



Full wwPDB EM Validation Report ⓘ

Apr 5, 2026 – 06:56 PM UTC

PDB ID : 9CX3 / pdb_00009cx3
EMDB ID : EMD-45977
Title : Structure of SH3 domain of Src in complex with beta-arrestin 1
Authors : Pakharukova, N.; Bansia, H.; des Georges, A.; Lefkowitz, R.J.
Deposited on : 2024-07-30
Resolution : 3.47 Å (reported)
Based on initial models : 6ni2, 2ptk

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

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

A user guide is available at

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

The types of validation reports are described at

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

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

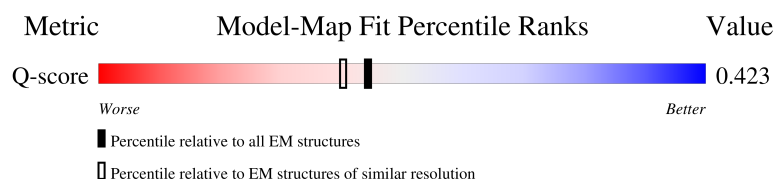
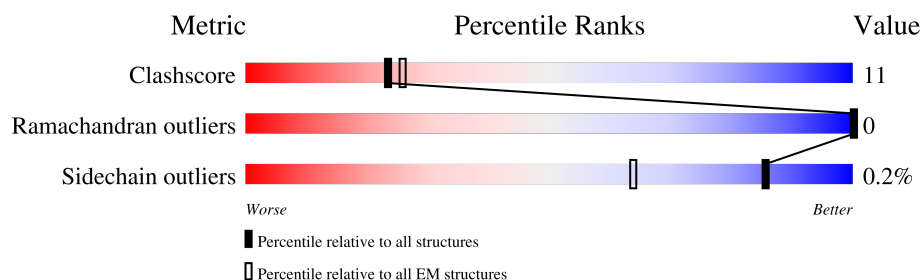
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13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

1 Overall quality at a glance

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

The reported resolution of this entry is 3.47 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	13733 (2.97 - 3.97)

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	124	 8% 69% 20% 10%
2	C	85	 8% 47% 18% 35%
3	H	237	 33% 15% 52%
4	V	29	 31% 10% 7% 52%

Continued on next page...

Mol	Chain	Length	Quality of chain
5	B	392	 62% 20% 18%
6	L	215	 38% 10% 52%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 5626 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 Nanobody 32.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	111	Total	C	N	O	S	0	0
			843	527	147	165	4		

- Molecule 2 is a protein called Proto-oncogene tyrosine-protein kinase Src.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	C	55	Total	C	N	O	S	0	0
			439	280	69	89	1		

There are 27 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	57	MET	-	initiating methionine	UNP P00523
C	58	GLY	-	expression tag	UNP P00523
C	59	SER	-	expression tag	UNP P00523
C	60	SER	-	expression tag	UNP P00523
C	61	HIS	-	expression tag	UNP P00523
C	62	HIS	-	expression tag	UNP P00523
C	63	HIS	-	expression tag	UNP P00523
C	64	HIS	-	expression tag	UNP P00523
C	65	HIS	-	expression tag	UNP P00523
C	66	HIS	-	expression tag	UNP P00523
C	67	ASP	-	expression tag	UNP P00523
C	68	TYR	-	expression tag	UNP P00523
C	69	ASP	-	expression tag	UNP P00523
C	70	ILE	-	expression tag	UNP P00523
C	71	PRO	-	expression tag	UNP P00523
C	72	THR	-	expression tag	UNP P00523
C	73	THR	-	expression tag	UNP P00523
C	74	GLU	-	expression tag	UNP P00523
C	75	ASN	-	expression tag	UNP P00523
C	76	LEU	-	expression tag	UNP P00523
C	77	TYR	-	expression tag	UNP P00523

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
C	78	PHE	-	expression tag	UNP P00523
C	79	GLN	-	expression tag	UNP P00523
C	80	GLY	-	expression tag	UNP P00523
C	81	HIS	-	expression tag	UNP P00523
C	82	MET	-	expression tag	UNP P00523
C	95	CYS	ARG	engineered mutation	UNP P00523

- Molecule 3 is a protein called Antibody fragment Fab30, heavy chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	H	114	Total	C	N	O	S	0	0
			872	556	143	170	3		

- Molecule 4 is a protein called Vasopressin V2 receptor.

Mol	Chain	Residues	Atoms					AltConf	Trace	
4	V	14	Total	C	N	O	P	S	0	0
			120	54	15	44	6	1		

- Molecule 5 is a protein called Beta-arrestin-1.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	B	323	Total	C	N	O	S	0	0
			2565	1645	442	474	4		

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	59	VAL	CYS	engineered mutation	UNP P29066
B	120	CYS	PRO	engineered mutation	UNP P29066
B	125	SER	CYS	engineered mutation	UNP P29066
B	140	LEU	CYS	engineered mutation	UNP P29066
B	150	VAL	CYS	engineered mutation	UNP P29066
B	242	VAL	CYS	engineered mutation	UNP P29066
B	251	VAL	CYS	engineered mutation	UNP P29066
B	269	SER	CYS	engineered mutation	UNP P29066

- Molecule 6 is a protein called Antibody fragment Fab30, light chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	L	103	Total	C	N	O	S	0	0
			787	498	130	156	3		

Chain V:  31% 10% 7% 52%

ALA ARG GLY ARG TPO PRO PRO SEP LEU PRO GLN D355 T359 T360 A361 S362 S363 S364 D368 THR SER SER

• Molecule 5: Beta-arrestin-1

Chain B:  62% 20% 18%

GLY ASP LYS GLY T6 R7 V8 K11 L18 L22 D29 G39 V40 V41 P45 L48 K49 E50 R51 R52 V53 Y63 G64 R65 GLU ASP LEU ASP VAL LEU G72 K77 N83 VAL GLN SER PHE PRO PRO ALA PRO GLU ASP LYS LYS P96 Q101 K107

L108 G109 I119 L123 T128 G137 E145 V146 K147 A148 A151 E152 N153 L154 E155 E156 K157 I158 R165 L166 R169 K170 V171 Q172 P175 P182 E185 T186 T187 R188 Q189 S193 D194 L197 H198 L203 D204 K205 I214 V220 T221 N222 N223 T224

T231 R232 I233 S234 V235 R236 Q237 Y238 F244 N245 T246 V253 A254 M255 E256 F257 A258 D259 P264 F268 T275 P276 N281 R282 E283 K284 R285 G286 G291 L305 L306 G309 L315 V323 K324 V325 K326 S330 ARG GLY LEU LEU GLY ASP LEU ALA

SER S941 D942 V943 A944 V945 E946 L947 M952 H953 P958 LYS GLU PRO PRO HIS ARG GLU VAL PRO GLU SER GLU THR PRO VAL ASP THR ASN LEU ILE GLU LEU ASP THR ASN ASP ASP ILE VAL PHE GLU ASP PHE ALA ARG

• Molecule 6: Antibody fragment Fab30, light chain

Chain L:  38% 10% 52%

SER ASP I3 Q4 T5 T6 Q7 SER PRO S10 A14 S15 D18 T21 I22 T23 C24 S32 S39 Q45 P45 L48 Y66 V69 R67 F72 T73 L74 T75 L79 Q91 Y92 K93 Y94 I107 LYS ARG THR VAL ALA ALA PRO SER VAL PHE THR

PRO PRO SER ASP SER GLN LEU LYS ASP THR ALA VAL THR TYR SER LEU SER SER ILE THR

LEU THR LEU SER LYS ALA ASP TYR LYS HIS VAL TYR CYS VAL GLU VAL THR HIS GLN TYR PRO ARG GLU ALA VAL THR LYS TRP PHE ASN ARG GLY ALA CYS

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	200270	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 ($e^-/\text{\AA}^2$)	58.5	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2400	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	8.027	Depositor
Minimum map value	-4.905	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.048	Depositor
Recommended contour level	0.65	Depositor
Map size ($\text{\AA}$)	414.72, 414.72, 414.72	wwPDB
Map dimensions	300, 300, 300	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	1.3824, 1.3824, 1.3824	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: TPO, SEP

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	0.15	0/859	0.35	0/1165
2	C	0.14	0/451	0.41	0/616
3	H	0.25	0/896	0.46	0/1220
4	V	0.33	0/54	0.36	0/66
5	B	0.20	0/2616	0.37	0/3545
6	L	0.15	0/803	0.36	0/1088
All	All	0.19	0/5679	0.39	0/7700

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	843	0	815	18	0
2	C	439	0	405	11	0
3	H	872	0	810	34	0
4	V	120	0	76	4	0
5	B	2565	0	2632	52	0
6	L	787	0	774	13	0
All	All	5626	0	5512	123	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 11.

