



Full wwPDB EM Validation Report ⓘ

Mar 20, 2026 – 06:16 AM UTC

PDB ID : 8QR1 / pdb_00008qr1
EMDB ID : EMD-18611
Title : Cryo-EM structure of the human Tip60 complex
Authors : Li, C.; Smirnova, E.; Schnitzler, C.; Crucifix, C.; Concordet, J.P.; Brion, A.;
Poterszman, A.; Schultz, P.; Papai, G.; Ben-Shem, A.
Deposited on : 2023-10-06
Resolution : 2.40 Å(reported)

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

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

A user guide is available at

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

The types of validation reports are described at

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

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

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

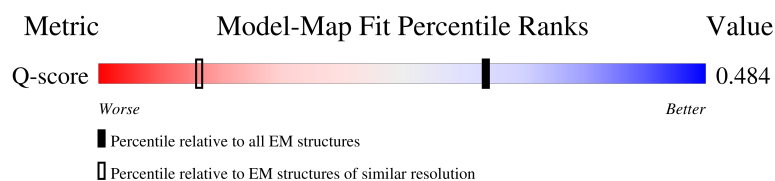
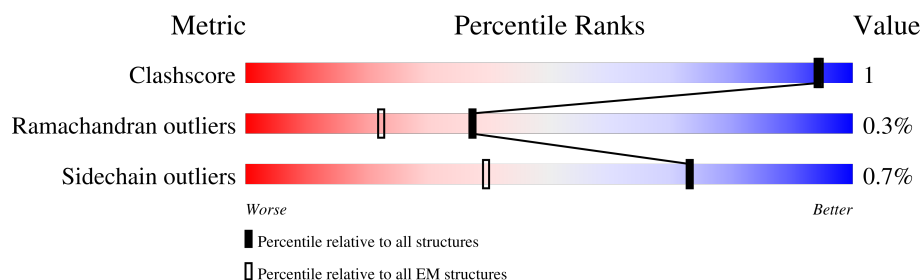
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 2.40 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	5628 (1.90 - 2.90)

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	3159	 8% 27% 72%
2	C	836	 9% 90%
3	F	467	 43% 54%
4	S	364	 26% 70%

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Mol	Chain	Length	Quality of chain
5	B	375	9% 96% .
5	G	375	94% 6% .
6	K	429	92% 5% .
7	E	456	94% . .
7	I	456	91% 5% .
7	L	456	5% 91% 5% .
8	D	463	90% 6% .
8	H	463	6% 91% 7% .
8	J	463	90% 5% 5% .

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 39353 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 E1A-binding protein p400.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	875	Total	C	N	O	S	0	0
			7059	4501	1270	1258	30		

- Molecule 2 is a protein called Enhancer of polycomb homolog 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	C	85	Total	C	N	O	S	0	0
			687	432	134	118	3		

- Molecule 3 is a protein called DNA methyltransferase 1-associated protein 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	F	213	Total	C	N	O	S	0	0
			1781	1138	322	317	4		

- Molecule 4 is a protein called Vacuolar protein sorting-associated protein 72 homolog.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	S	108	Total	C	N	O	S	0	0
			823	534	142	145	2		

- Molecule 5 is a protein called Actin, cytoplasmic 1, N-terminally processed.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	B	375	Total	C	N	O	S	0	0
			2925	1850	491	561	23		
5	G	354	Total	C	N	O	S	0	0
			2764	1755	459	530	20		

- Molecule 6 is a protein called Actin-like protein 6A.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	K	406	Total	C	N	O	S	0	0
			3164	2001	538	601	24		

- Molecule 7 is a protein called RuvB-like 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	L	439	Total	C	N	O	S	0	0
			3374	2124	579	654	17		
7	I	434	Total	C	N	O	S	0	0
			3329	2096	573	644	16		
7	E	440	Total	C	N	O	S	0	0
			3383	2129	581	656	17		

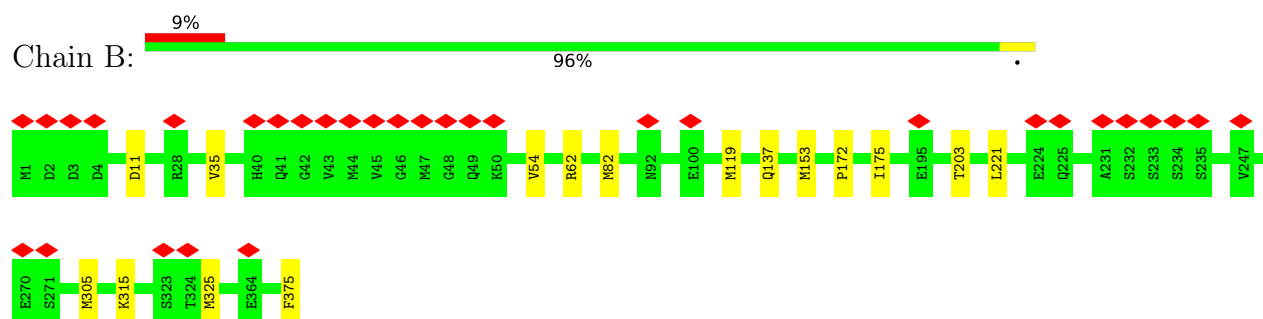
- Molecule 8 is a protein called RuvB-like 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	H	429	Total	C	N	O	S	0	0
			3299	2065	575	644	15		
8	D	434	Total	C	N	O	S	0	0
			3362	2102	590	655	15		
8	J	439	Total	C	N	O	S	0	0
			3403	2125	597	665	16		

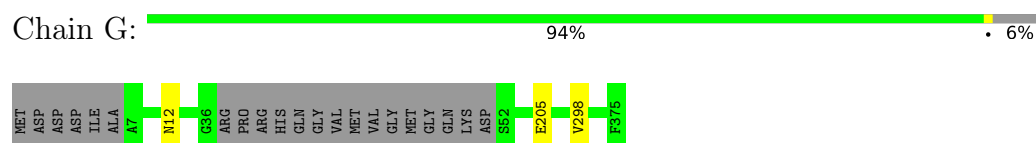
[illegible]



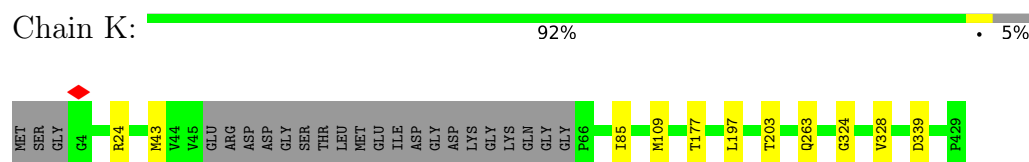
- Molecule 5: Actin, cytoplasmic 1, N-terminally processed



