



Full wwPDB EM Validation Report ⓘ

Mar 14, 2026 – 04:33 AM UTC

PDB ID : 9L1P / pdb_00009l1p
EMDB ID : EMD-62751
Title : Cryo-EM structure of the ICT01-BTN2A1/BTN3A1/BTN3A2 complex, local refinement
Authors : Xin, W.Z.; Huang, B.D.; Zhou, Q.
Deposited on : 2024-12-15
Resolution : 3.50 Å(reported)

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

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

A user guide is available at

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

The types of validation reports are described at

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

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

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

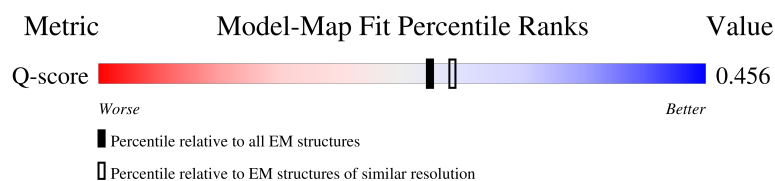
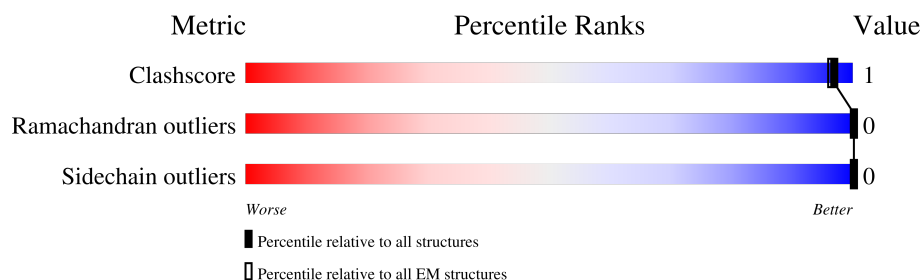
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.50 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	13950 (3.00 - 4.00)

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	B	534	 17% 5% 78%
2	E	340	 26% 8% 66%
3	H	255	 7% 64% 24% 12%
4	L	236	 69% 21% 9%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5173 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 Butyrophilin subfamily 2 member A1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	B	115	Total	C	N	O	S	0	0
			920	577	163	176	4		

There are 35 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	7	MET	-	initiating methionine	UNP Q7KYR7
B	8	ASP	-	expression tag	UNP Q7KYR7
B	9	MET	-	expression tag	UNP Q7KYR7
B	10	ARG	-	expression tag	UNP Q7KYR7
B	11	VAL	-	expression tag	UNP Q7KYR7
B	12	PRO	-	expression tag	UNP Q7KYR7
B	13	ALA	-	expression tag	UNP Q7KYR7
B	14	GLN	-	expression tag	UNP Q7KYR7
B	15	LEU	-	expression tag	UNP Q7KYR7
B	16	LEU	-	expression tag	UNP Q7KYR7
B	17	GLY	-	expression tag	UNP Q7KYR7
B	18	LEU	-	expression tag	UNP Q7KYR7
B	19	LEU	-	expression tag	UNP Q7KYR7
B	20	LEU	-	expression tag	UNP Q7KYR7
B	21	LEU	-	expression tag	UNP Q7KYR7
B	22	TRP	-	expression tag	UNP Q7KYR7
B	23	LEU	-	expression tag	UNP Q7KYR7
B	24	SER	-	expression tag	UNP Q7KYR7
B	25	GLY	-	expression tag	UNP Q7KYR7
B	26	ALA	-	expression tag	UNP Q7KYR7
B	27	ARG	-	expression tag	UNP Q7KYR7
B	28	CYS	-	expression tag	UNP Q7KYR7
B	528	GLY	-	expression tag	UNP Q7KYR7
B	529	SER	-	expression tag	UNP Q7KYR7
B	530	SER	-	expression tag	UNP Q7KYR7
B	531	GLY	-	expression tag	UNP Q7KYR7
B	532	MET	-	expression tag	UNP Q7KYR7
B	533	ASP	-	expression tag	UNP Q7KYR7