All (123) 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)
3:H:55:SER:HB3	3:H:104:GLN:NE2	1.57	1.17
3:H:55:SER:HB3	3:H:104:GLN:HE22	1.12	0.95
5:B:305:LEU:HG	5:B:352:MET:HE1	1.57	0.84
6:L:39:GLN:HB3	6:L:45:PRO:HA	1.66	0.77
3:H:55:SER:CB	3:H:104:GLN:NE2	2.44	0.76
5:B:233:ILE:HG23	5:B:325:VAL:HG12	1.69	0.74
5:B:77:LYS:HE2	5:B:315:LEU:HD12	1.69	0.73
3:H:50:TRP:NE1	3:H:53:SER:OG	2.25	0.69
5:B:153:ASN:HB3	5:B:156:GLU:HG3	1.76	0.67
1:A:35:ALA:HB2	1:A:98:PHE:HD1	1.60	0.66
3:H:36:SER:HB2	3:H:104:GLN:HB2	1.77	0.66
3:H:58:TYR:O	5:B:282:ARG:NH2	2.28	0.66
3:H:86:MET:HE1	3:H:119:VAL:HG11	1.78	0.66
5:B:175:PRO:HD3	5:B:305:LEU:HD11	1.76	0.66
5:B:123:LEU:O	5:B:170:LYS:NZ	2.29	0.65
5:B:235:VAL:HG22	5:B:323:VAL:HG22	1.79	0.65
2:C:87:VAL:HB	2:C:138:ALA:HB3	1.78	0.64
5:B:234:SER:HB3	5:B:255:MET:HG2	1.78	0.63
6:L:22:ILE:HD11	6:L:74:LEU:HD22	1.80	0.63
1:A:39:ARG:NH2	4:V:364:SEP:O3P	2.33	0.61
1:A:38:ARG:NH2	1:A:88:GLU:O	2.34	0.61
5:B:305:LEU:O	5:B:306:LEU:HG	2.02	0.60
1:A:67:PHE:HZ	1:A:82:MET:HE2	1.67	0.60
6:L:91:GLN:OE1	6:L:93:LYS:N	2.35	0.59
3:H:92:GLU:N	3:H:92:GLU:OE1	2.34	0.59
5:B:169:ARG:NH1	5:B:291:GLY:O	2.31	0.58
1:A:70:SER:HB2	1:A:79:TYR:HB2	1.85	0.58
5:B:145:GLU:OE1	5:B:147:LYS:NZ	2.37	0.58
5:B:185:GLU:OE1	5:B:198:HIS:NE2	2.37	0.58
3:H:104:GLN:O	3:H:104:GLN:HG2	2.04	0.57
1:A:88:GLU:OE2	1:A:88:GLU:N	2.32	0.57
1:A:30:ASP:O	1:A:71:ARG:NH2	2.40	0.55
1:A:75:LYS:O	1:A:75:LYS:HD3	2.07	0.55
3:H:60:TYR:OH	5:B:282:ARG:NH1	2.40	0.55
5:B:236:ARG:NH2	5:B:238:TYR:OH	2.37	0.55
3:H:55:SER:CB	3:H:104:GLN:HE22	2.02	0.55
5:B:45:PRO:HA	5:B:48:LEU:HB3	1.89	0.55

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:B:182:PRO:HG2	5:B:203:LEU:HB2	1.89	0.55
1:A:22:CYS:SG	1:A:23:VAL:N	2.80	0.54
3:H:102:SER:HB3	3:H:110:LEU:HA	1.89	0.54
3:H:100:ALA:HB2	3:H:113:TRP:CZ3	2.43	0.54
6:L:18:ASP:H	6:L:79:LEU:HB2	1.73	0.53
1:A:39:ARG:HB2	1:A:45:ARG:HG2	1.90	0.53
6:L:5:MET:SD	6:L:91:GLN:HB3	2.49	0.53
3:H:41:ARG:HD3	3:H:97:TYR:CZ	2.44	0.52
1:A:67:PHE:CZ	1:A:82:MET:HE2	2.44	0.52
2:C:128:GLN:N	2:C:128:GLN:OE1	2.42	0.52
5:B:41:VAL:HG11	5:B:53:VAL:HG11	1.90	0.51
6:L:7:GLN:HA	6:L:24:CYS:HA	1.92	0.51
3:H:43:ALA:HB3	3:H:46:LYS:HB3	1.93	0.51
5:B:22:LEU:HD12	5:B:166:LEU:HD23	1.93	0.50
5:B:222:ASN:OD1	5:B:224:THR:N	2.38	0.50
2:C:136:TYR:OH	5:B:77:LYS:HG2	2.12	0.50
3:H:9:GLU:OE2	3:H:117:THR:OG1	2.20	0.50
3:H:38:HIS:HB2	3:H:100:ALA:HB3	1.92	0.50
5:B:189:GLN:HB3	5:B:193:SER:HB2	1.95	0.49
6:L:48:LEU:O	6:L:56:TYR:N	2.45	0.49
2:C:86:PHE:O	2:C:107:ARG:N	2.47	0.48
2:C:106:GLU:N	2:C:106:GLU:OE2	2.46	0.48
6:L:14:ALA:H	6:L:107:ILE:HB	1.78	0.48
5:B:342:ASP:N	5:B:342:ASP:OD1	2.46	0.48
2:C:100:LEU:HD12	2:C:101:SER:N	2.28	0.48
5:B:50:GLU:N	5:B:50:GLU:OE2	2.47	0.48
5:B:155:GLU:N	5:B:155:GLU:OE2	2.46	0.48
5:B:224:THR:O	5:B:264:PRO:HB3	2.14	0.48
1:A:97:THR:HG22	1:A:99:VAL:H	1.78	0.48
5:B:187:THR:HA	5:B:197:LEU:O	2.14	0.47
3:H:39:TRP:HD1	3:H:73:ILE:HD11	1.79	0.47
5:B:214:ILE:HG21	5:B:323:VAL:HG21	1.97	0.47
1:A:38:ARG:O	1:A:46:GLU:N	2.33	0.47
6:L:91:GLN:HE22	6:L:94:TYR:H	1.63	0.47
3:H:119:VAL:HG12	3:H:119:VAL:O	2.15	0.47
3:H:67:VAL:HG12	3:H:70:ARG:NH2	2.30	0.46
3:H:11:GLY:HA2	3:H:23:LEU:HD13	1.97	0.46
5:B:29:ASP:HB3	5:B:172:GLN:HG2	1.98	0.46
3:H:70:ARG:NH1	3:H:90:ARG:HH12	2.13	0.45
1:A:47:LEU:HD22	1:A:60:ALA:HB3	1.98	0.45
3:H:36:SER:HB2	3:H:104:GLN:CB	2.46	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:133:PRO:HG2	2:C:136:TYR:CD1	2.52	0.45
5:B:220:VAL:HG12	5:B:268:PHE:HB3	1.97	0.45
5:B:283:GLU:N	5:B:283:GLU:OE2	2.49	0.45
5:B:222:ASN:OD1	5:B:223:ASN:N	2.50	0.45
3:H:39:TRP:CZ3	3:H:97:TYR:HB3	2.52	0.45
5:B:18:LEU:HD21	5:B:148:ALA:HB3	1.97	0.45
2:C:90:TYR:HD1	2:C:136:TYR:HD2	1.65	0.44
5:B:77:LYS:HE2	5:B:315:LEU:CD1	2.41	0.44
5:B:52:ARG:NH2	5:B:83:ASN:O	2.42	0.44
6:L:56:TYR:HB3	6:L:59:VAL:HG22	2.00	0.44
6:L:32:SER:HB2	6:L:67:ARG:CZ	2.48	0.43
5:B:326:LYS:HG3	5:B:344:ALA:HB2	2.00	0.43
5:B:63:TYR:CD2	5:B:244:PHE:HE2	2.36	0.43
5:B:11:LYS:NZ	5:B:165:ARG:O	2.29	0.43
5:B:119:ILE:HG23	5:B:123:LEU:HD12	2.00	0.43
5:B:235:VAL:HG11	5:B:276:PRO:HG3	1.99	0.43
3:H:33:TYR:HB2	5:B:275:THR:HG21	2.01	0.43
3:H:67:VAL:HG12	3:H:70:ARG:HH21	1.84	0.42
4:V:363:SEP:N	5:B:8:VAL:O	2.53	0.42
5:B:305:LEU:O	5:B:306:LEU:CG	2.67	0.42
3:H:36:SER:HB2	3:H:104:GLN:CG	2.49	0.42
3:H:38:HIS:O	3:H:100:ALA:N	2.47	0.41
6:L:32:SER:HA	6:L:72:PHE:CE1	2.55	0.41
3:H:105:PHE:O	3:H:106:TRP:C	2.63	0.41
5:B:231:ILE:HG23	5:B:258:ALA:HB3	2.01	0.41
5:B:285:ARG:HD2	5:B:286:GLY:N	2.36	0.41
3:H:70:ARG:HD2	3:H:88:SER:O	2.20	0.41
3:H:86:MET:SD	3:H:89:LEU:HD21	2.60	0.41
5:B:204:ASP:OD1	5:B:205:LYS:HG2	2.21	0.41
3:H:38:HIS:CD2	3:H:102:SER:OG	2.73	0.41
5:B:39:GLY:N	5:B:101:GLN:HE22	2.19	0.41
5:B:128:THR:HG22	5:B:291:GLY:HA2	2.02	0.41
1:A:90:THR:HG23	1:A:111:THR:HA	2.02	0.41
5:B:253:VAL:HG13	5:B:281:ASN:OD1	2.21	0.41
1:A:45:ARG:NH2	4:V:363:SEP:O1P	2.54	0.41
1:A:64:LYS:HA	1:A:64:LYS:HD2	1.82	0.41
5:B:346:GLU:C	5:B:347:LEU:HD12	2.46	0.41
2:C:120:LEU:HG	2:C:131:TYR:CE1	2.56	0.40
2:C:136:TYR:CD1	2:C:136:TYR:N	2.90	0.40
2:C:91:ASP:OD1	2:C:103:LYS:HG2	2.21	0.40
4:V:364:SEP:O	5:B:107:LYS:NZ	2.32	0.40