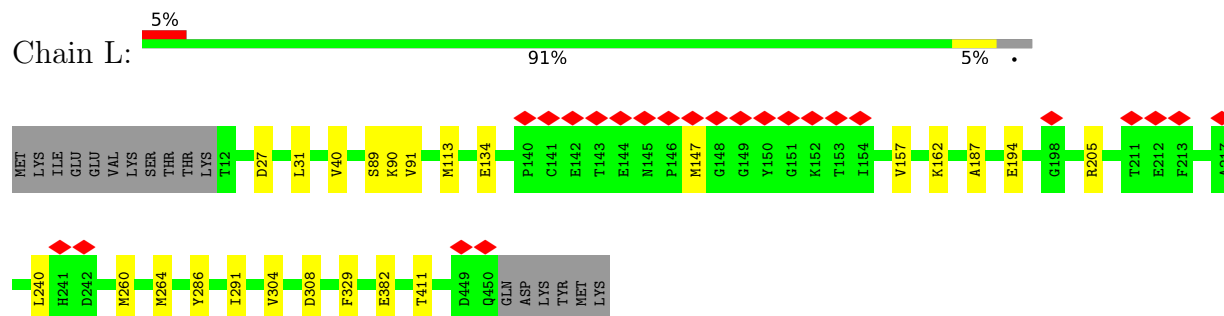
- Molecule 5: Actin, cytoplasmic 1, N-terminally processed



- Molecule 6: Actin-like protein 6A

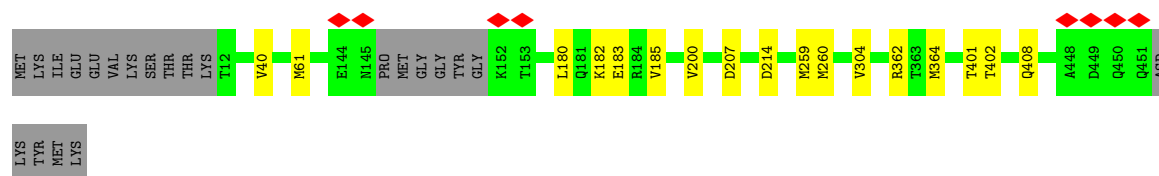


- Molecule 7: RuvB-like 1

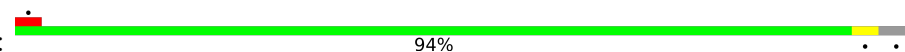


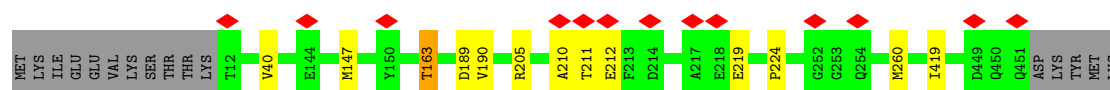
- Molecule 7: RuvB-like 1

Chain I:  91% 5%

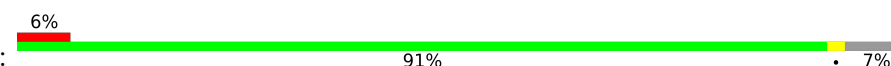


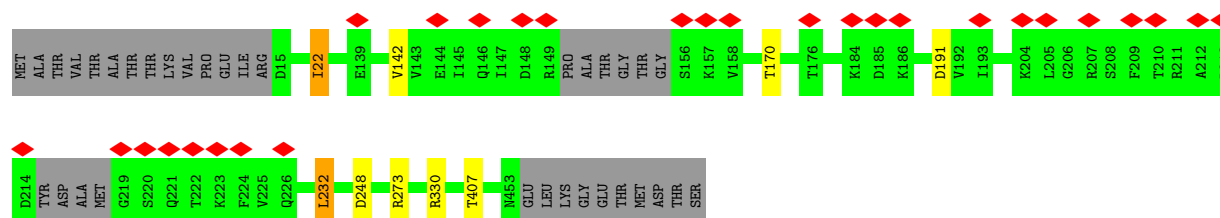
- Molecule 7: RuvB-like 1

Chain E:  94% 5%



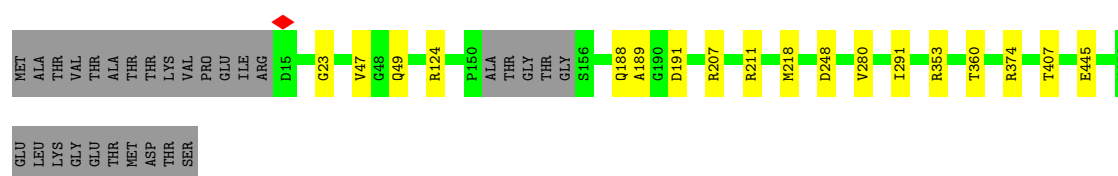
- Molecule 8: RuvB-like 2

Chain H:  91% 6% 7%

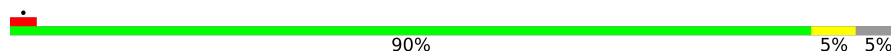


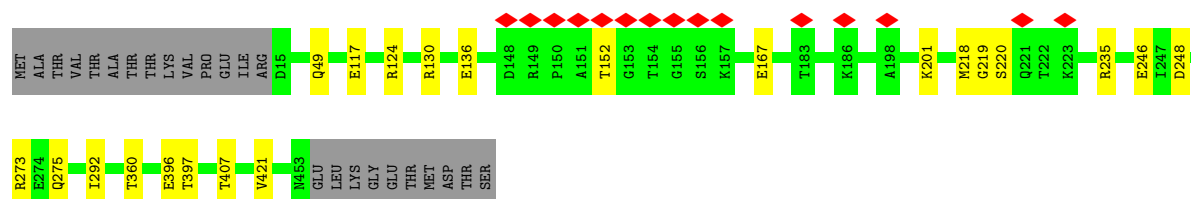
- Molecule 8: RuvB-like 2

Chain D:  90% 6%



- Molecule 8: RuvB-like 2

Chain J:  90% 5% 5%



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	180397	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	40, 52	Depositor
Minimum defocus (nm)	1200	Depositor
Maximum defocus (nm)	2800	Depositor
Magnification	Not provided	
Image detector	FEI FALCON IV (4k x 4k), GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	74.696	Depositor
Minimum map value	-28.889	Depositor
Average map value	-0.003	Depositor
Map value standard deviation	1.069	Depositor
Recommended contour level	6	Depositor
Map size ($\text{\AA}$)	500.78, 500.78, 500.78	wwPDB
Map dimensions	686, 686, 686	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	0.73, 0.73, 0.73	Depositor

5 Model quality

5.1 Standard geometry

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	0.22	0/7210	0.45	0/9766
2	C	0.29	0/709	0.57	0/961
3	F	0.28	0/1830	0.56	0/2472
4	S	0.34	0/855	0.71	1/1177 (0.1%)
5	B	0.22	0/2988	0.43	1/4045 (0.0%)
5	G	0.25	0/2824	0.44	0/3825
6	K	0.27	0/3236	0.45	0/4387
7	E	0.27	0/3428	0.55	1/4622 (0.0%)
7	I	0.24	0/3371	0.49	0/4545
7	L	0.25	0/3419	0.50	0/4610
8	D	0.24	0/3401	0.47	0/4580
8	H	0.24	0/3335	0.49	0/4492
8	J	0.26	0/3443	0.52	0/4636
All	All	0.25	0/40049	0.49	3/54118 (0.0%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
2	C	0	1
4	S	0	1
All	All	0	2

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	119	MET	CA-CB-CG	5.35	124.80	114.10
7	E	212	GLU	CA-CB-CG	5.33	124.75	114.10
4	S	343	PRO	N-CA-C	5.25	117.10	110.70

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	C	449	GLN	Peptide
4	S	307	ASP	Peptide

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	7059	0	7067	23	0
2	C	687	0	654	4	0
3	F	1781	0	1727	7	0
4	S	823	0	798	4	0
5	B	2925	0	2891	9	0
5	G	2764	0	2732	0	0
6	K	3164	0	3105	6	0
7	E	3383	0	3482	5	0
7	I	3329	0	3418	11	0
7	L	3374	0	3474	10	0
8	D	3362	0	3414	11	0
8	H	3299	0	3337	4	0
8	J	3403	0	3472	11	0
All	All	39353	0	39571	91	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 1.