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Chain	Residue	Modelled	Actual	Comment	Reference
B	534	TYR	-	expression tag	UNP Q7KYR7
B	535	LYS	-	expression tag	UNP Q7KYR7
B	536	ASP	-	expression tag	UNP Q7KYR7
B	537	ASP	-	expression tag	UNP Q7KYR7
B	538	ASP	-	expression tag	UNP Q7KYR7
B	539	ASP	-	expression tag	UNP Q7KYR7
B	540	LYS	-	expression tag	UNP Q7KYR7

- Molecule 2 is a protein called Butyrophilin subfamily 3 member A2.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	E	116	Total	C	N	O	S	0	0
			889	561	151	172	5		

There are 35 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	8	MET	-	initiating methionine	UNP P78410
E	9	ASP	-	expression tag	UNP P78410
E	10	MET	-	expression tag	UNP P78410
E	11	ARG	-	expression tag	UNP P78410
E	12	VAL	-	expression tag	UNP P78410
E	13	PRO	-	expression tag	UNP P78410
E	14	ALA	-	expression tag	UNP P78410
E	15	GLN	-	expression tag	UNP P78410
E	16	LEU	-	expression tag	UNP P78410
E	17	LEU	-	expression tag	UNP P78410
E	18	GLY	-	expression tag	UNP P78410
E	19	LEU	-	expression tag	UNP P78410
E	20	LEU	-	expression tag	UNP P78410
E	21	LEU	-	expression tag	UNP P78410
E	22	LEU	-	expression tag	UNP P78410
E	23	TRP	-	expression tag	UNP P78410
E	24	LEU	-	expression tag	UNP P78410
E	25	SER	-	expression tag	UNP P78410
E	26	GLY	-	expression tag	UNP P78410
E	27	ALA	-	expression tag	UNP P78410
E	28	ARG	-	expression tag	UNP P78410
E	29	CYS	-	expression tag	UNP P78410
E	335	GLY	-	expression tag	UNP P78410
E	336	SER	-	expression tag	UNP P78410
E	337	SER	-	expression tag	UNP P78410

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Chain	Residue	Modelled	Actual	Comment	Reference
E	338	GLY	-	expression tag	UNP P78410
E	339	MET	-	expression tag	UNP P78410
E	340	ASP	-	expression tag	UNP P78410
E	341	TYR	-	expression tag	UNP P78410
E	342	LYS	-	expression tag	UNP P78410
E	343	ASP	-	expression tag	UNP P78410
E	344	ASP	-	expression tag	UNP P78410
E	345	ASP	-	expression tag	UNP P78410
E	346	ASP	-	expression tag	UNP P78410
E	347	LYS	-	expression tag	UNP P78410

- Molecule 3 is a protein called ICT01 heavy chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	H	225	Total	C	N	O	S	0	0
			1710	1080	285	337	8		