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:106:TRP:HZ2	5:B:353:HIS:HB3	1.87	0.40
1:A:39:ARG:O	1:A:91:ALA:HB1	2.22	0.40
3:H:15:VAL:HG12	3:H:121:VAL:HG21	2.03	0.40
6:L:21:THR:HG22	6:L:75:THR:OG1	2.21	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	109/124 (88%)	102 (94%)	7 (6%)	0	100	100
2	C	53/85 (62%)	50 (94%)	3 (6%)	0	100	100
3	H	112/237 (47%)	108 (96%)	4 (4%)	0	100	100
4	V	6/29 (21%)	5 (83%)	1 (17%)	0	100	100
5	B	315/392 (80%)	304 (96%)	11 (4%)	0	100	100
6	L	99/215 (46%)	88 (89%)	11 (11%)	0	100	100
All	All	694/1082 (64%)	657 (95%)	37 (5%)	0	100	100

There are no Ramachandran outliers to report.

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	90/103 (87%)	89 (99%)	1 (1%)	65	74
2	C	48/76 (63%)	48 (100%)	0	100	100
3	H	89/200 (44%)	89 (100%)	0	100	100
4	V	6/16 (38%)	6 (100%)	0	100	100
5	B	287/349 (82%)	287 (100%)	0	100	100
6	L	89/190 (47%)	89 (100%)	0	100	100
All	All	609/934 (65%)	608 (100%)	1 (0%)	85	84

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	29	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	73	ASN
1	A	106	GLN
5	B	30	HIS
6	L	28	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

6 non-standard protein/DNA/RNA residues are 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
4	TPO	V	359	4	8,10,11	0.81	0	10,14,16	1.01	1 (10%)
4	SEP	V	362	4	8,9,10	1.58	1 (12%)	7,12,14	1.26	1 (14%)
4	SEP	V	364	4	8,9,10	1.57	1 (12%)	7,12,14	1.05	0
4	SEP	V	357	4	8,9,10	0.68	0	7,12,14	0.65	0
4	SEP	V	363	4	8,9,10	1.57	1 (12%)	7,12,14	1.24	1 (14%)
4	TPO	V	360	4	8,10,11	0.79	0	10,14,16	1.05	1 (10%)

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 Torsions and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	TPO	V	359	4	-	2/9/11/13	-
4	SEP	V	362	4	-	0/6/8/10	-
4	SEP	V	364	4	-	2/6/8/10	-
4	SEP	V	357	4	-	5/6/8/10	-
4	SEP	V	363	4	-	2/6/8/10	-
4	TPO	V	360	4	-	1/9/11/13	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	V	362	SEP	P-O1P	3.47	1.61	1.50
4	V	364	SEP	P-O1P	3.46	1.61	1.50
4	V	363	SEP	P-O1P	3.45	1.61	1.50

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	V	362	SEP	OG-CB-CA	2.63	110.70	108.14
4	V	363	SEP	OG-CB-CA	2.56	110.63	108.14
4	V	360	TPO	O-C-CA	-2.55	118.22	124.77
4	V	359	TPO	O-C-CA	-2.53	118.28	124.77

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	V	357	SEP	N-CA-CB-OG
4	V	357	SEP	C-CA-CB-OG
4	V	357	SEP	CB-OG-P-O1P
4	V	357	SEP	CB-OG-P-O2P
4	V	357	SEP	CB-OG-P-O3P
4	V	363	SEP	N-CA-CB-OG
4	V	364	SEP	N-CA-CB-OG
4	V	364	SEP	C-CA-CB-OG
4	V	363	SEP	CB-OG-P-O1P
4	V	360	TPO	CB-OG1-P-O3P
4	V	359	TPO	C-CA-CB-CG2
4	V	359	TPO	CB-OG1-P-O1P

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	V	364	SEP	2	0
4	V	363	SEP	2	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

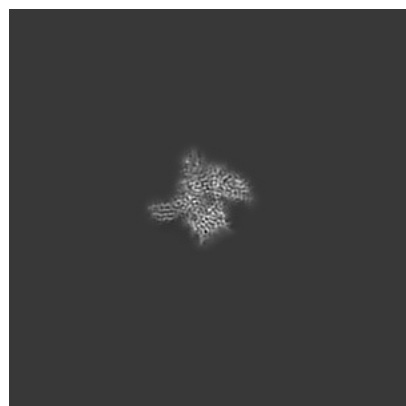
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-45977. These allow visual inspection of the internal detail of the map and identification of artifacts.