All (91) 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)
5:B:62:ARG:O	5:B:203:THR:OG1	2.11	0.69
8:J:167:GLU:N	8:J:167:GLU:OE1	2.28	0.66
7:L:147:MET:SD	7:L:147:MET:N	2.71	0.63
5:B:305:MET:SD	5:B:305:MET:N	2.73	0.61
1:A:1205:LEU:O	1:A:1209:THR:OG1	2.18	0.61
7:I:362:ARG:NH2	7:I:364:MET:SD	2.75	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:E:189:ASP:OD1	7:E:205:ARG:NH1	2.34	0.60
5:B:11:ASP:OD1	5:B:137:GLN:NE2	2.34	0.60
1:A:1752:GLN:OE1	8:J:201:LYS:NZ	2.26	0.59
8:D:211:ARG:NH1	8:D:218:MET:O	2.36	0.59
8:H:170:THR:HG23	8:H:232:LEU:HD22	1.83	0.58
8:J:124:ARG:NE	8:J:248:ASP:OD2	2.36	0.58
5:B:221:LEU:O	5:B:315:LYS:NZ	2.27	0.57
1:A:1170:SER:O	1:A:1181:LYS:NZ	2.36	0.57
8:D:280:VAL:CG1	8:D:291:ILE:HD11	2.36	0.56
8:H:22:ILE:HD11	7:I:61:MET:HE1	1.87	0.54
7:I:183:GLU:HB2	7:I:185:VAL:HG23	1.91	0.53
3:F:51:GLU:N	3:F:51:GLU:OE2	2.41	0.53
8:J:218:MET:SD	8:J:219:GLY:N	2.82	0.52
1:A:824:THR:HG21	1:A:2248:TRP:CH2	2.45	0.51
7:I:180:LEU:HD23	7:I:200:VAL:HG11	1.92	0.51
6:K:203:THR:OG1	6:K:263:GLN:NE2	2.44	0.51
8:D:124:ARG:NE	8:D:248:ASP:OD2	2.39	0.51
7:L:187:ALA:O	7:L:205:ARG:NH1	2.43	0.50
1:A:1233:GLU:O	1:A:1237:THR:OG1	2.28	0.50
1:A:1877:ARG:O	7:I:182:LYS:NZ	2.42	0.50
7:I:401:THR:O	7:I:402:THR:OG1	2.27	0.50
1:A:1092:ARG:NH1	1:A:1313:GLU:OE2	2.45	0.49
8:D:280:VAL:HG11	8:D:291:ILE:HD11	1.95	0.49
3:F:226:ARG:NH2	6:K:339:ASP:OD1	2.45	0.49
5:B:35:VAL:HG22	5:B:54:VAL:HG22	1.95	0.48
7:L:90:LYS:HD2	7:L:91:VAL:HG23	1.95	0.48
8:J:130:ARG:HG3	8:J:292:ILE:HD11	1.96	0.48
7:I:260:MET:HA	7:I:260:MET:HE2	1.96	0.48
3:F:64:ASP:OD1	3:F:65:LYS:N	2.47	0.48
1:A:1838:VAL:O	1:A:1839:VAL:HG12	2.14	0.47
8:J:218:MET:SD	8:J:220:SER:N	2.87	0.47
7:L:286:TYR:CD2	7:L:291:ILE:HD11	2.50	0.47
1:A:1915:ARG:NE	1:A:1986:ASP:OD2	2.48	0.46
8:D:49:GLN:NE2	8:D:360:THR:O	2.48	0.46
4:S:214:TYR:OH	8:D:188:GLN:OE1	2.34	0.46
4:S:270:ASP:O	4:S:272:ALA:N	2.47	0.46
8:J:49:GLN:NE2	8:J:360:THR:O	2.47	0.46
7:L:382:GLU:OE1	7:L:411:THR:OG1	2.33	0.46
1:A:1877:ARG:HA	1:A:1880:THR:HG22	1.97	0.46
1:A:1934:LEU:CD1	1:A:1970:ILE:HD11	2.46	0.46
5:B:153:MET:HE3	5:B:153:MET:C	2.41	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:147:LEU:HD13	3:F:148:HIS:N	2.30	0.45
7:L:134:GLU:N	7:L:162:LYS:O	2.48	0.45
1:A:1749:ARG:NH2	8:J:246:GLU:OE2	2.50	0.45
1:A:1934:LEU:HD12	1:A:1970:ILE:HD11	1.99	0.45
6:K:43:MET:HE2	6:K:85:ILE:HD13	1.97	0.45
7:L:27:ASP:OD1	7:L:31:LEU:N	2.50	0.45
8:D:191:ASP:OD1	8:D:207:ARG:NE	2.46	0.45
7:E:147:MET:HE3	7:E:147:MET:H	1.82	0.44
7:E:210:ALA:O	7:E:211:THR:OG1	2.31	0.44
8:J:117:GLU:OE2	8:J:273:ARG:NH2	2.50	0.44
1:A:1133:LEU:O	1:A:1137:GLU:N	2.50	0.44
1:A:1941:TYR:OH	7:I:207:ASP:OD1	2.21	0.44
7:E:163:THR:OG1	7:E:224:PRO:O	2.35	0.44
7:L:304:VAL:HG21	7:L:329:PHE:CD1	2.53	0.43
1:A:862:GLU:OE1	1:A:865:ARG:NH2	2.50	0.43
8:D:189:ALA:O	8:D:207:ARG:NH2	2.50	0.43
7:I:408:GLN:NE2	8:D:353:ARG:O	2.49	0.43
1:A:1965:ARG:NH1	7:I:214:ASP:OD1	2.49	0.43
1:A:2052:MET:HE2	1:A:2052:MET:N	2.34	0.43
5:B:82:MET:SD	5:B:82:MET:N	2.91	0.43
5:B:172:PRO:HA	5:B:175:ILE:HD12	2.01	0.43
8:J:136:GLU:OE2	8:J:235:ARG:NH2	2.51	0.43
7:L:260:MET:HE2	7:E:260:MET:SD	2.59	0.43
1:A:866:LEU:HD13	5:B:375:PHE:CE1	2.54	0.42
7:L:113:MET:HA	7:L:113:MET:HE2	2.01	0.42
3:F:147:LEU:HD13	3:F:148:HIS:H	1.83	0.42
3:F:98:MET:HE2	3:F:120:GLU:HA	2.01	0.42
2:C:397:ASP:OD2	2:C:397:ASP:N	2.51	0.42
2:C:430:ASP:N	2:C:430:ASP:OD2	2.53	0.42
8:H:248:ASP:OD2	8:H:273:ARG:NH1	2.51	0.42
4:S:308:ILE:HD11	4:S:319:ILE:HD12	2.01	0.42
8:J:396:GLU:HG3	8:J:397:THR:HG23	2.01	0.41
1:A:1133:LEU:O	1:A:1138:GLY:N	2.51	0.41
2:C:425:TRP:O	2:C:426:THR:OG1	2.36	0.41
4:S:218:THR:CG2	8:D:189:ALA:HB1	2.50	0.41
6:K:177:THR:OG1	6:K:197:LEU:O	2.37	0.41
8:D:23:GLY:O	8:D:374:ARG:NH2	2.47	0.41
6:K:324:GLY:O	6:K:328:VAL:HG23	2.21	0.41
1:A:821:TRP:NE1	2:C:459:VAL:O	2.54	0.41
1:A:1880:THR:HG21	7:I:259:MET:HG2	2.03	0.41
6:K:109:MET:HE2	6:K:109:MET:HA	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:H:142:VAL:N	8:H:191:ASP:O	2.47	0.40
3:F:195:TYR:HA	3:F:198:ILE:HG22	2.02	0.40
1:A:1826:ASP:OD1	1:A:1830:ARG:NH1	2.51	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	865/3159 (27%)	821 (95%)	40 (5%)	4 (0%)	24	37
2	C	83/836 (10%)	75 (90%)	7 (8%)	1 (1%)	10	16
3	F	211/467 (45%)	192 (91%)	18 (8%)	1 (0%)	24	37
4	S	104/364 (29%)	84 (81%)	17 (16%)	3 (3%)	3	3
5	B	373/375 (100%)	367 (98%)	6 (2%)	0	100	100
5	G	350/375 (93%)	342 (98%)	8 (2%)	0	100	100
6	K	402/429 (94%)	397 (99%)	5 (1%)	0	100	100
7	E	438/456 (96%)	418 (95%)	19 (4%)	1 (0%)	43	58
7	I	430/456 (94%)	423 (98%)	7 (2%)	0	100	100
7	L	437/456 (96%)	423 (97%)	13 (3%)	1 (0%)	43	58
8	D	430/463 (93%)	418 (97%)	12 (3%)	0	100	100
8	H	423/463 (91%)	407 (96%)	14 (3%)	2 (0%)	24	37
8	J	437/463 (94%)	421 (96%)	15 (3%)	1 (0%)	43	58
All	All	4983/8762 (57%)	4788 (96%)	181 (4%)	14 (0%)	37	50

