



Full wwPDB EM Validation Report ⓘ

Mar 9, 2026 – 12:02 PM UTC

PDB ID : 9M1H / pdb_00009m1h
EMDB ID : EMD-63571
Title : Cryo-EM structure of PGE2-EP1-Gq complex
Authors : Meng, X.; Xu, Y.; Xu, H.E.
Deposited on : 2025-02-26
Resolution : 2.55 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:

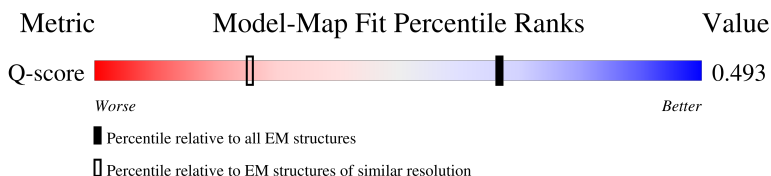
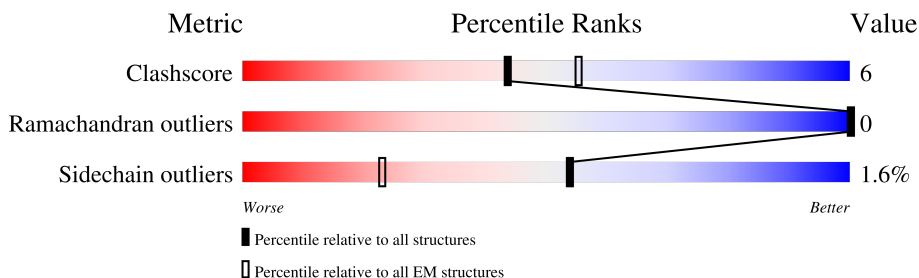
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data 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) : 2.49

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 2.55 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	7475 (2.05 - 3.05)

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	402	
2	B	361	
3	C	345	
4	G	71	

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 7055 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 Prostaglandin E2 receptor EP1 subtype.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	286	Total	C	N	O	S	0	0
			2123	1382	390	336	15		

- Molecule 2 is a protein called Guanine nucleotide-binding protein G(i) subunit alpha-1, Guanine nucleotide-binding protein G(s) subunit alpha isoforms short, Guanine nucleotide-binding protein G(q) subunit alpha.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	232	Total	C	N	O	S	0	0
			1908	1207	338	355	8		

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	32	ARG	ALA	conflict	UNP P63092
B	34	LEU	HIS	conflict	UNP P63092
B	42	ASP	GLY	conflict	UNP P63092
B	43	ASN	GLU	conflict	UNP P63092
B	56	TYR	LEU	conflict	UNP P63092
B	203	ALA	GLY	conflict	UNP P63092
B	226	ASP	ALA	conflict	UNP P63092
B	229	ASP	-	linker	UNP P63092
B	230	TYR	-	linker	UNP P63092
B	239	ASP	LEU	conflict	UNP P63092
B	310	LYS	ASP	conflict	UNP P63092
B	313	VAL	LEU	conflict	UNP P63092
B	314	ASP	ARG	conflict	UNP P63092
B	325	ILE	TYR	conflict	UNP P63092
B	333	SER	ALA	conflict	UNP P63092
B	339	ALA	ILE	conflict	UNP P63092
B	342	ILE	VAL	conflict	UNP P63092
B	349	ILE	THR	conflict	UNP P50148
B	353	MET	LEU	conflict	UNP P50148

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Chain	Residue	Modelled	Actual	Comment	Reference
B	356	ARG	LYS	conflict	UNP P50148

- Molecule 3 is a protein called Guanine nucleotide-binding protein G(I)/G(S)/G(T) subunit beta-1.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	338	Total	C	N	O	S	0	0
			2591	1597	467	506	21		

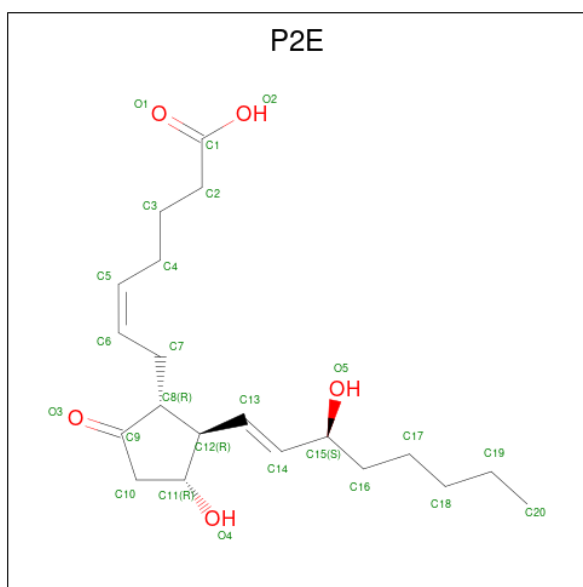
There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	1	MET	-	initiating methionine	UNP P62873
C	2	GLY	-	expression tag	UNP P62873
C	3	SER	-	expression tag	UNP P62873
C	4	LEU	-	expression tag	UNP P62873
C	5	LEU	-	expression tag	UNP P62873
C	6	GLN	-	expression tag	UNP P62873

- Molecule 4 is a protein called Guanine nucleotide-binding protein G(I)/G(S)/G(O) subunit gamma-2.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	G	54	Total	C	N	O	S	0	0
			407	253	71	80	3		

- Molecule 5 is (Z)-7-[(1R,2R,3R)-3-hydroxy-2-[(E,3S)-3-hydroxyoct-1-enyl]-5-oxo-cyclopentyl]hept-5-enoic acid (CCD ID: P2E) (formula: C₂₀H₃₂O₅).



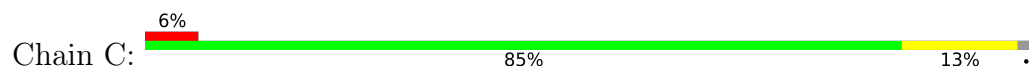
Mol	Chain	Residues	Atoms			AltConf
5	A	1	Total	C	O	0
			25	20	5	

- Molecule 6 is water.

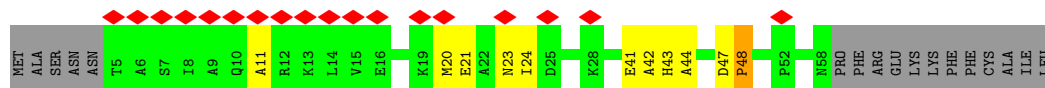
Mol	Chain	Residues	Atoms		AltConf
6	A	1	Total	O	0
			1	1	



- Molecule 3: Guanine nucleotide-binding protein G(I)/G(S)/G(T) subunit beta-1



- Molecule 4: Guanine nucleotide-binding protein G(I)/G(S)/G(O) subunit gamma-2



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	543885	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	30	Depositor
Minimum defocus (nm)	1200	Depositor
Maximum defocus (nm)	5000	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	4.689	Depositor
Minimum map value	-2.439	Depositor
Average map value	-0.001	Depositor
Map value standard deviation	0.071	Depositor
Recommended contour level	0.354	Depositor
Map size ($\text{\AA}$)	263.68, 263.68, 263.68	wwPDB
Map dimensions	320, 320, 320	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	0.824, 0.824, 0.824	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: P2E

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	1.05	26/2165 (1.2%)	0.76	7/2950 (0.2%)
2	B	0.41	3/1942 (0.2%)	0.44	3/2614 (0.1%)
3	C	0.11	0/2637	0.30	0/3576
4	G	1.22	7/411 (1.7%)	0.68	2/555 (0.4%)
All	All	0.69	36/7155 (0.5%)	0.53	12/9695 (0.1%)