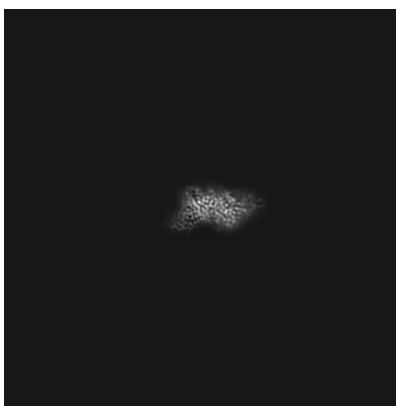
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

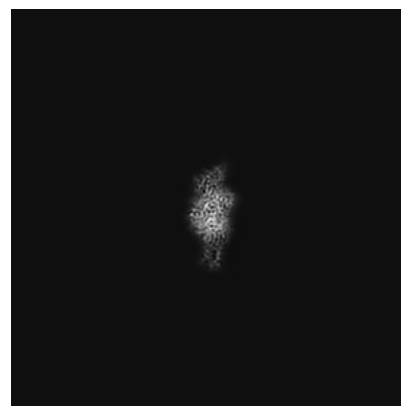
6.1.1 Primary map



X

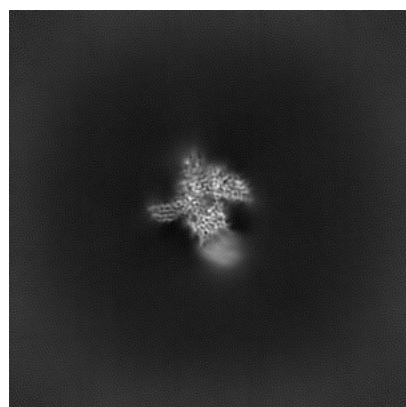


Y

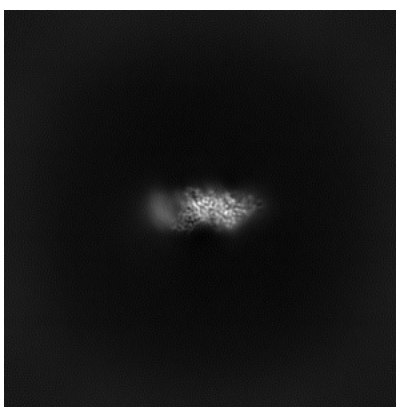


Z

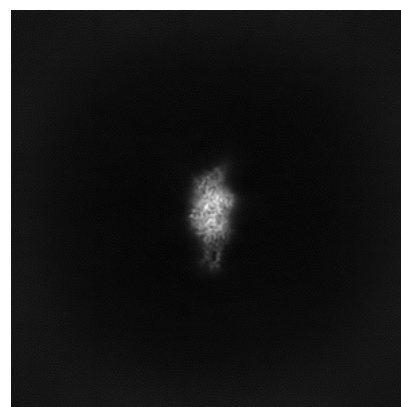
6.1.2 Raw map



X



Y

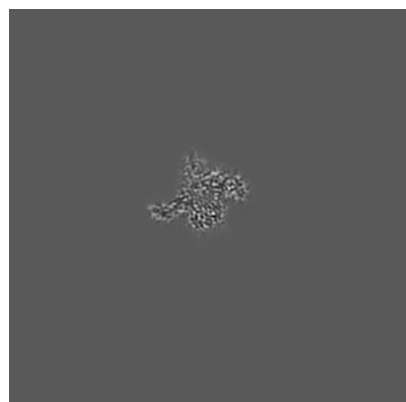


Z

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

6.2 Central slices [i](#)

6.2.1 Primary map



X Index: 150

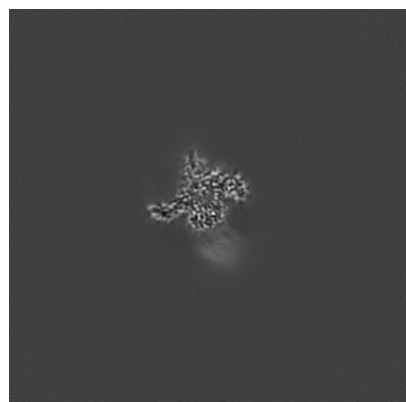


Y Index: 150

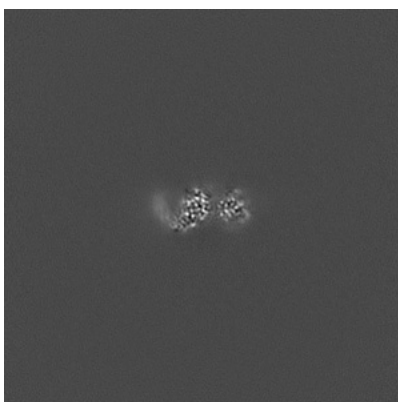


Z Index: 150

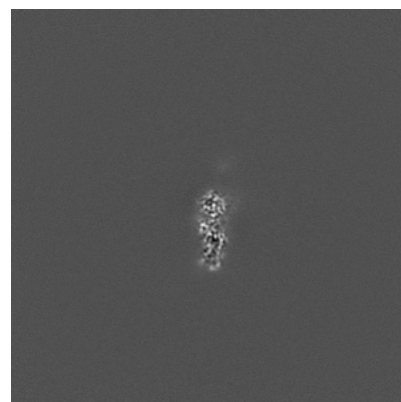
6.2.2 Raw map



X Index: 150



Y Index: 150

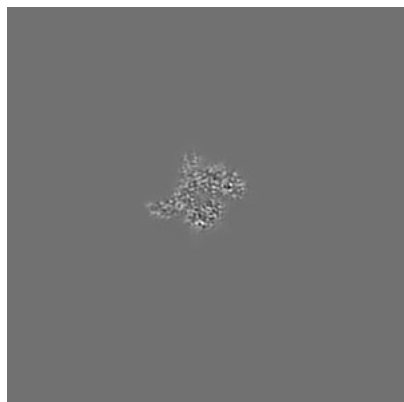


Z Index: 150

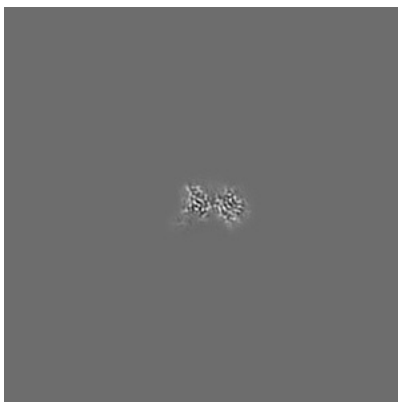
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

6.3.1 Primary map



X Index: 149

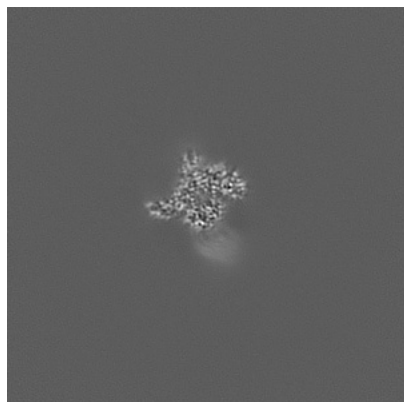


Y Index: 154

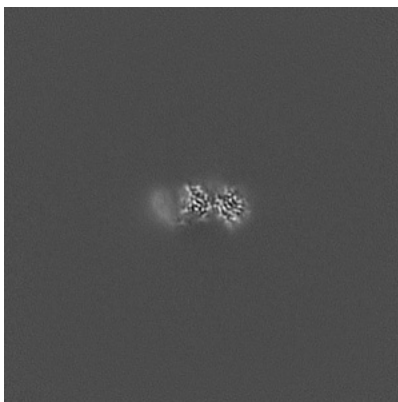


Z Index: 146

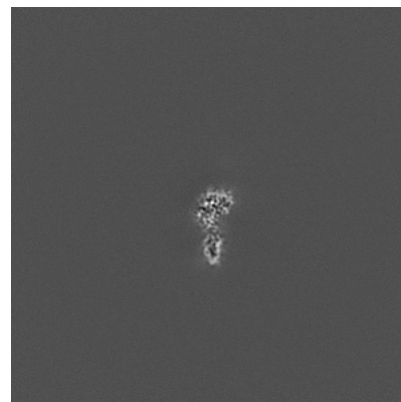
6.3.2 Raw map



X Index: 149



Y Index: 154

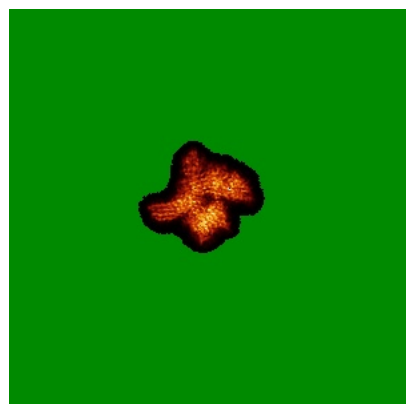


Z Index: 146

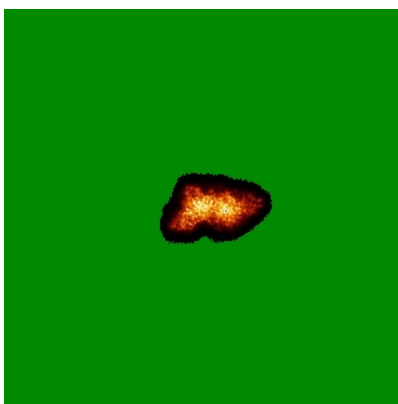
The images above show the largest variance slices of the map in three orthogonal directions.