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
8	H	330	ARG
1	A	1839	VAL
1	A	1751	ASP
4	S	271	ASP
8	H	232	LEU
7	E	219	GLU
1	A	2238	VAL
8	J	152	THR
2	C	450	ARG
3	F	188	VAL
4	S	212	ILE
7	L	157	VAL
4	S	345	PRO
1	A	1887	PHE

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	749/2663 (28%)	748 (100%)	1 (0%)	88	95
2	C	70/738 (10%)	70 (100%)	0	100	100
3	F	182/400 (46%)	181 (100%)	1 (0%)	81	91
4	S	86/312 (28%)	84 (98%)	2 (2%)	44	66
5	B	318/318 (100%)	317 (100%)	1 (0%)	86	93
5	G	301/318 (95%)	298 (99%)	3 (1%)	68	84
6	K	347/364 (95%)	346 (100%)	1 (0%)	86	93
7	E	371/387 (96%)	367 (99%)	4 (1%)	65	82
7	I	363/387 (94%)	361 (99%)	2 (1%)	78	89
7	L	370/387 (96%)	364 (98%)	6 (2%)	55	76
8	D	362/390 (93%)	359 (99%)	3 (1%)	73	86
8	H	353/390 (90%)	351 (99%)	2 (1%)	78	89
8	J	369/390 (95%)	366 (99%)	3 (1%)	73	86
All	All	4241/7444 (57%)	4212 (99%)	29 (1%)	73	88

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1422	ILE
3	F	47	PHE
4	S	214	TYR
4	S	349	LEU
5	B	325	MET
5	G	12	ASN
5	G	205	GLU
5	G	298	VAL
6	K	24	ARG
7	L	40	VAL
7	L	89	SER
7	L	194	GLU
7	L	240	LEU
7	L	264	MET
7	L	308	ASP
8	H	22	ILE
8	H	407	THR
7	I	40	VAL
7	I	304	VAL
8	D	47	VAL
8	D	407	THR
8	D	445	GLU
7	E	40	VAL
7	E	163	THR
7	E	190	VAL
7	E	419	ILE
8	J	275	GLN
8	J	407	THR
8	J	421	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (31) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	848	HIS
1	A	883	ASN
1	A	1297	HIS
1	A	1981	ASN
1	A	1996	ASN
1	A	2200	ASN
3	F	78	GLN
3	F	134	GLN

Continued on next page...

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Mol	Chain	Res	Type
3	F	144	GLN
5	B	12	ASN
5	B	40	HIS
5	B	353	GLN
5	G	59	GLN
5	G	121	GLN
5	G	263	GLN
6	K	190	GLN
6	K	205	GLN
6	K	263	GLN
6	K	354	ASN
7	L	241	HIS
7	L	420	ASN
7	I	335	ASN
8	D	27	HIS
8	D	313	ASN
7	E	18	HIS
7	E	247	ASN
7	E	251	GLN
7	E	348	HIS
7	E	393	HIS
8	J	25	HIS
8	J	221	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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.

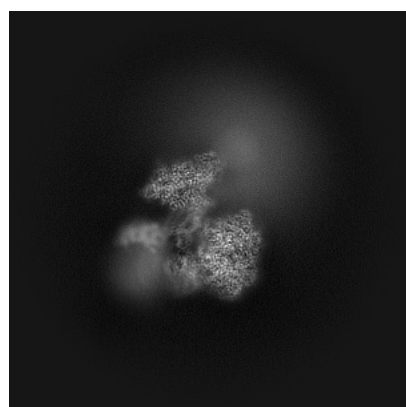
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-18611. These allow visual inspection of the internal detail of the map and identification of artifacts.