All (36) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	334	PHE	C-O	-8.15	1.14	1.24
4	G	44	ALA	CA-C	-7.79	1.42	1.52
1	A	135	ARG	C-O	-7.59	1.14	1.24
1	A	347	ASP	C-O	-7.57	1.17	1.24
4	G	48	PRO	CA-CB	-7.47	1.43	1.53
1	A	220	VAL	C-O	-7.37	1.15	1.24
4	G	47	ASP	C-O	-7.33	1.19	1.25
2	B	251	THR	N-CA	-7.30	1.35	1.45
1	A	298	GLN	C-O	-7.12	1.15	1.24
1	A	221	CYS	C-O	-6.98	1.15	1.24
1	A	334	PHE	N-CA	-6.62	1.38	1.46
1	A	352	ILE	C-N	6.56	1.42	1.34
1	A	225	SER	C-O	-6.52	1.15	1.24
1	A	333	LEU	C-O	-6.44	1.15	1.24
4	G	43	HIS	CA-C	-6.18	1.44	1.52
1	A	137	VAL	CA-C	-6.02	1.44	1.52
2	B	250	ARG	CA-C	-5.86	1.45	1.52
1	A	137	VAL	C-O	-5.84	1.16	1.24
4	G	42	ALA	CA-C	-5.83	1.44	1.52
1	A	135	ARG	N-CA	-5.82	1.38	1.46
1	A	167	ALA	N-CA	-5.77	1.39	1.46
1	A	298	GLN	CA-C	-5.76	1.45	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	333	LEU	CA-C	-5.74	1.44	1.52
1	A	163	ALA	CA-C	-5.67	1.45	1.52
1	A	166	LEU	CA-C	-5.65	1.45	1.52
1	A	178	ARG	CA-C	-5.60	1.45	1.52
1	A	334	PHE	CA-C	-5.43	1.45	1.52
2	B	249	LEU	CA-C	-5.41	1.45	1.52
1	A	333	LEU	CA-CB	-5.24	1.45	1.53
4	G	47	ASP	CA-C	-5.21	1.46	1.52
1	A	117	MET	C-O	-5.16	1.18	1.24
1	A	167	ALA	CA-C	-5.16	1.46	1.52
4	G	41	GLU	CA-C	-5.16	1.45	1.52
1	A	117	MET	N-CA	5.10	1.52	1.46
1	A	137	VAL	N-CA	-5.09	1.40	1.46
1	A	165	ALA	N-CA	-5.03	1.40	1.46

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	334	PHE	N-CA-C	-9.70	99.76	111.69
2	B	251	THR	N-CA-C	-8.79	100.00	113.89
1	A	137	VAL	N-CA-C	-7.09	103.27	111.00
2	B	251	THR	CA-C-N	-7.03	114.07	122.93
2	B	251	THR	C-N-CA	-7.03	114.07	122.93
1	A	340	ALA	O-C-N	-6.36	114.78	122.22
1	A	163	ALA	N-CA-C	5.36	116.93	111.14
4	G	47	ASP	CA-C-N	-5.19	113.35	119.84
4	G	47	ASP	C-N-CA	-5.19	113.35	119.84
1	A	223	THR	N-CA-C	-5.13	105.32	112.45
1	A	135	ARG	N-CA-C	5.12	119.37	113.18
1	A	333	LEU	N-CA-C	5.09	119.55	113.23

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2123	0	2270	35	0
2	B	1908	0	1905	24	0
3	C	2591	0	2496	30	0
4	G	407	0	419	4	0
5	A	25	0	30	5	0
6	A	1	0	0	0	0
All	All	7055	0	7120	92	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 6.

All (92) 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:119:PHE:HE2	1:A:166:LEU:CD1	1.58	1.17
1:A:119:PHE:CE2	1:A:166:LEU:HD12	1.83	1.13
1:A:119:PHE:CE2	1:A:166:LEU:CD1	2.31	1.13
1:A:119:PHE:HE2	1:A:166:LEU:HD12	1.18	0.97
1:A:119:PHE:CE2	1:A:166:LEU:HD11	2.06	0.90
1:A:59:GLN:OE1	1:A:360:ARG:HG2	1.71	0.90
2:B:323:ARG:NH1	2:B:324:HIS:HE1	1.75	0.84
1:A:93:ALA:O	1:A:96:LEU:HG	1.89	0.72
2:B:323:ARG:NH1	2:B:324:HIS:CE1	2.59	0.69
2:B:244:TRP:CH2	2:B:324:HIS:CD2	2.81	0.69
1:A:119:PHE:CD2	1:A:166:LEU:HD12	2.26	0.68
2:B:51:LYS:O	2:B:54:ARG:NH1	2.27	0.68
3:C:272:ASP:OD1	3:C:273:ASN:ND2	2.28	0.67
3:C:231:GLU:N	3:C:231:GLU:OE1	2.29	0.66
4:G:20:MET:O	4:G:23:ASN:ND2	2.29	0.66
2:B:323:ARG:HH11	2:B:324:HIS:CE1	2.15	0.64
1:A:162:ALA:O	1:A:166:LEU:HD13	1.98	0.63
1:A:95:VAL:HG21	5:A:501:P2E:H7	1.82	0.61
3:C:12:LEU:HD13	4:G:11:ALA:HB1	1.83	0.61
2:B:318:ALA:O	2:B:319:SER:OG	2.16	0.60
3:C:12:LEU:HD12	3:C:13:ARG:N	2.17	0.60
2:B:323:ARG:HH12	2:B:324:HIS:HE1	1.54	0.56
2:B:281:GLU:N	2:B:281:GLU:OE1	2.38	0.55
3:C:56:LEU:HD23	3:C:92:THR:HG23	1.89	0.54
3:C:119:CYS:SG	3:C:127:SER:OG	2.65	0.54
2:B:46:LYS:O	2:B:50:VAL:HG23	2.07	0.54
1:A:311:SER:O	1:A:315:VAL:HG22	2.08	0.53
2:B:310:LYS:HA	2:B:313:VAL:HG22	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:344:GLN:NE2	5:A:501:P2E:H16A	2.23	0.53
3:C:168:ASP:O	3:C:169:THR:OG1	2.23	0.53
3:C:291:LEU:HD22	3:C:301:VAL:HG22	1.91	0.53
1:A:293:VAL:O	1:A:296:VAL:HG22	2.09	0.52
3:C:291:LEU:CD2	3:C:301:VAL:HG13	2.40	0.52
1:A:59:GLN:OE1	1:A:360:ARG:CG	2.53	0.52
1:A:170:LEU:O	1:A:171:LEU:C	2.52	0.51
1:A:96:LEU:HD12	1:A:97:ARG:N	2.26	0.51
3:C:318:ASN:OD1	3:C:319:ARG:N	2.40	0.51
1:A:178:ARG:HG3	1:A:191:GLY:HA3	1.92	0.51
3:C:291:LEU:HD23	3:C:301:VAL:HG13	1.92	0.51
2:B:250:ARG:O	2:B:323:ARG:HD2	2.11	0.50
3:C:281:VAL:HG13	3:C:290:LEU:HD21	1.94	0.50
2:B:244:TRP:CZ3	2:B:324:HIS:HD2	2.30	0.49
3:C:79:SER:OG	3:C:81:ASP:OD1	2.29	0.49
3:C:329:GLY:O	4:G:48:PRO:HB2	2.14	0.48
3:C:76:VAL:HG23	3:C:86:ILE:CD1	2.43	0.48
2:B:244:TRP:CZ2	2:B:324:HIS:CD2	3.02	0.47
2:B:244:TRP:CH2	2:B:324:HIS:HD2	2.28	0.47
2:B:262:ASP:OD1	2:B:263:LEU:N	2.48	0.47
3:C:230:HIS:NE2	3:C:248:THR:O	2.48	0.47
2:B:52:GLN:NE2	2:B:338:ASN:OD1	2.47	0.46
3:C:140:VAL:HG13	3:C:140:VAL:O	2.15	0.46
2:B:309:ARG:O	2:B:313:VAL:HG13	2.15	0.46
1:A:209:SER:O	1:A:213:VAL:HG23	2.16	0.46
3:C:76:VAL:HG23	3:C:86:ILE:HD13	1.97	0.46
1:A:90:ILE:HB	1:A:91:PRO:HD3	1.98	0.46
3:C:144:LEU:HD12	3:C:144:LEU:O	2.16	0.46
1:A:187:TRP:CH2	1:A:337:VAL:HG11	2.52	0.45
3:C:154:CYS:O	3:C:155:ARG:NH1	2.48	0.45
3:C:160:ASN:ND2	3:C:175:ASP:OD1	2.50	0.45
1:A:117:MET:HE2	5:A:501:P2E:H17A	1.99	0.44
1:A:157:ALA:O	1:A:161:VAL:HG23	2.16	0.44
3:C:130:ASN:OD1	3:C:131:LEU:N	2.51	0.44
3:C:178:THR:HG23	3:C:180:GLN:H	1.82	0.44
3:C:194:SER:OG	3:C:237:ILE:HG22	2.17	0.44
3:C:278:ILE:HD12	3:C:278:ILE:N	2.33	0.44
1:A:215:LEU:O	1:A:219:LEU:HD13	2.18	0.43
2:B:313:VAL:O	2:B:317:THR:HG23	2.19	0.43
2:B:207:GLU:OE2	2:B:210:LYS:N	2.46	0.43
2:B:275:ILE:HD12	2:B:275:ILE:N	2.34	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A:501:P2E:H6	5:A:501:P2E:H12	1.89	0.42
1:A:340:ALA:O	1:A:341:SER:C	2.62	0.42
1:A:44:THR:O	1:A:48:VAL:HG23	2.19	0.42
1:A:168:VAL:HA	1:A:171:LEU:HG	2.02	0.42
2:B:221:ILE:O	2:B:255:ILE:N	2.41	0.42
3:C:67:HIS:O	3:C:76:VAL:HG12	2.20	0.42
3:C:128:ILE:N	3:C:128:ILE:HD12	2.34	0.42
4:G:21:GLU:O	4:G:24:ILE:HG22	2.20	0.42
2:B:329:HIS:NE2	2:B:345:ASP:OD2	2.52	0.42
1:A:81:LEU:HD21	1:A:352:ILE:HD12	2.02	0.42
1:A:131:MET:O	1:A:135:ARG:HG3	2.20	0.42
3:C:92:THR:HG22	3:C:92:THR:O	2.20	0.42
1:A:177:GLY:HA2	1:A:200:GLN:HA	2.02	0.41
3:C:325:VAL:HG13	3:C:332:VAL:HG22	2.03	0.41
1:A:156:LEU:C	1:A:156:LEU:HD23	2.45	0.41
3:C:17:GLU:O	3:C:21:ASN:ND2	2.47	0.41
1:A:355:ARG:HB3	1:A:356:GLN:H	1.74	0.41
2:B:297:GLU:O	2:B:298:ASP:C	2.63	0.41
1:A:37:ALA:O	1:A:40:ILE:HG12	2.21	0.41
1:A:329:LEU:O	1:A:333:LEU:HB2	2.21	0.40
2:B:275:ILE:HD12	2:B:275:ILE:H	1.86	0.40
1:A:77:VAL:HG23	1:A:78:ALA:N	2.37	0.40
1:A:344:GLN:HE21	5:A:501:P2E:H16A	1.86	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	282/402 (70%)	264 (94%)	18 (6%)	0	100	100
2	B	228/361 (63%)	222 (97%)	6 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	C	336/345 (97%)	329 (98%)	7 (2%)	0	100	100
4	G	52/71 (73%)	50 (96%)	2 (4%)	0	100	100
All	All	898/1179 (76%)	865 (96%)	33 (4%)	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	211/296 (71%)	202 (96%)	9 (4%)	26	39
2	B	210/316 (66%)	208 (99%)	2 (1%)	68	80
3	C	279/287 (97%)	278 (100%)	1 (0%)	84	90
4	G	43/58 (74%)	43 (100%)	0	100	100
All	All	743/957 (78%)	731 (98%)	12 (2%)	54	73