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

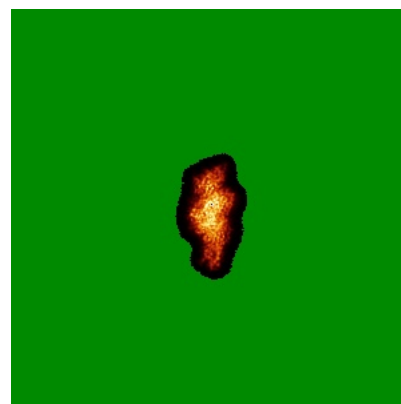
6.4.1 Primary map



X

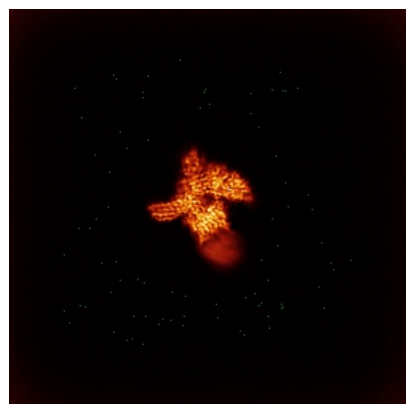


Y

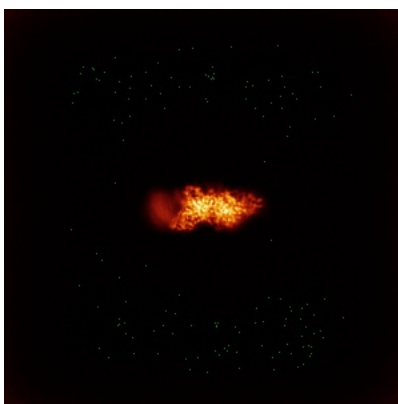


Z

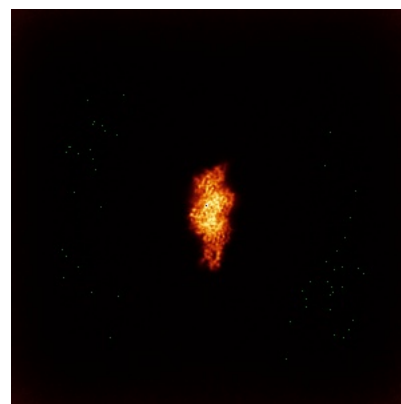
6.4.2 Raw map



X



Y

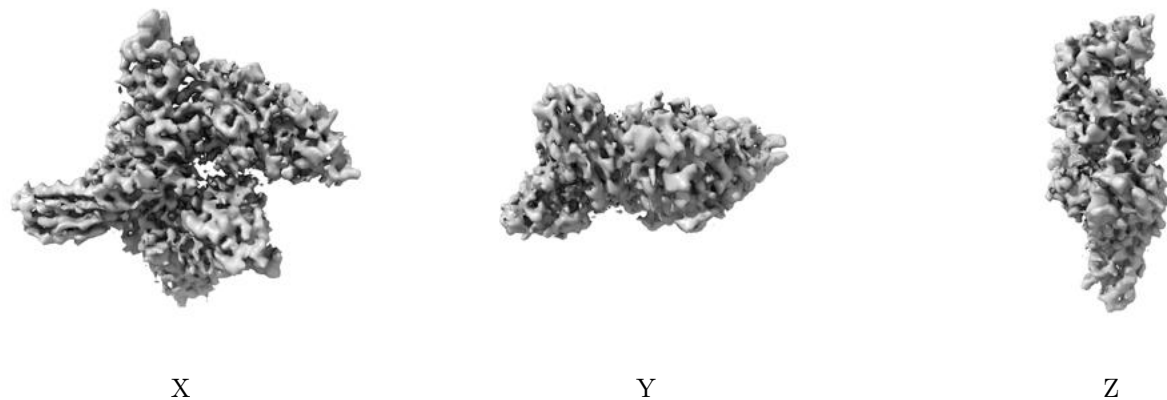


Z

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.

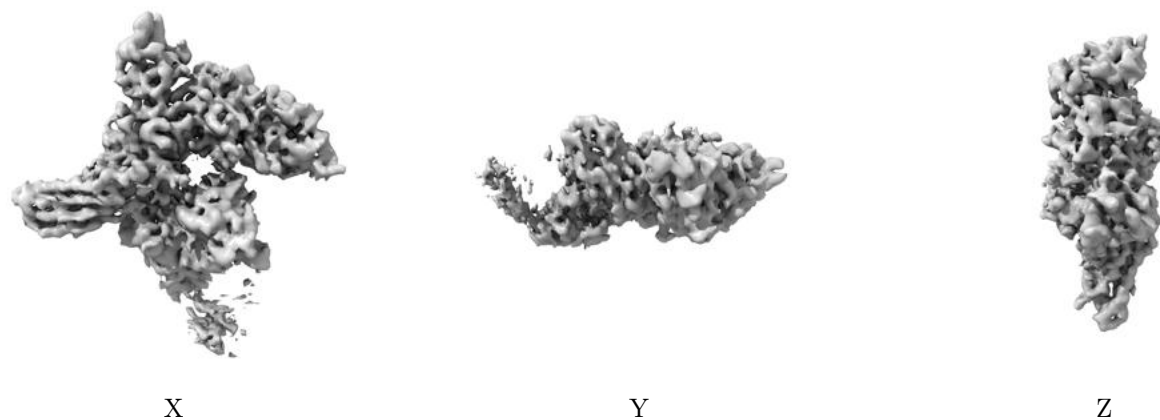
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.65. 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.