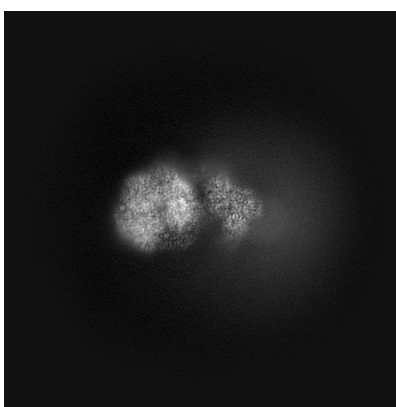
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

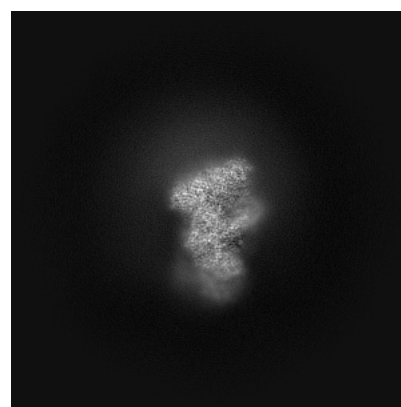
6.1.1 Primary map



X



Y

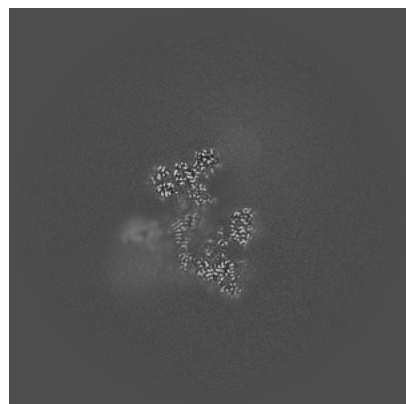


Z

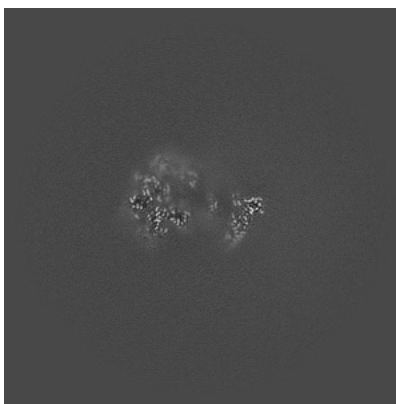
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

6.2.1 Primary map



X Index: 343



Y Index: 343

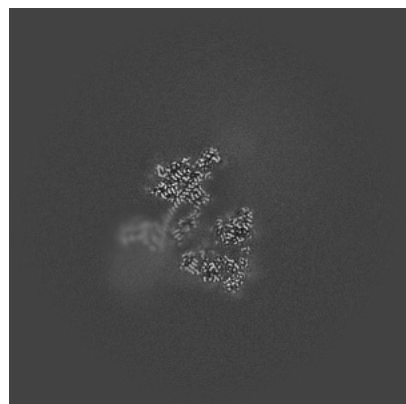


Z Index: 343

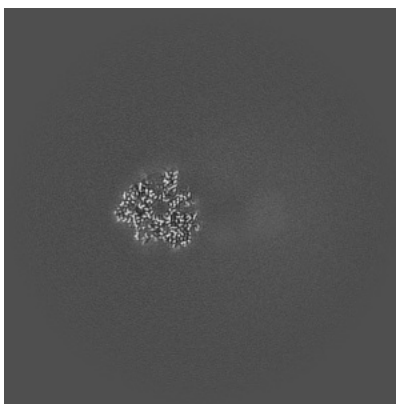
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

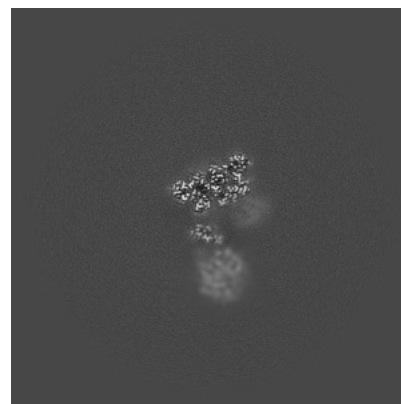
6.3.1 Primary map



X Index: 352



Y Index: 377



Z Index: 292

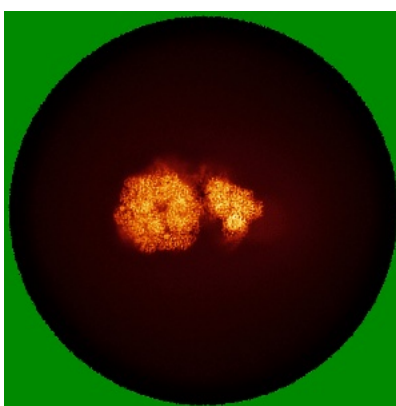
The images above show the largest variance slices of the map in three orthogonal directions.

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

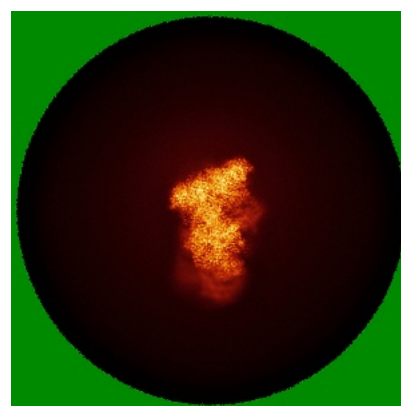
6.4.1 Primary map



X



Y

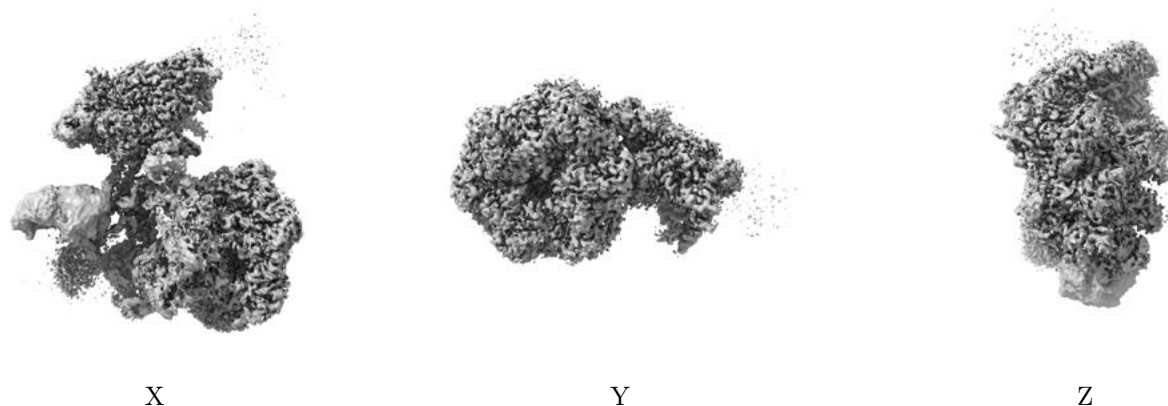


Z

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

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 6.0. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