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	59	GLN
1	A	63	ARG
1	A	65	ARG
1	A	117	MET
1	A	178	ARG
1	A	223	THR
1	A	224	LEU
1	A	333	LEU
1	A	347	ASP
2	B	360	LEU
2	B	361	VAL
3	C	64	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	111	HIS
1	A	145	HIS
1	A	298	GLN
1	A	343	ASN
1	A	344	GLN
2	B	231	ASN
2	B	324	HIS
4	G	23	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

1 ligand 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	$\# Z > 2$
5	P2E	A	501	-	25,25,25	3.68	9 (36%)	25,31,31	1.84	5 (20%)

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
5	P2E	A	501	-	-	10/19/35/35	0/1/1/1

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	501	P2E	C10-C9	-10.77	1.36	1.51
5	A	501	P2E	C12-C8	-9.51	1.38	1.55
5	A	501	P2E	O4-C11	-7.58	1.27	1.43
5	A	501	P2E	C8-C9	4.40	1.59	1.52
5	A	501	P2E	C10-C11	3.82	1.57	1.52
5	A	501	P2E	C12-C11	3.15	1.62	1.55
5	A	501	P2E	O5-C15	-3.06	1.38	1.43
5	A	501	P2E	O3-C9	-2.04	1.18	1.21
5	A	501	P2E	C15-C14	2.00	1.54	1.50

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	501	P2E	C7-C6-C5	-4.50	109.68	126.21
5	A	501	P2E	C15-C14-C13	-4.20	114.82	124.74
5	A	501	P2E	C12-C13-C14	-3.30	113.80	124.86
5	A	501	P2E	C16-C15-C14	-2.35	107.44	111.94
5	A	501	P2E	C3-C2-C1	-2.31	108.47	114.51

There are no chirality outliers.

All (10) torsion outliers are listed below:

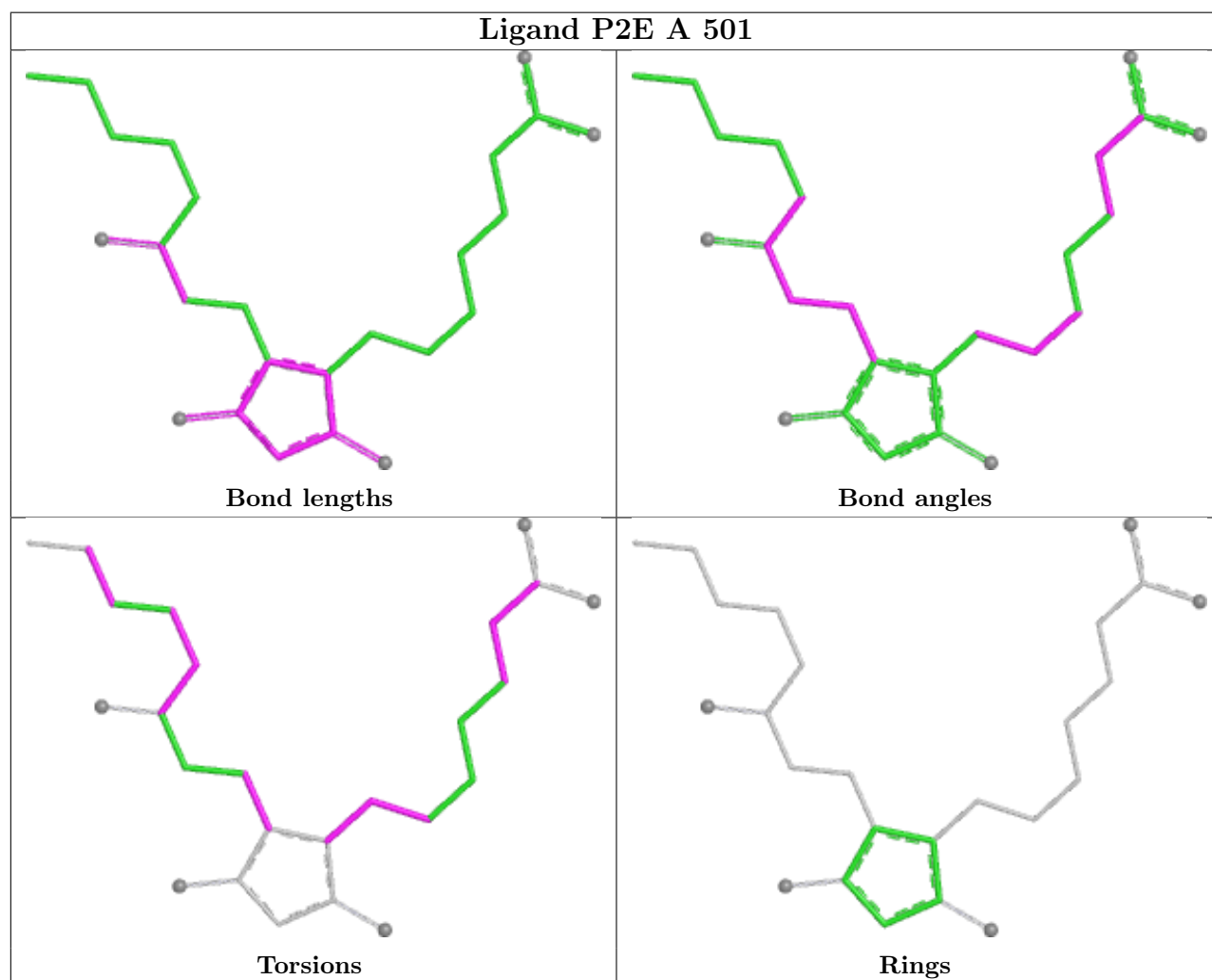
Mol	Chain	Res	Type	Atoms
5	A	501	P2E	C14-C15-C16-C17
5	A	501	P2E	O5-C15-C16-C17
5	A	501	P2E	C1-C2-C3-C4
5	A	501	P2E	C17-C18-C19-C20
5	A	501	P2E	C8-C12-C13-C14
5	A	501	P2E	C5-C6-C7-C8
5	A	501	P2E	O1-C1-C2-C3
5	A	501	P2E	O2-C1-C2-C3
5	A	501	P2E	C6-C7-C8-C12
5	A	501	P2E	C15-C16-C17-C18