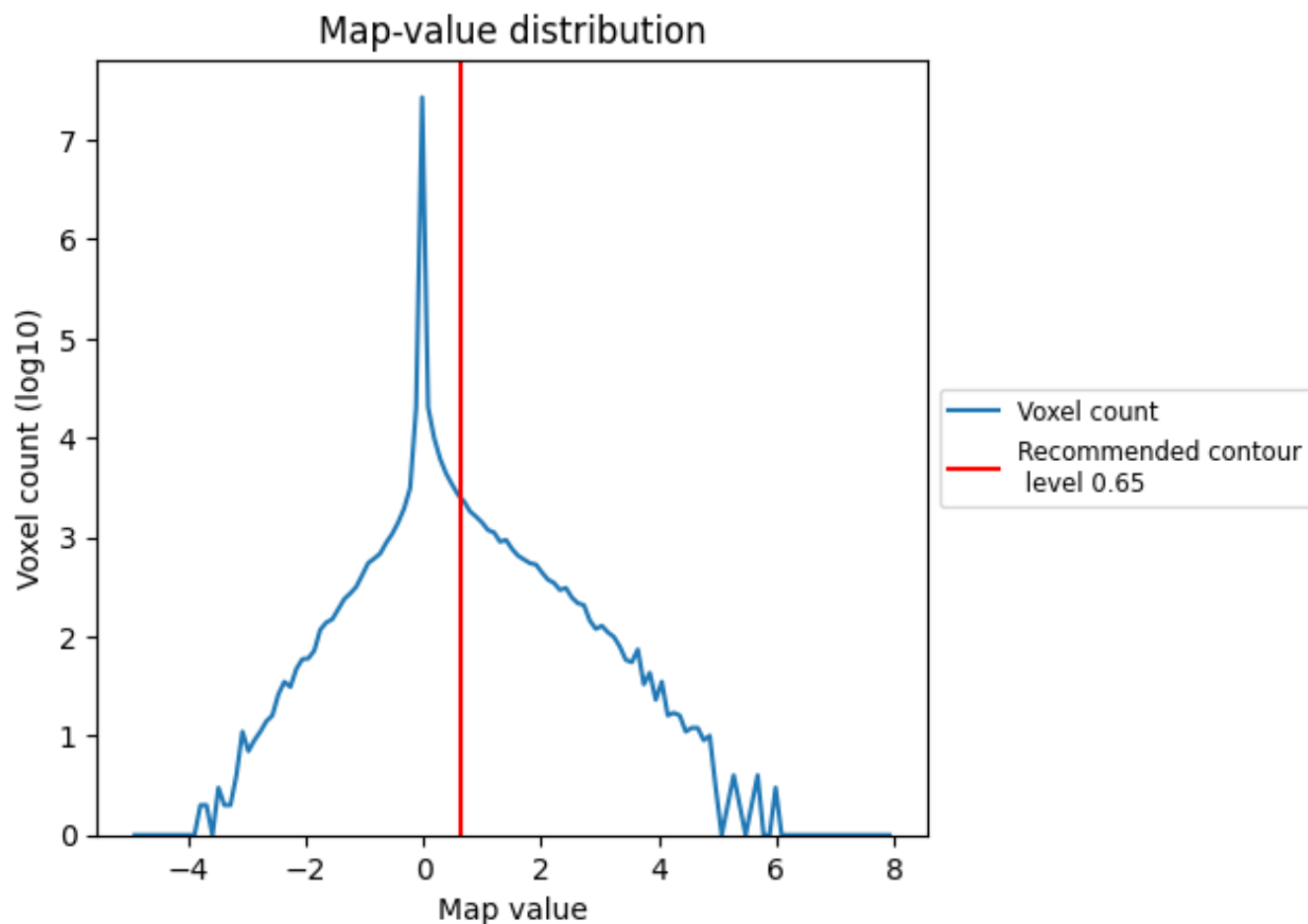
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

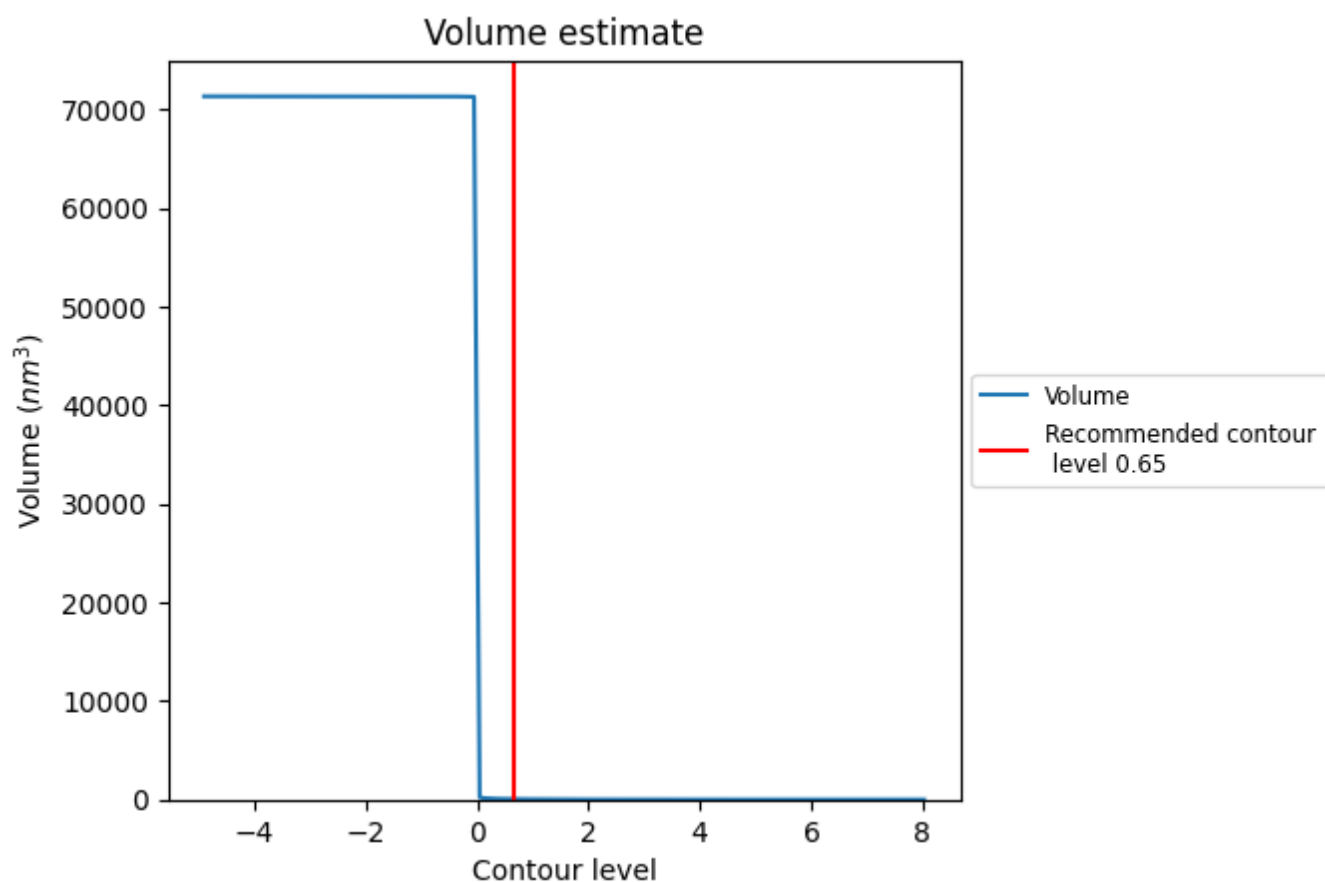
This section contains the results of statistical analysis of the map.

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.