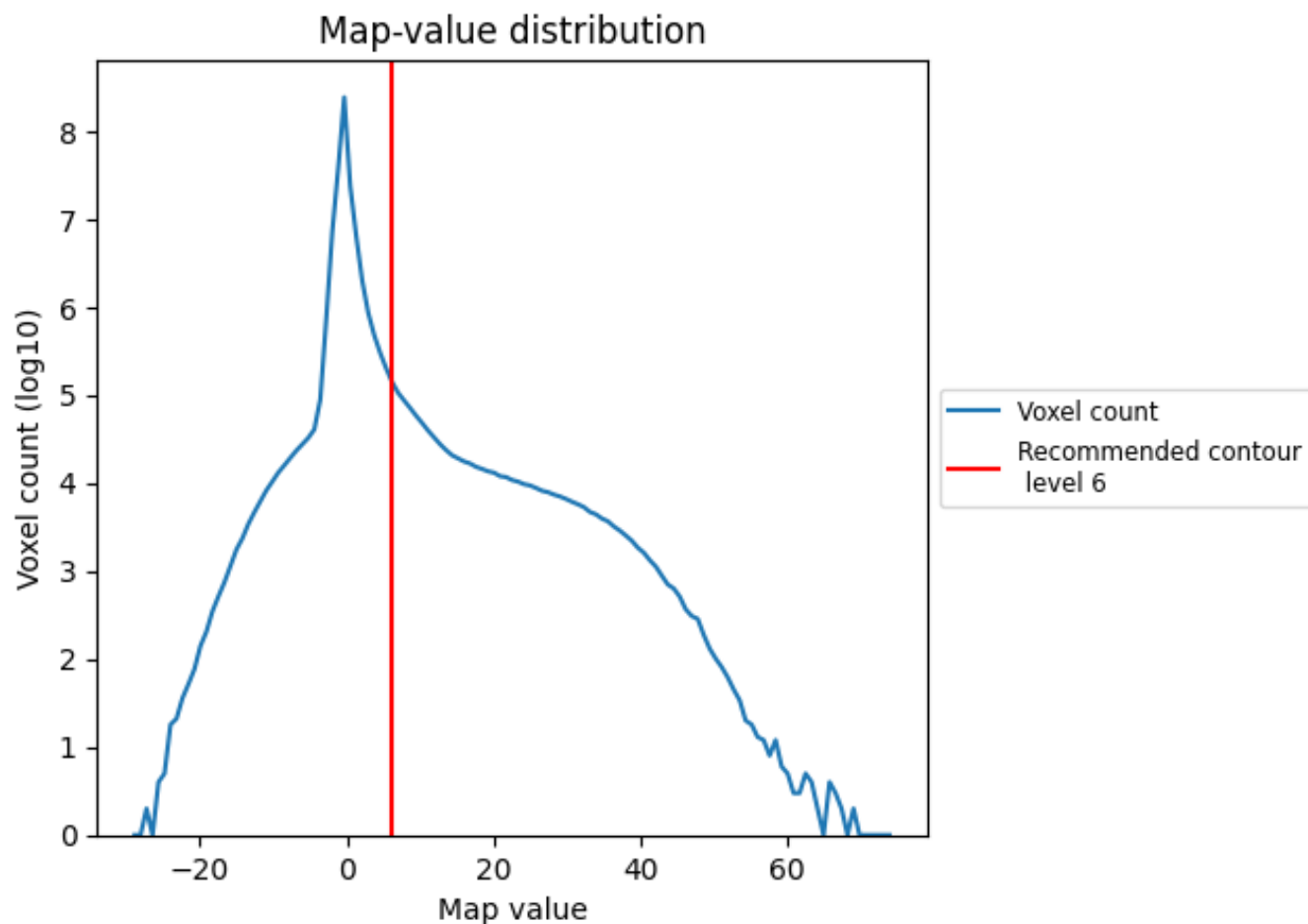
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

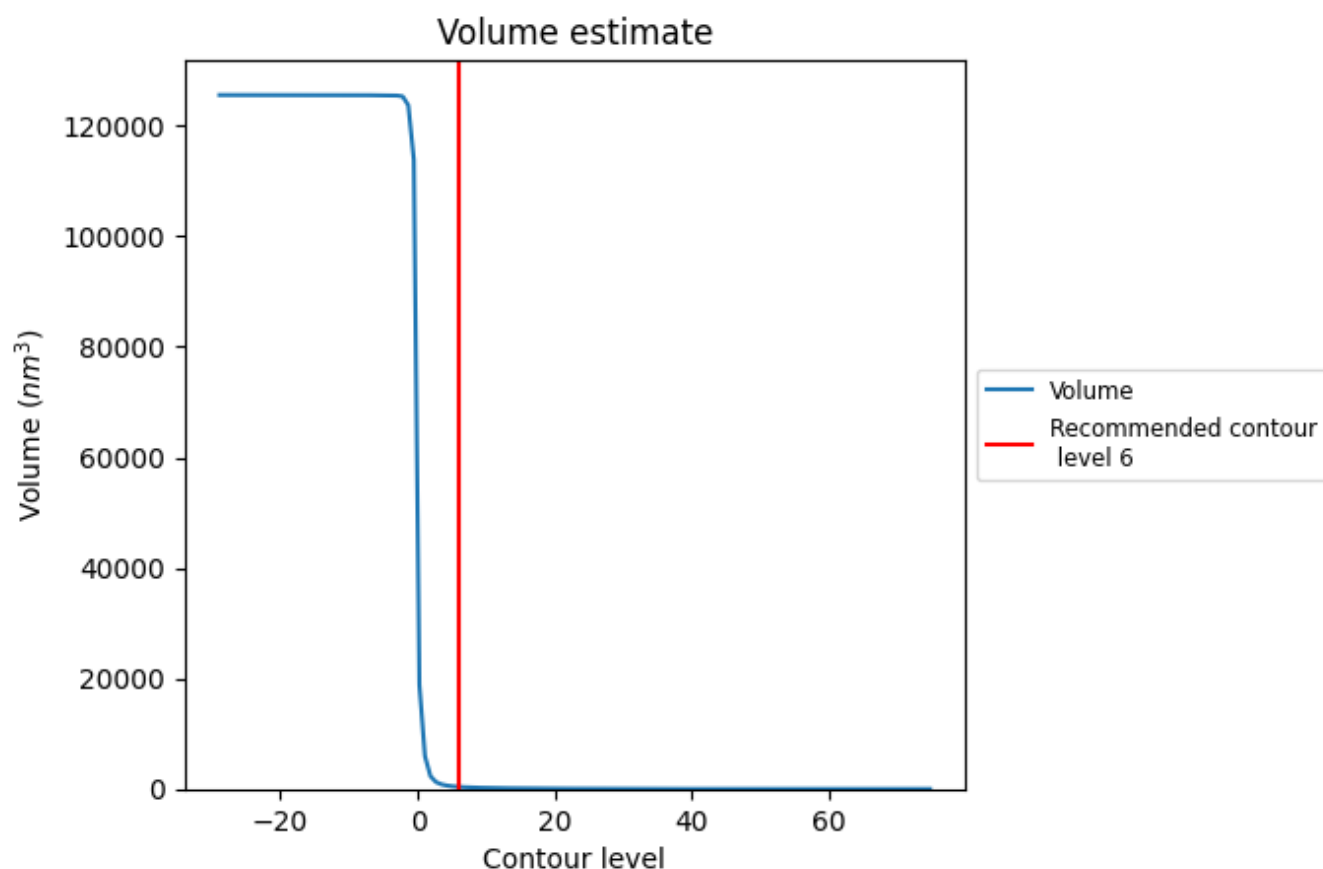
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