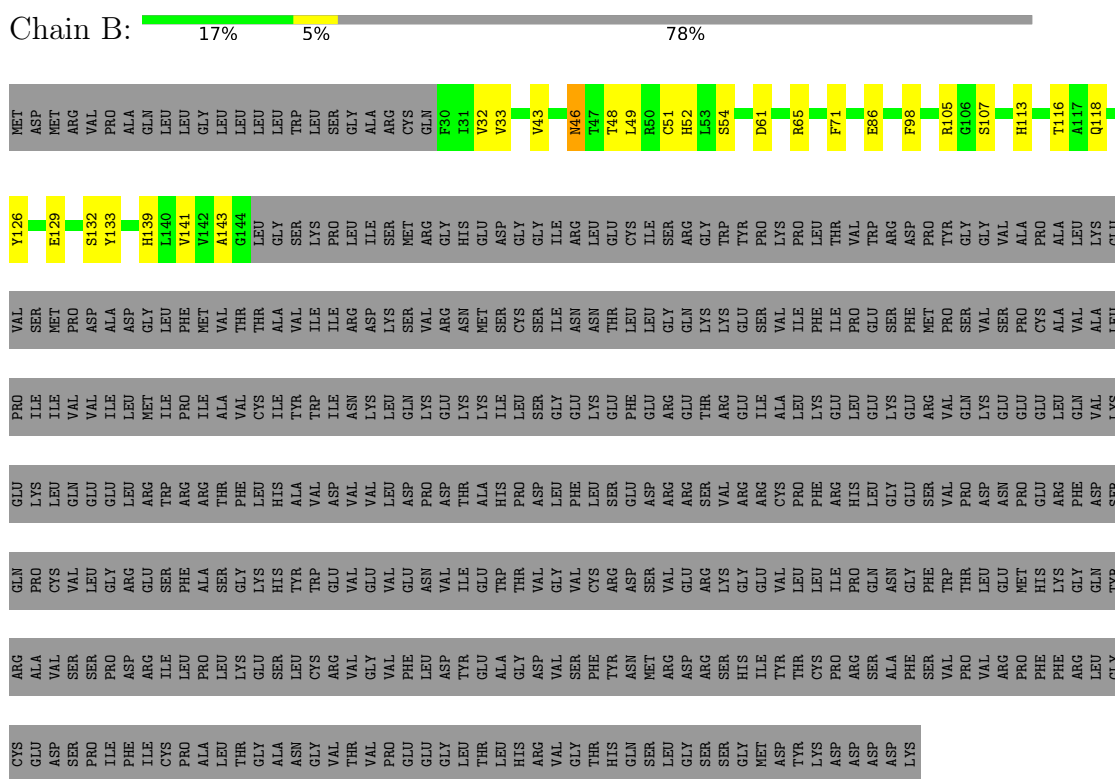
- Molecule 4 is a protein called ICT01 light chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	L	214	Total	C	N	O	S	0	0
			1654	1037	278	333	6		

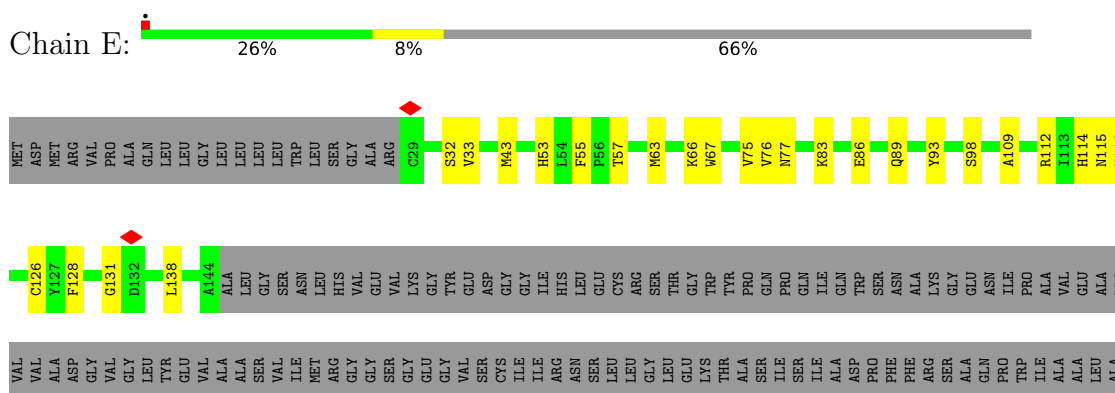
3 Residue-property plots

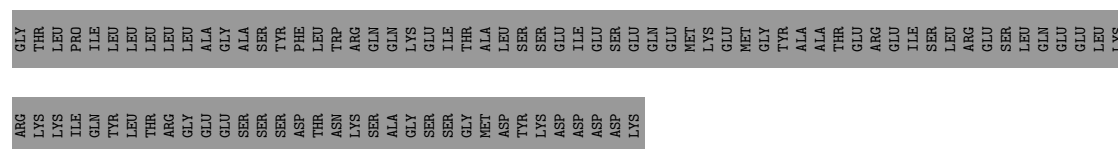
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion 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: Butyrophilin subfamily 2 member A1