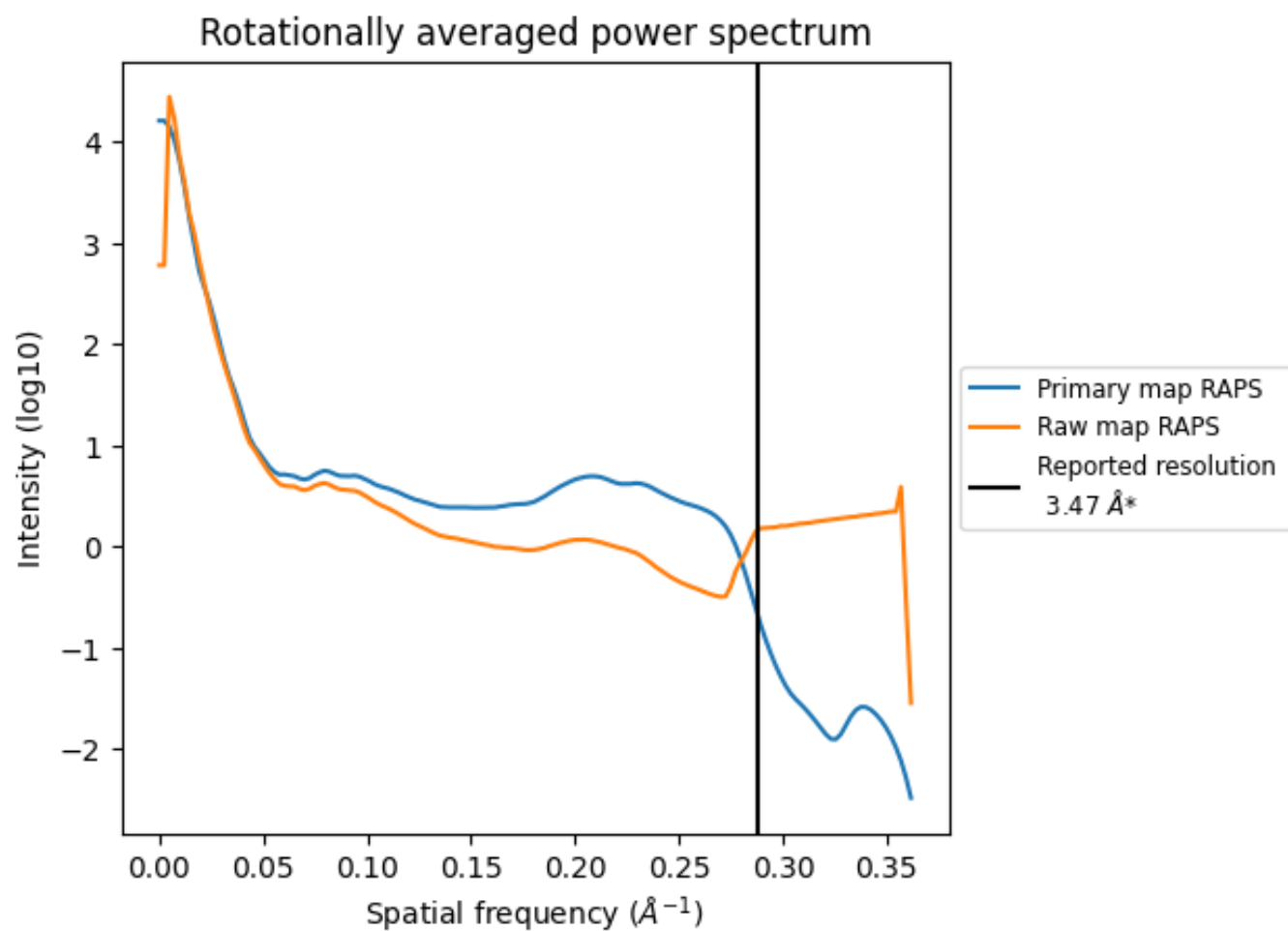
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 50 nm^3 ; this corresponds to an approximate mass of 45 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 ⓘ

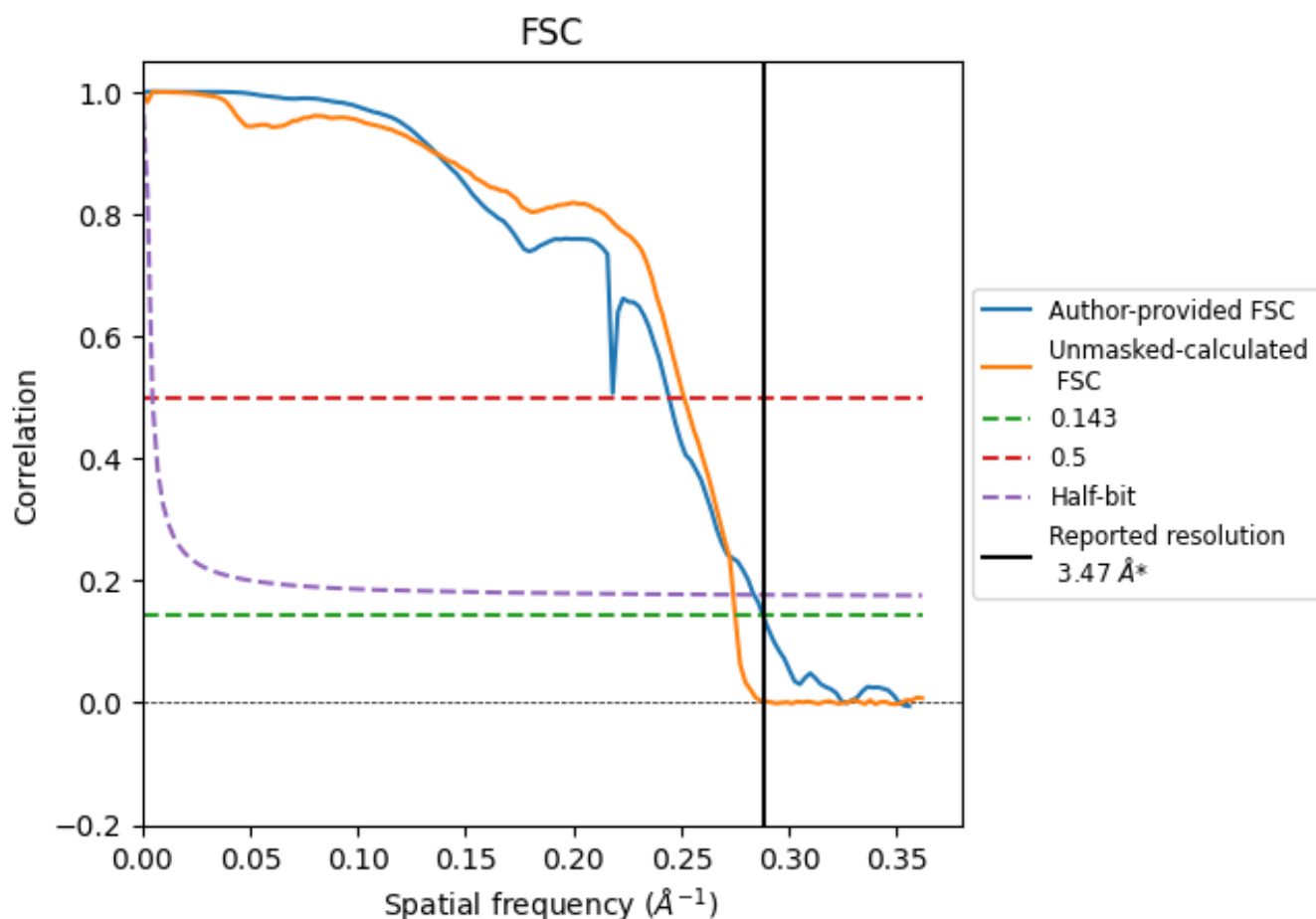


*Reported resolution corresponds to spatial frequency of 0.288 Å⁻¹

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 $0.288 \text{ \AA}^{-1}$

8.2 Resolution estimates [i](#)

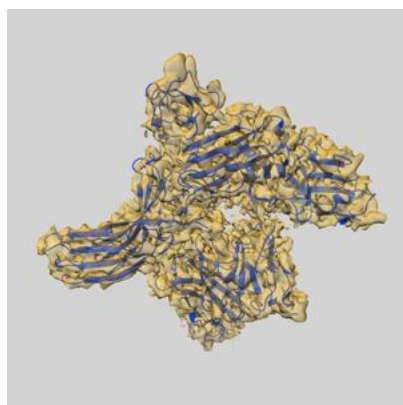
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.47	-	-
Author-provided FSC curve	3.47	4.09	3.52
Unmasked-calculated*	3.64	3.98	3.65

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

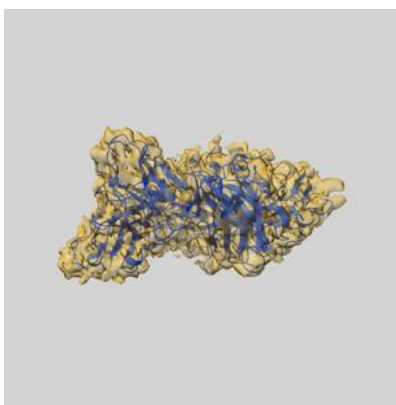
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-45977 and PDB model 9CX3. Per-residue inclusion information can be found in section [3](#) on page [7](#).