There are no ring outliers.

1 monomer is involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	501	P2E	5	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

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

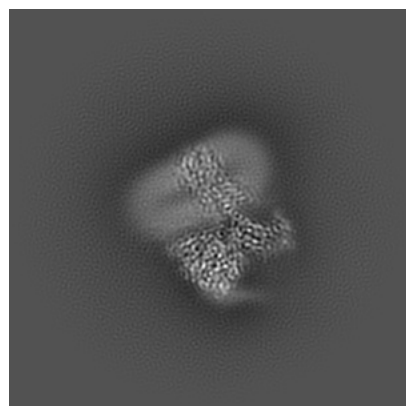
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-63571. These allow visual inspection of the internal detail of the map and identification of artifacts.

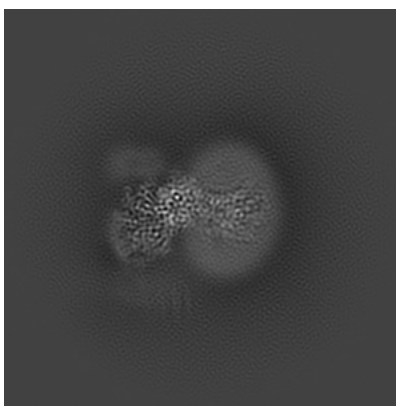
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

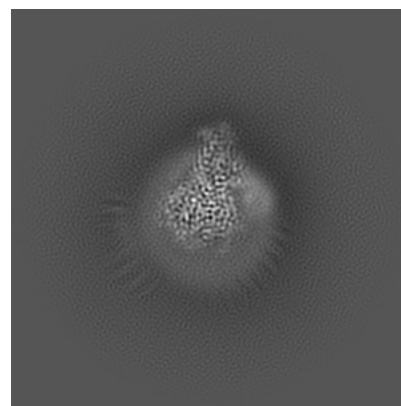
6.1.1 Primary map



X

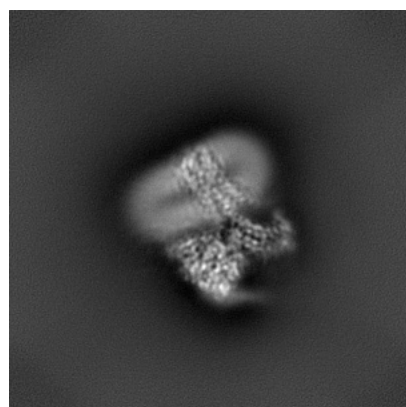


Y

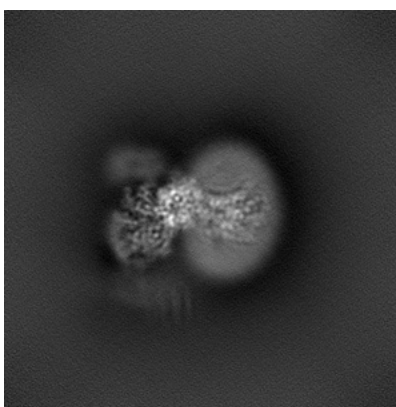


Z

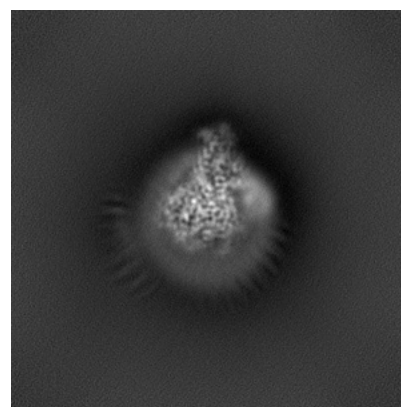
6.1.2 Raw map



X



Y

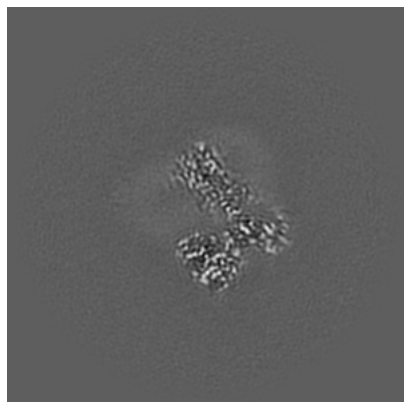


Z

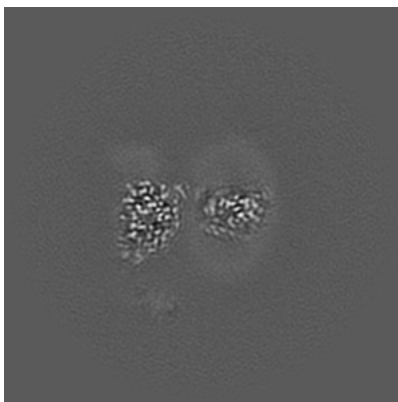
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