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

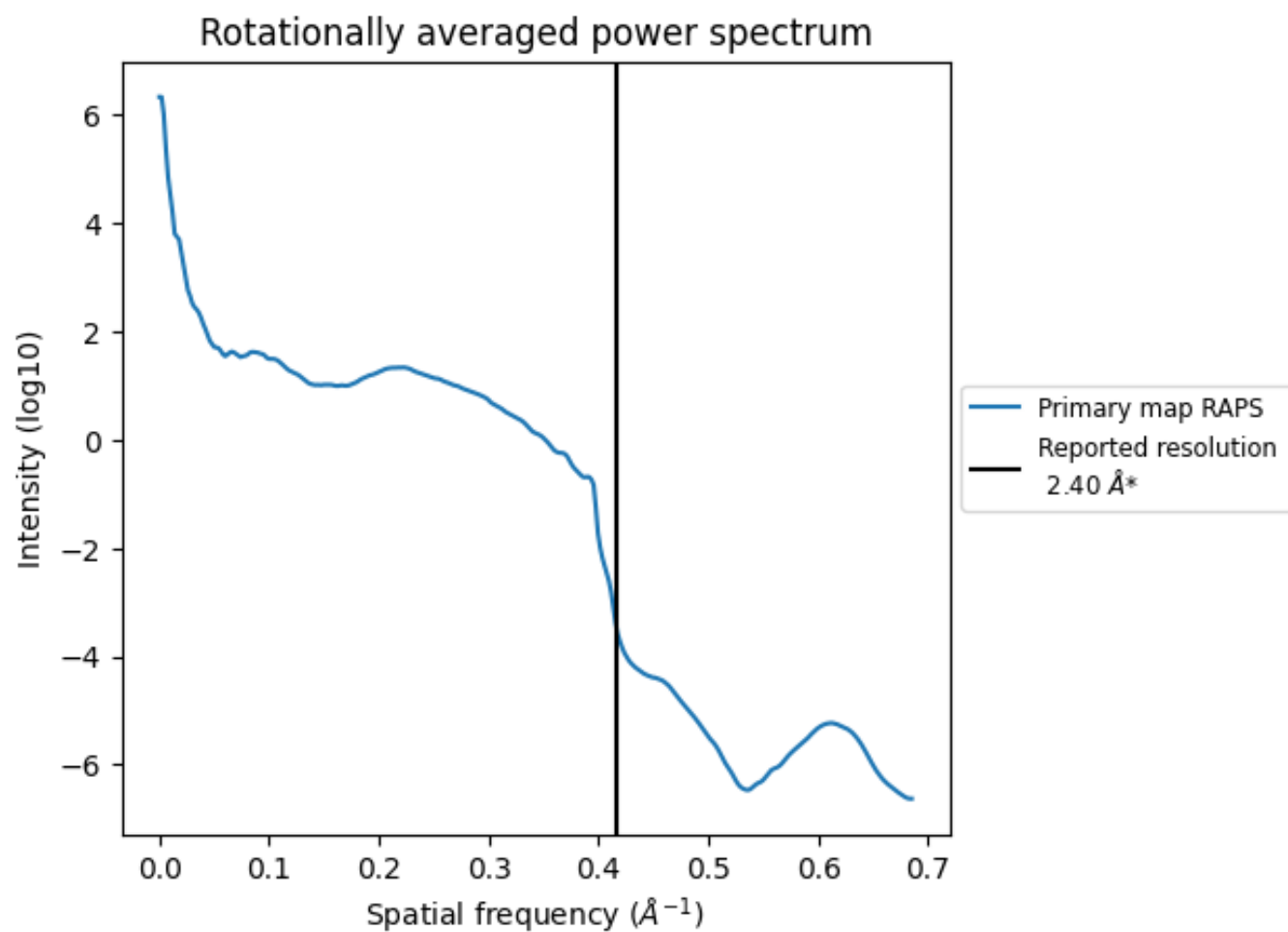
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 377 nm^3 ; this corresponds to an approximate mass of 341 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 0.417 Å⁻¹

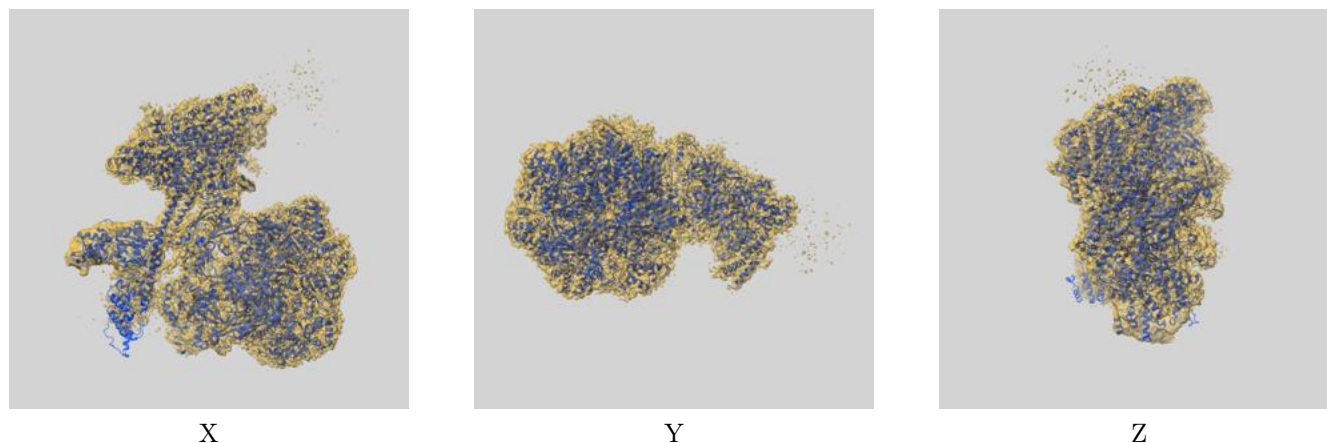
8 Fourier-Shell correlation ⓘ

This section was not generated. No FSC curve or half-maps provided.

9 Map-model fit [i](#)

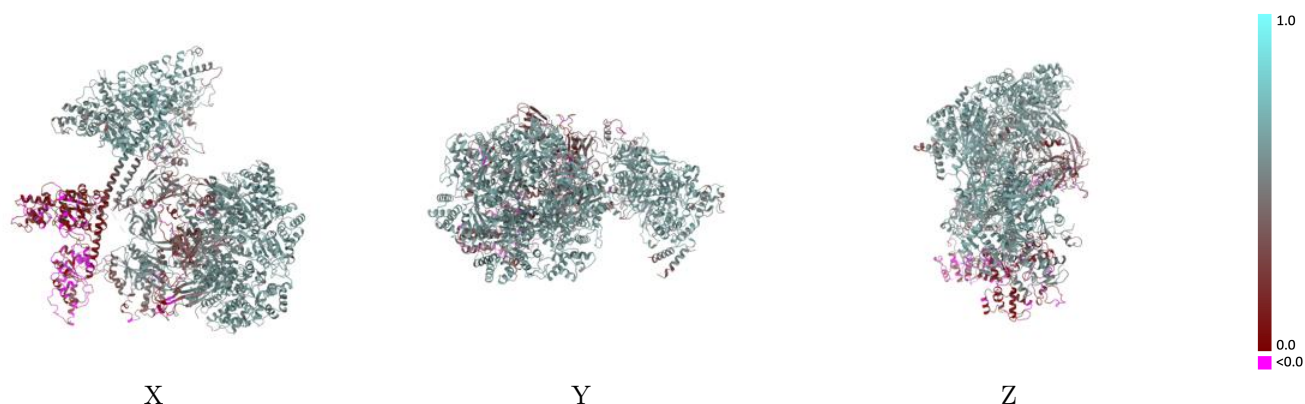
This section contains information regarding the fit between EMDB map EMD-18611 and PDB model 8QR1. Per-residue inclusion information can be found in section [3](#) on page [6](#).

