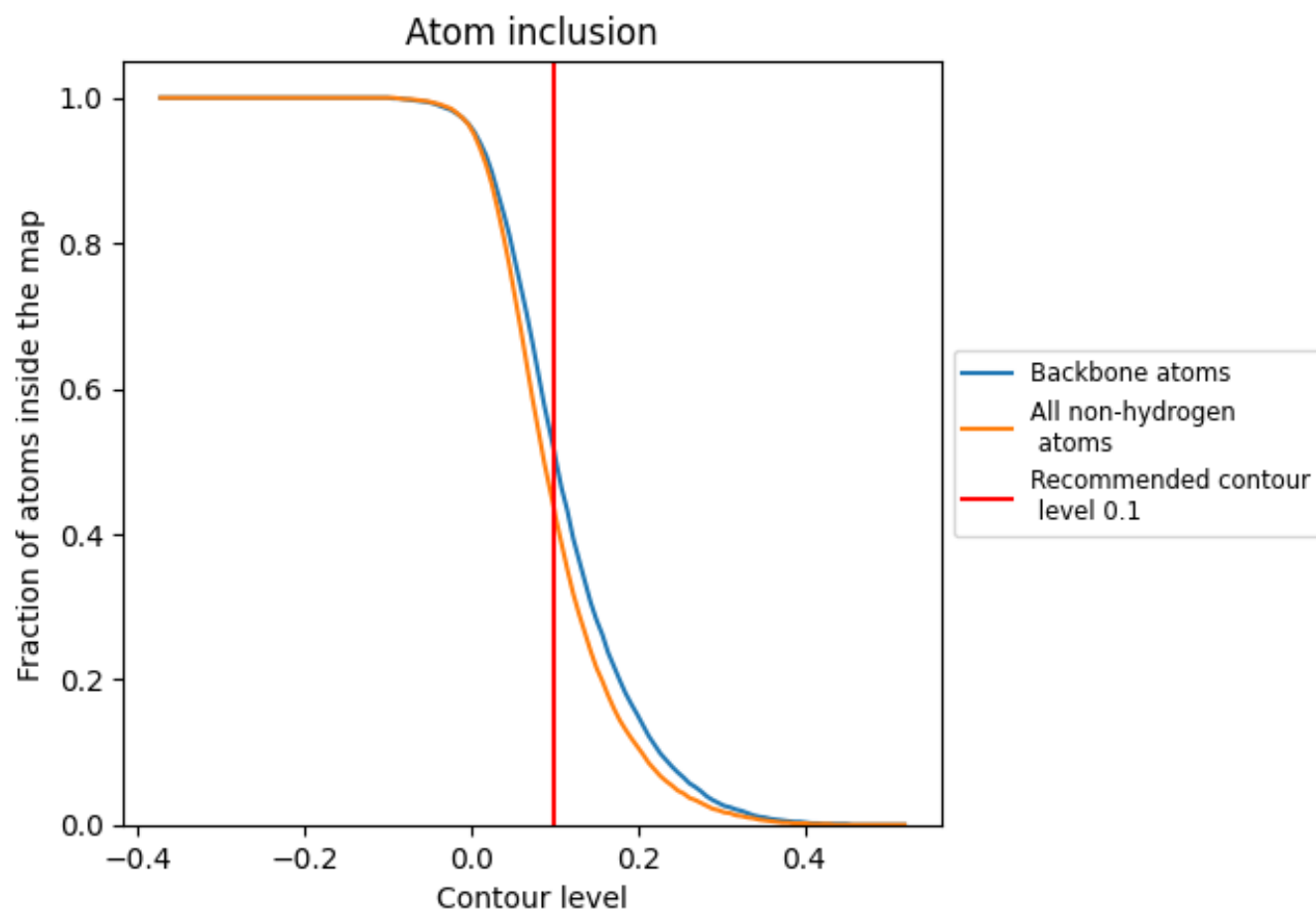
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

9.3 Atom inclusion mapped to coordinate model [i](#)



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.1).

9.4 Atom inclusion [i](#)



At the recommended contour level, 51% of all backbone atoms, 43% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary ⓘ

The table lists the average atom inclusion at the recommended contour level (0.1) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.4330	0.4020
A	0.2780	0.3500
B	0.2600	0.3420
C	0.3390	0.3920
D	0.4740	0.4110
P	0.3740	0.3970
R	0.6460	0.4520