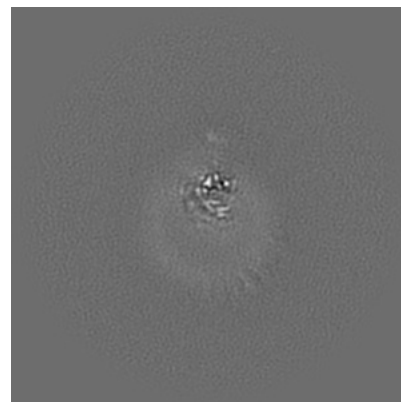
6.2.1 Primary map



X Index: 160

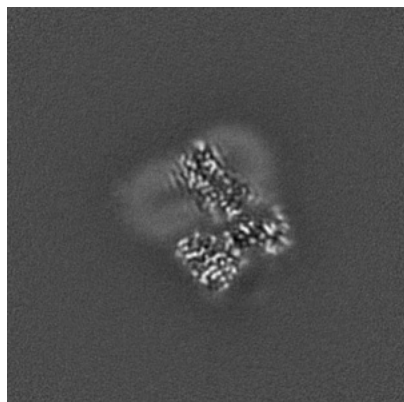


Y Index: 160

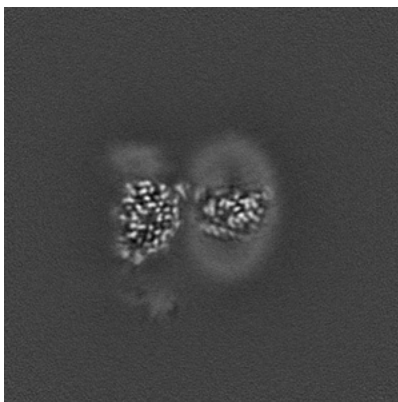


Z Index: 160

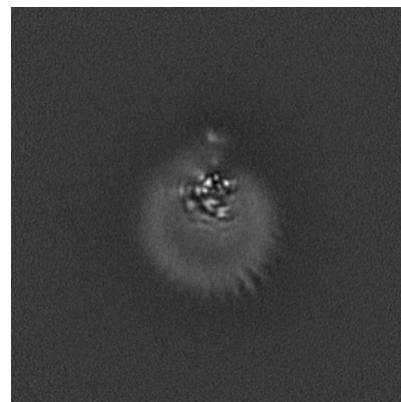
6.2.2 Raw map



X Index: 160



Y Index: 160

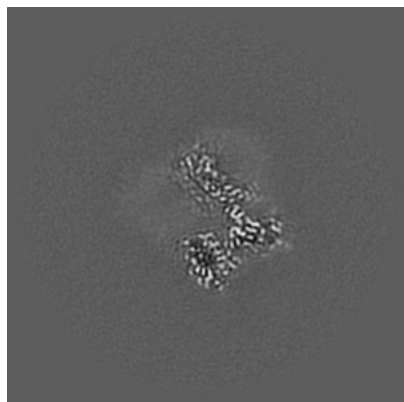


Z Index: 160

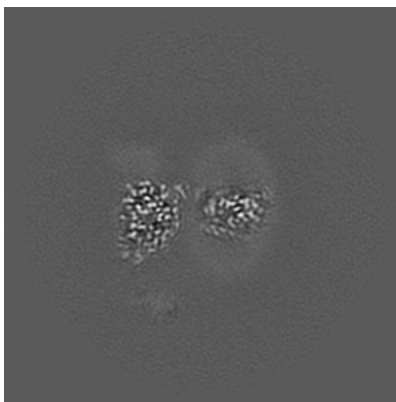
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

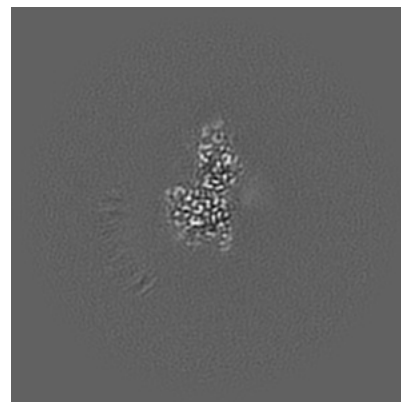
6.3.1 Primary map



X Index: 164

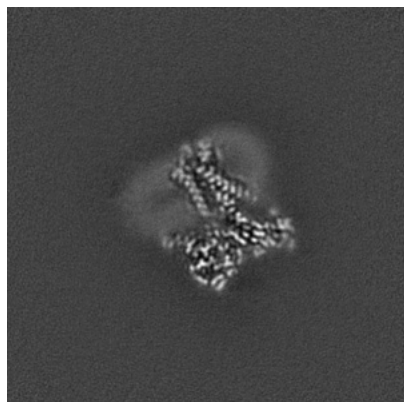


Y Index: 160

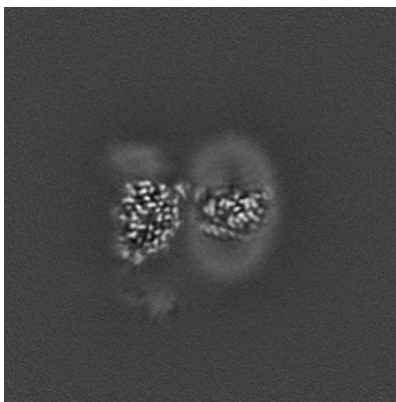


Z Index: 129

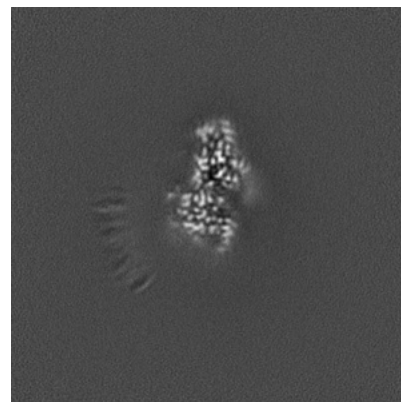
6.3.2 Raw map



X Index: 168



Y Index: 160

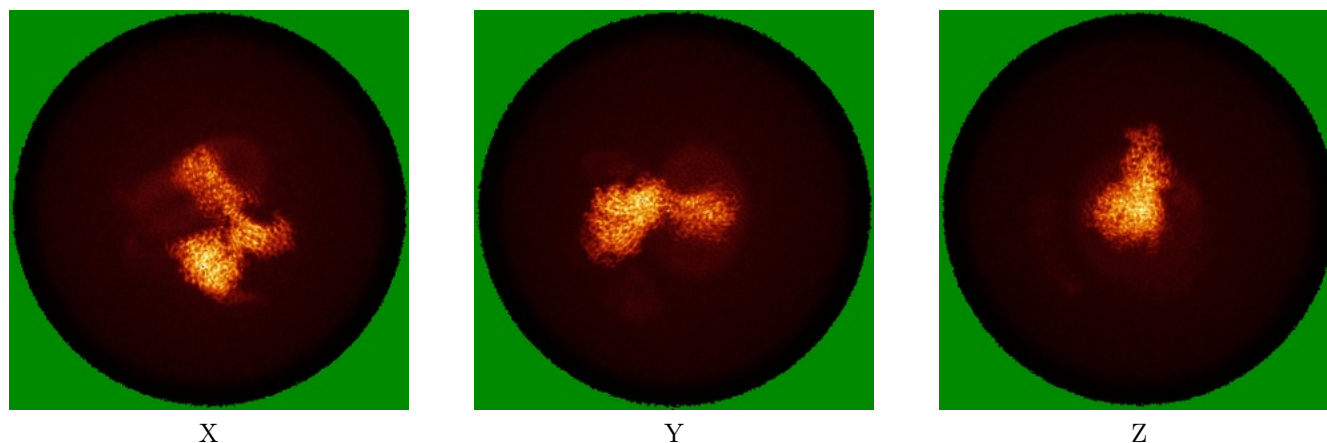


Z Index: 133

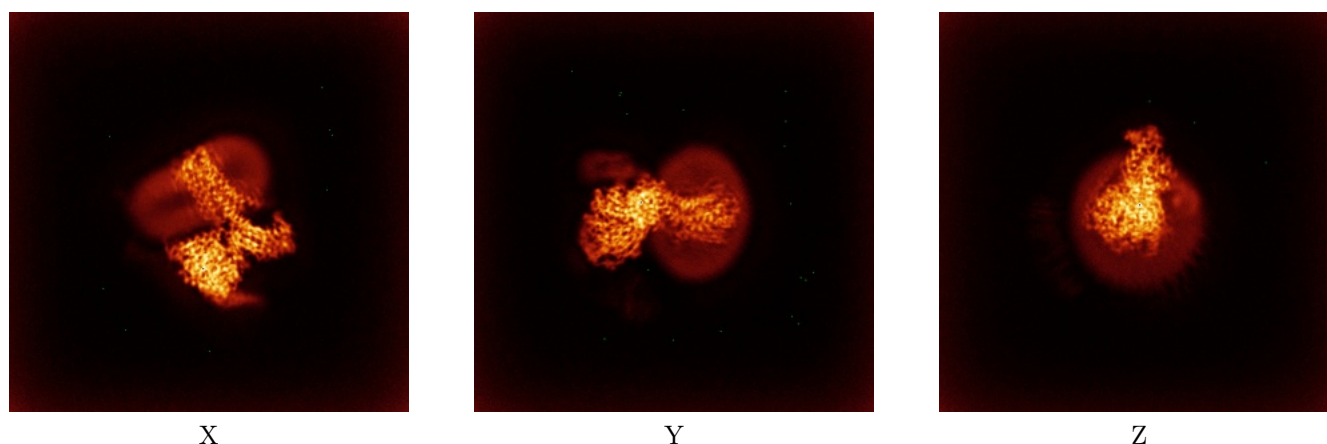
The images above show the largest variance slices of the map in three orthogonal directions.

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

6.4.1 Primary map



6.4.2 Raw map



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.