9.1 Map-model overlay [i](#)



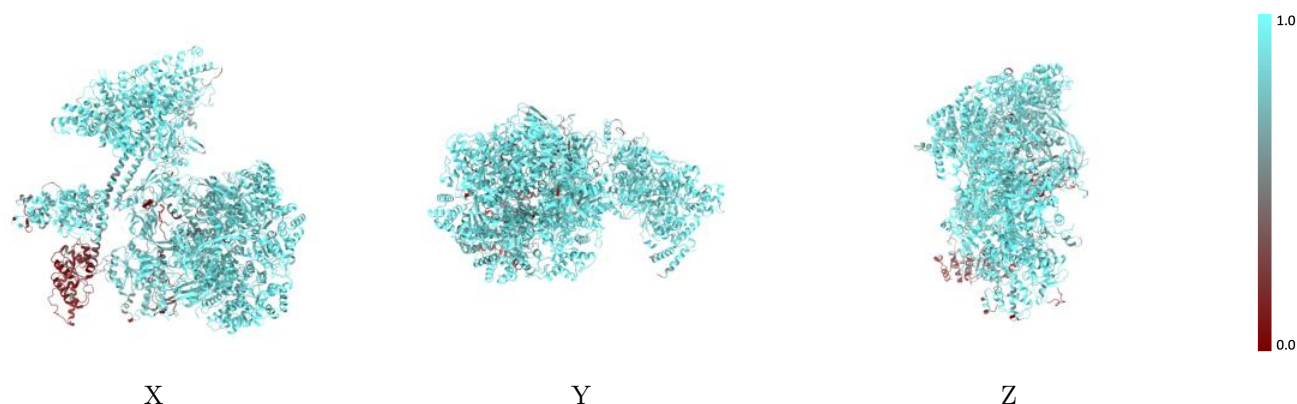
The images above show the 3D surface view of the map at the recommended contour level 6.0 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



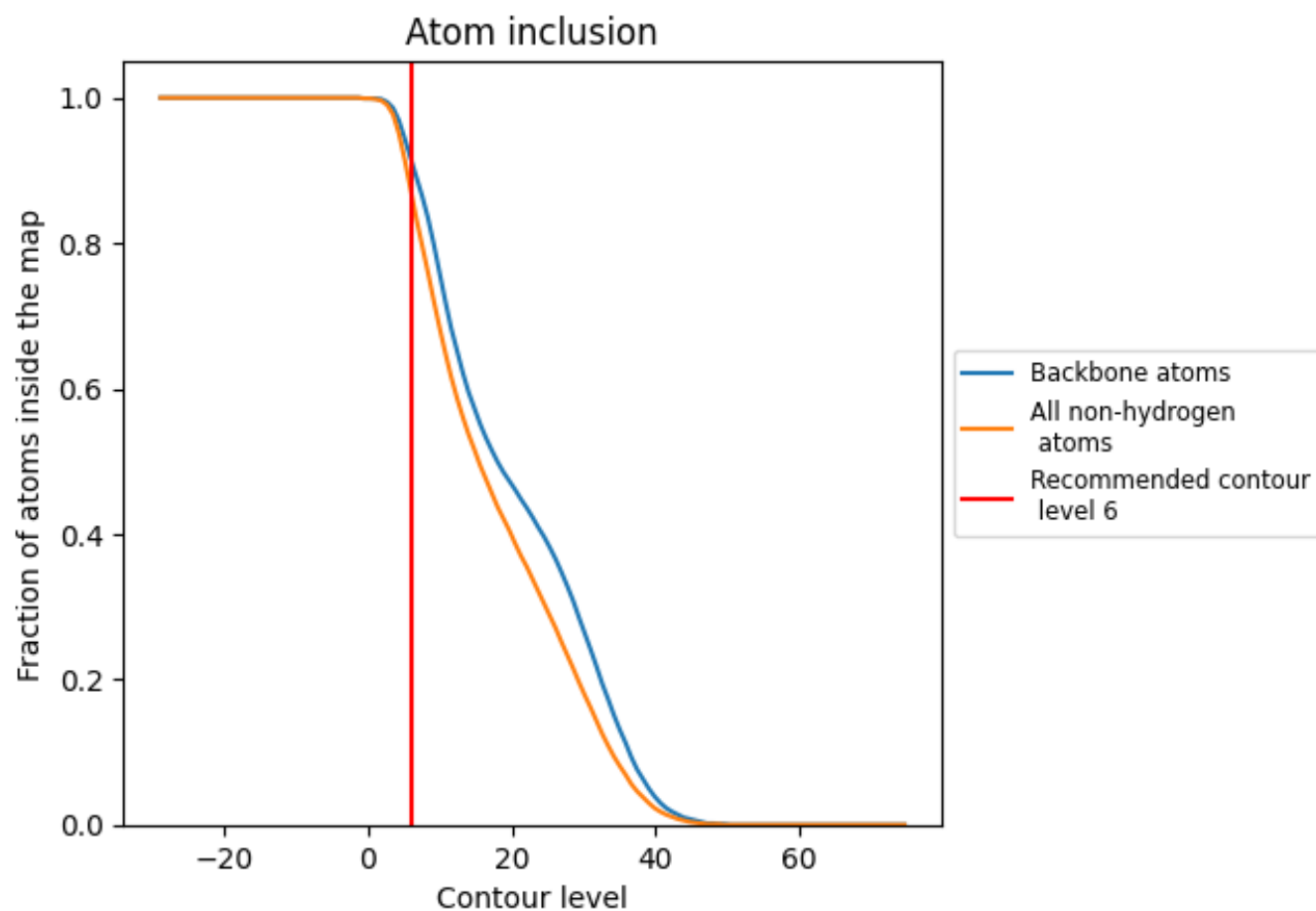
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (6).

9.4 Atom inclusion ⓘ



At the recommended contour level, 92% of all backbone atoms, 87% 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 (6) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8700	 0.4840
A	 0.6420	 0.3310
B	 0.8420	 0.1250
C	 0.9330	 0.5740
D	 0.9580	 0.5950
E	 0.9090	 0.5570
F	 0.9230	 0.5560
G	 0.9680	 0.5910
H	 0.8920	 0.5180
I	 0.9540	 0.5860
J	 0.8970	 0.5290
K	 0.9760	 0.6090
L	 0.8940	 0.5070
S	 0.8430	 0.4020