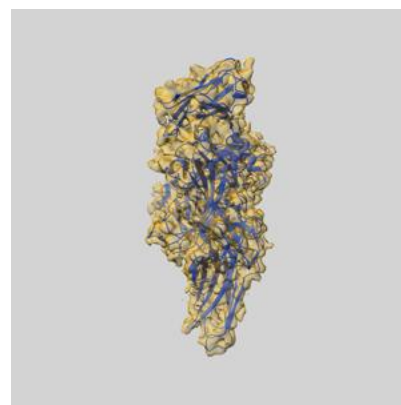
9.1 Map-model overlay [i](#)



X



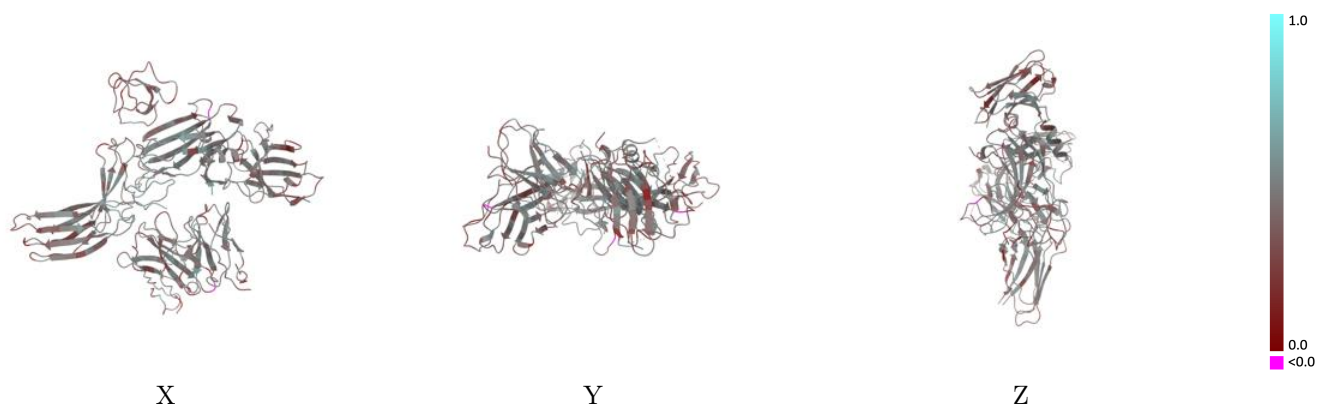
Y



Z

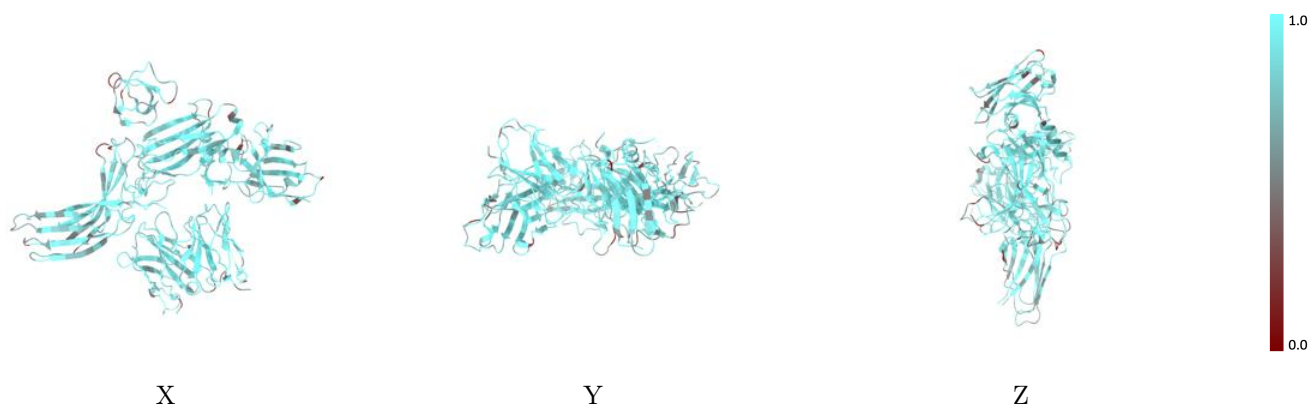
The images above show the 3D surface view of the map at the recommended contour level 0.65 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



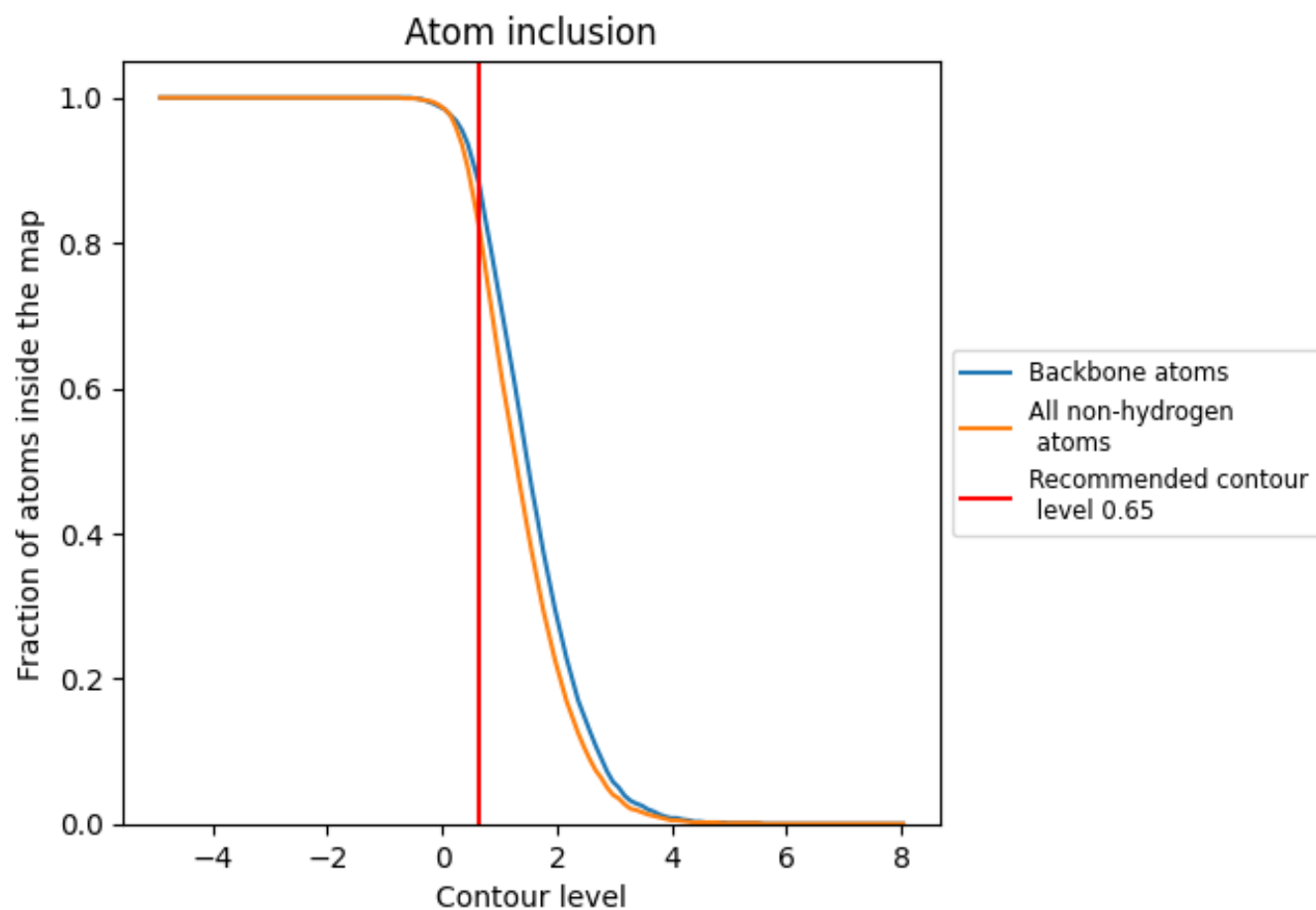
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.65).

9.4 Atom inclusion [i](#)



At the recommended contour level, 88% of all backbone atoms, 82% 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 (0.65) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8200	0.4230
A	0.8090	0.4040
B	0.8130	0.4430
C	0.7050	0.3630
H	0.8450	0.3980
L	0.8900	0.4410
V	0.8330	0.4130

1.0

0.0

<0.0