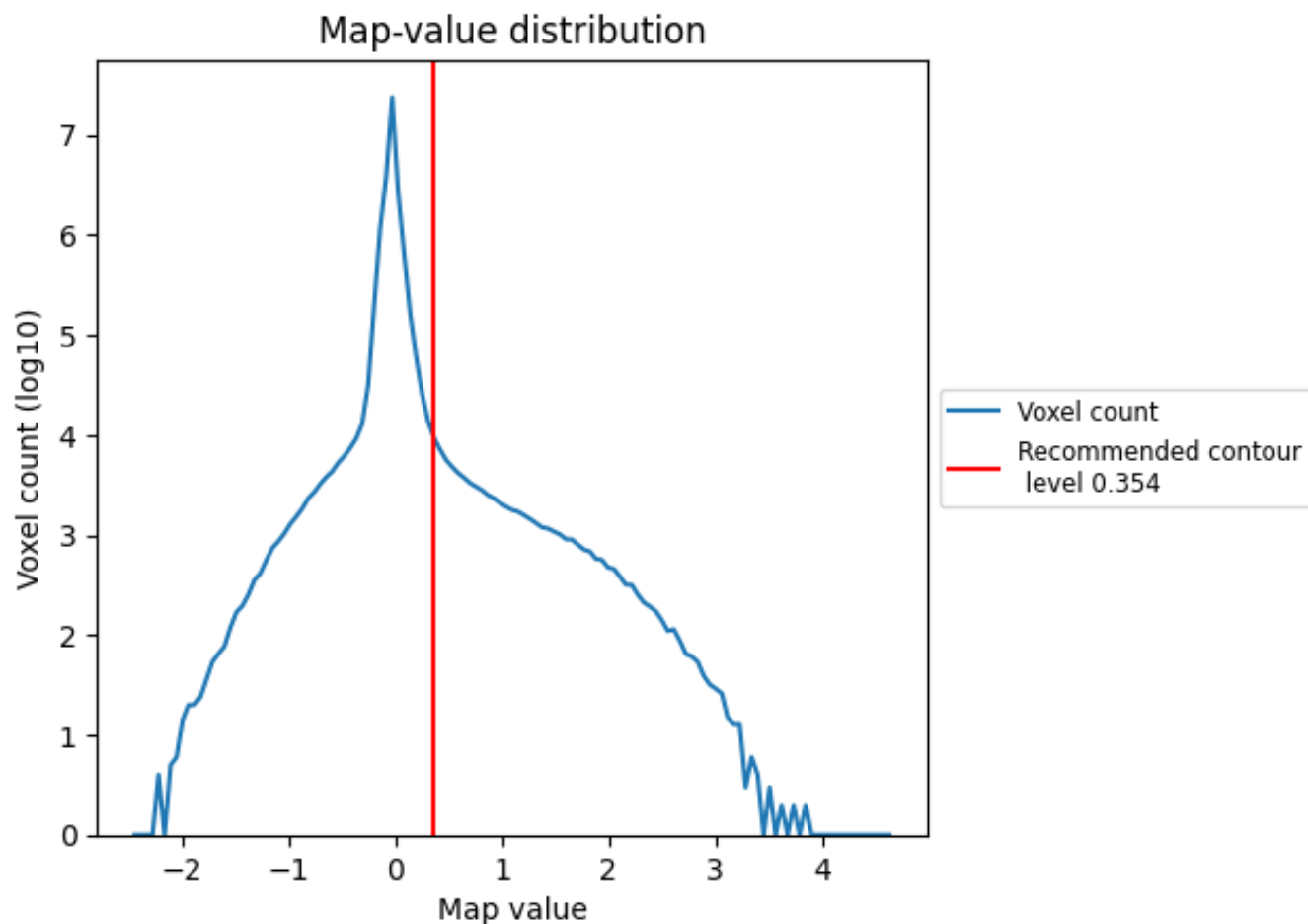
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

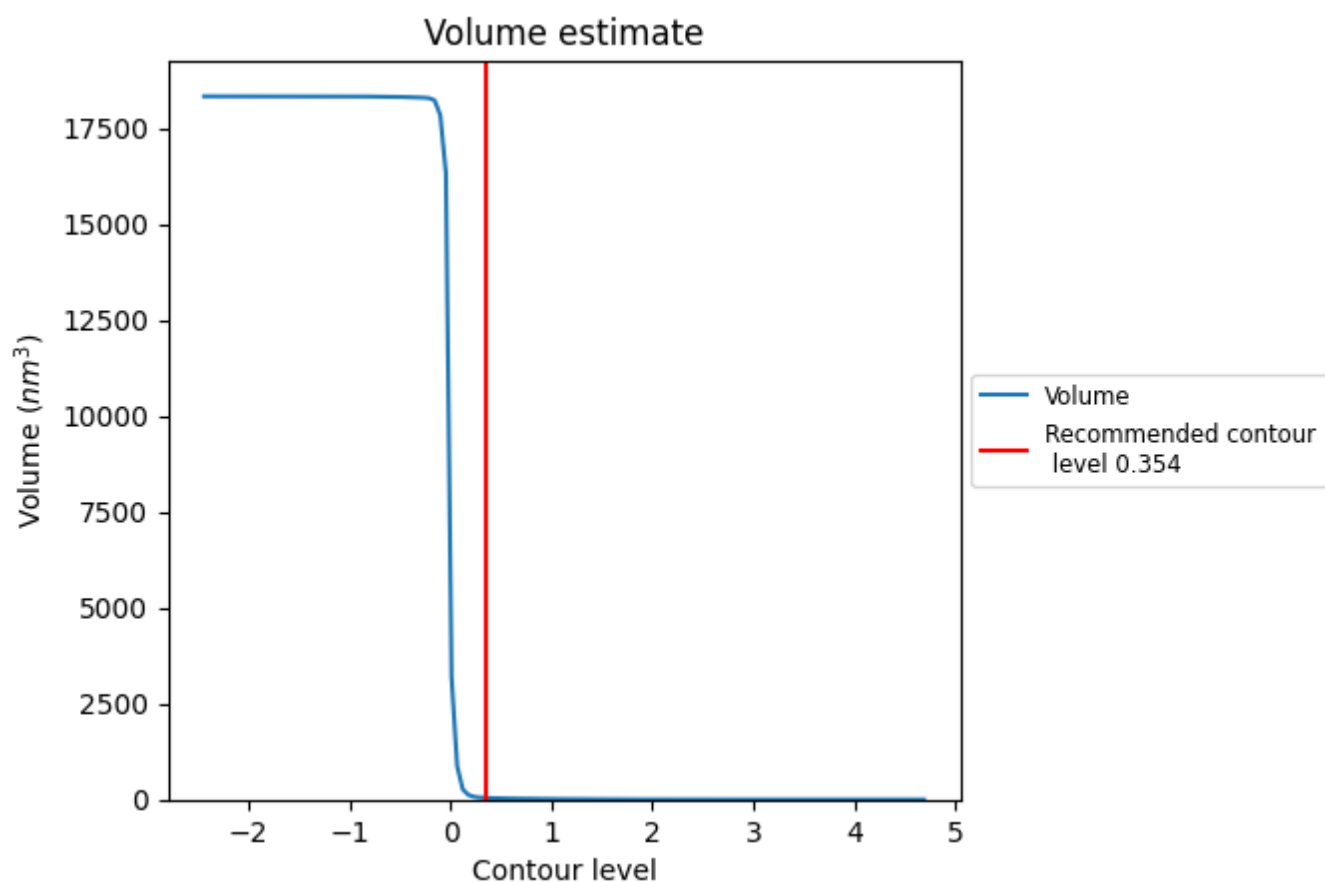
This section contains the results of statistical analysis of the map.

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.

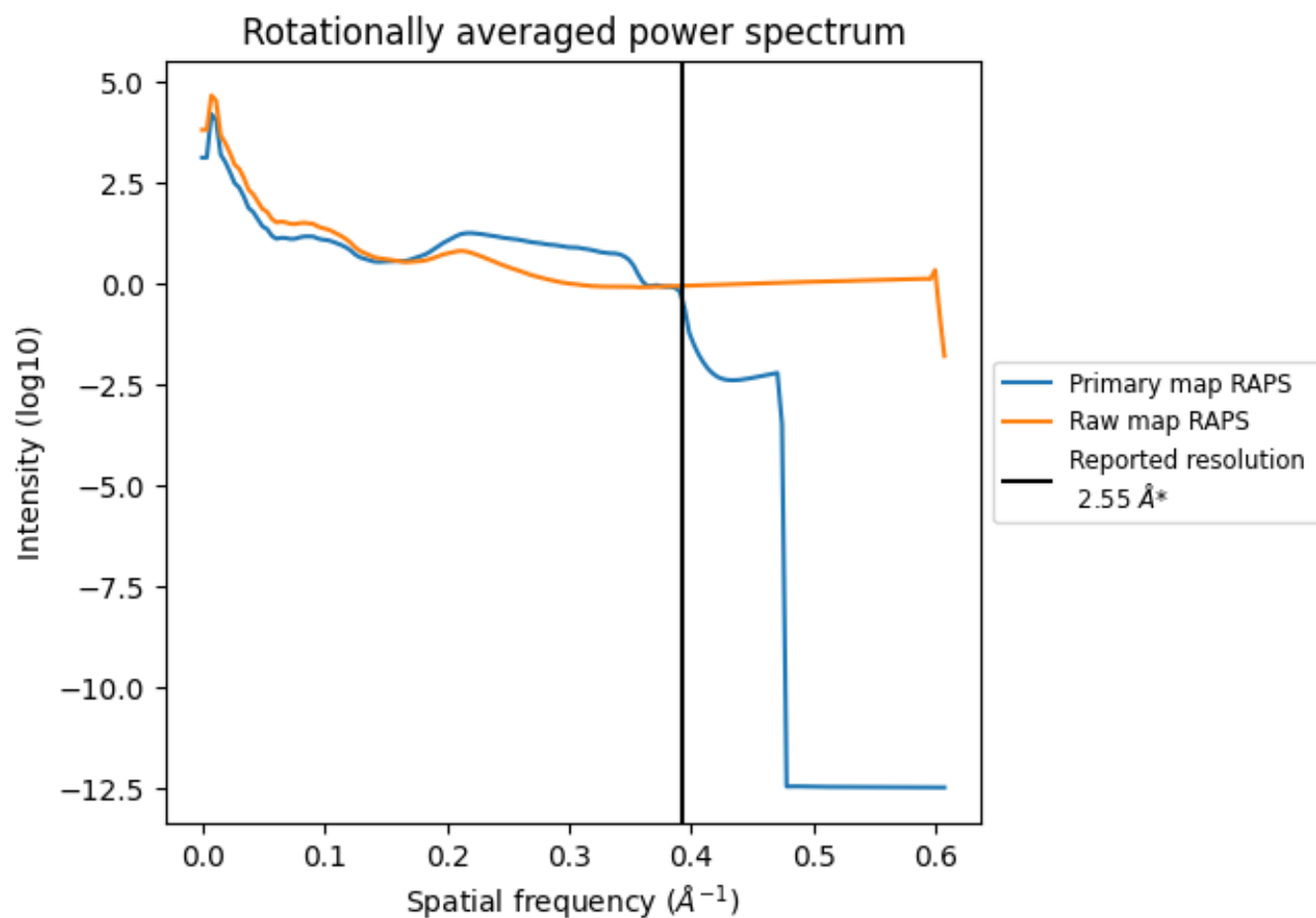
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 43 nm³; this corresponds to an approximate mass of 39 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 ⓘ

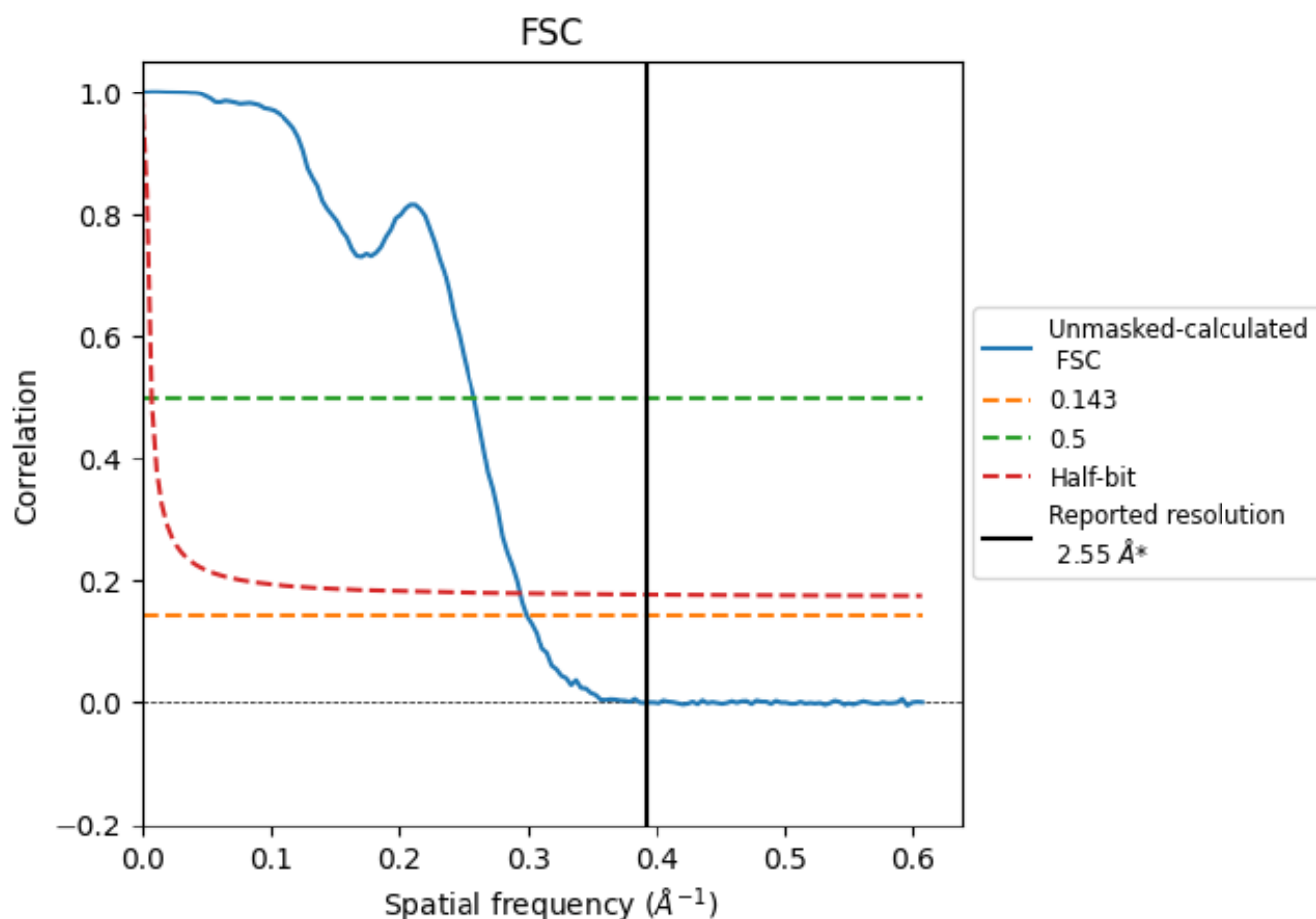


*Reported resolution corresponds to spatial frequency of 0.392 Å⁻¹

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 $0.392 \text{ \AA}^{-1}$

8.2 Resolution estimates [i](#)

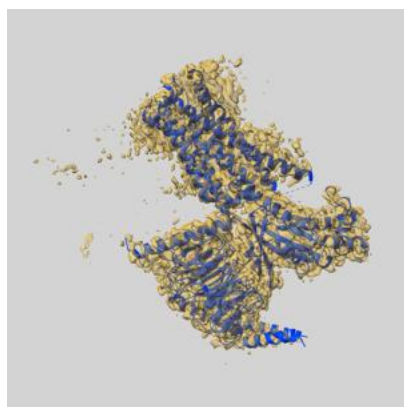
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.55	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	3.34	3.87	3.40

*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 3.34 differs from the reported value 2.55 by more than 10 %

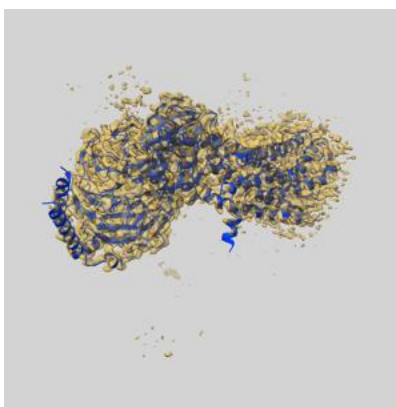
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-63571 and PDB model 9M1H. Per-residue inclusion information can be found in section [3](#) on page [6](#).

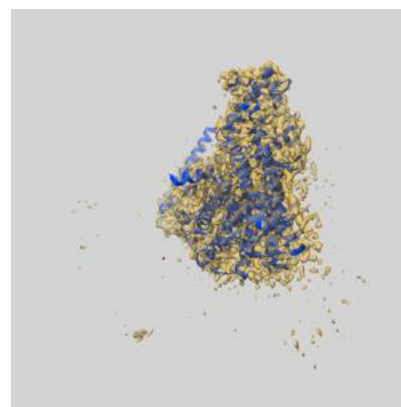
9.1 Map-model overlay [i](#)



X



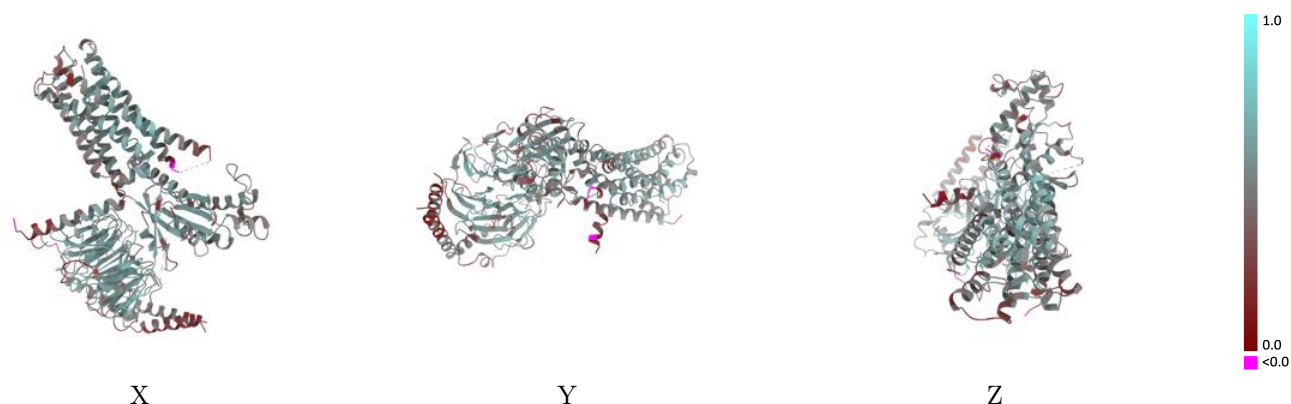
Y



Z

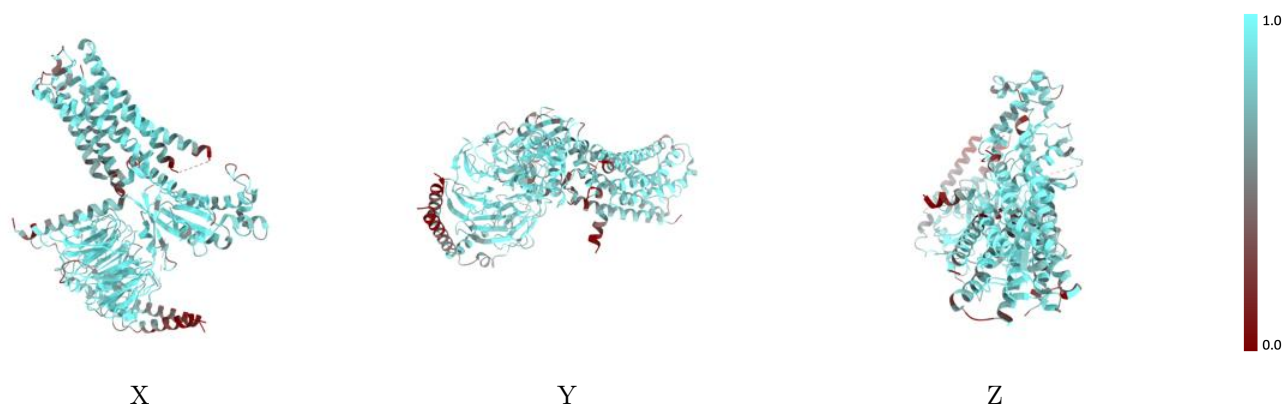
The images above show the 3D surface view of the map at the recommended contour level 0.354 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



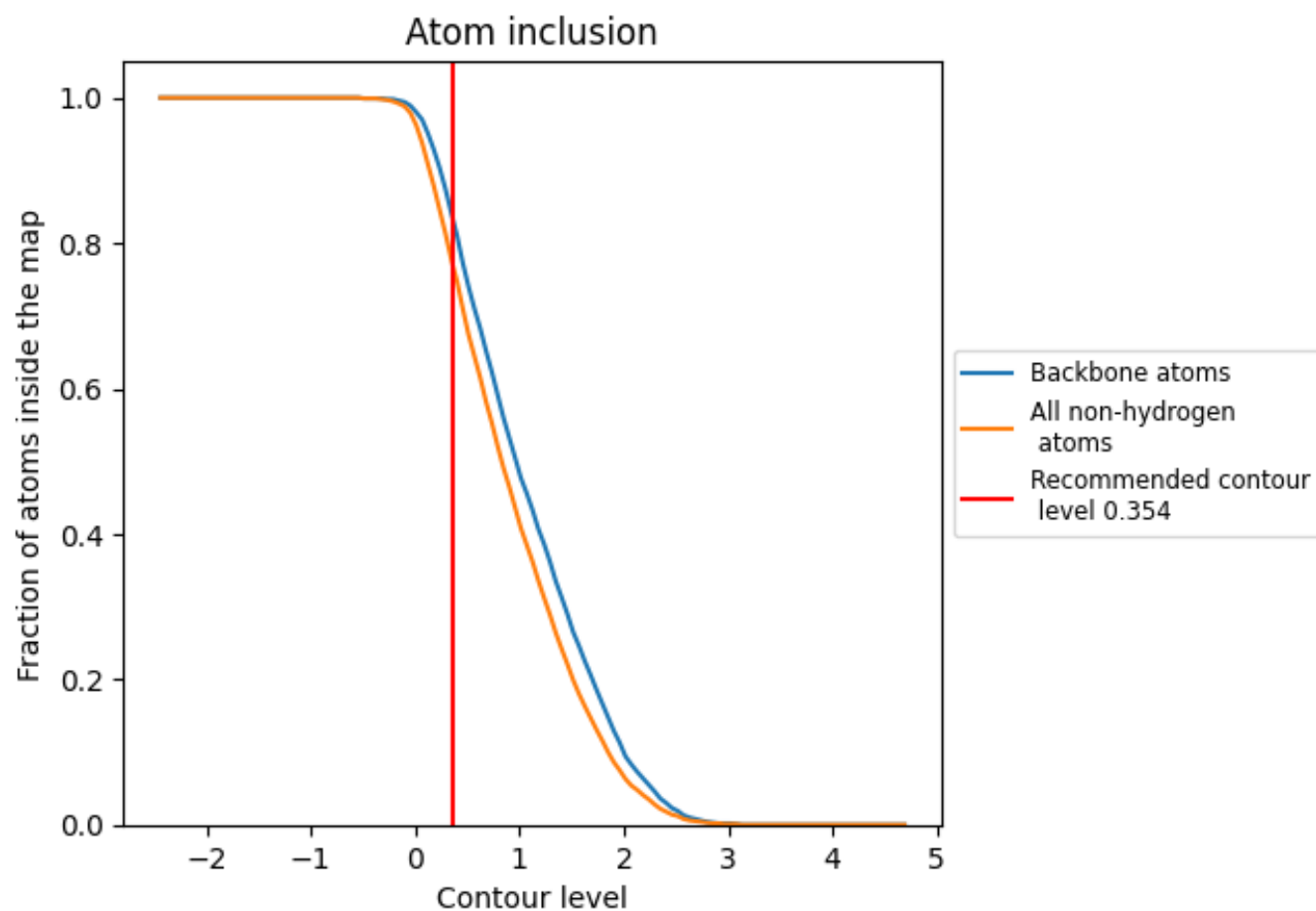
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.354).

9.4 Atom inclusion [i](#)



At the recommended contour level, 84% of all backbone atoms, 77% 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 (0.354) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7730	0.4930
A	0.7520	0.4850
B	0.7880	0.4900
C	0.8250	0.5210
G	0.5570	0.3780