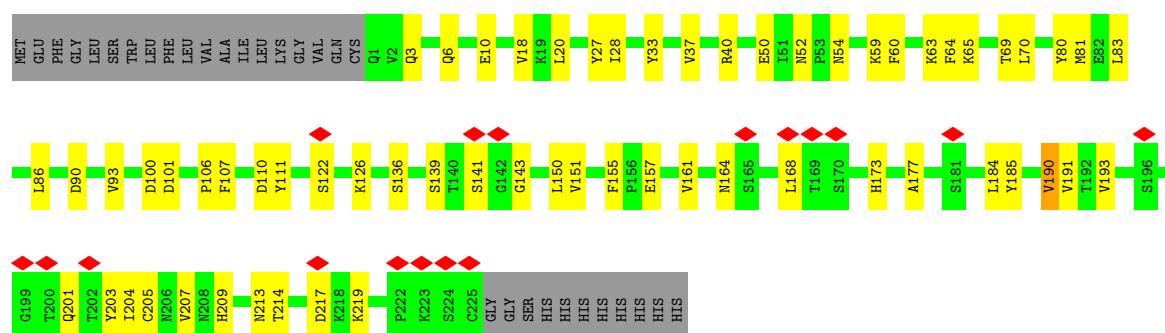


• Molecule 2: Butyrophilin subfamily 3 member A2

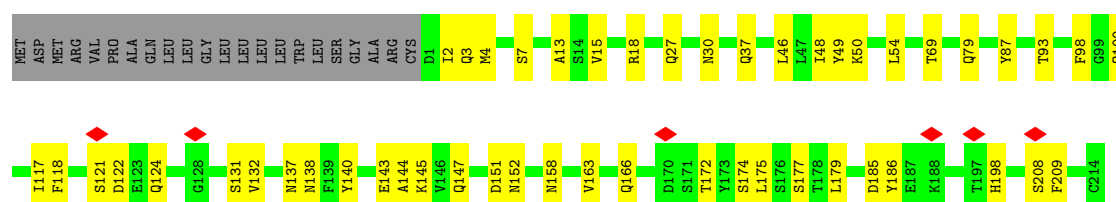




• Molecule 3: ICT01 heavy chain



• Molecule 4: ICT01 light chain



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	354724	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 ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	1200	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.273	Depositor
Minimum map value	-0.161	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.003	Depositor
Recommended contour level	0.02	Depositor
Map size ($\text{\AA}$)	391.32, 391.32, 391.32	wwPDB
Map dimensions	360, 360, 360	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	1.087, 1.087, 1.087	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	B	2.09	12/939 (1.3%)	1.74	15/1270 (1.2%)
2	E	2.23	15/906 (1.7%)	1.72	13/1226 (1.1%)
3	H	2.19	42/1751 (2.4%)	1.62	25/2380 (1.1%)
4	L	2.15	31/1692 (1.8%)	1.70	27/2299 (1.2%)
All	All	2.17	100/5288 (1.9%)	1.68	80/7175 (1.1%)

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
1	B	0	1
2	E	0	1
3	H	0	1
4	L	0	1
All	All	0	4

All (100) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	53	HIS	CD2-NE2	8.30	1.47	1.37
3	H	110	ASP	CA-C	8.19	1.60	1.52
4	L	198	HIS	ND1-CE1	8.15	1.40	1.32
4	L	175	LEU	CA-C	8.02	1.61	1.52
4	L	186	TYR	CA-C	7.84	1.61	1.52
3	H	122	SER	CA-C	-7.67	1.44	1.53
3	H	219	LYS	CA-C	7.14	1.60	1.52
4	L	140	TYR	CA-C	7.00	1.59	1.52
2	E	75	VAL	CA-C	-6.97	1.46	1.53
4	L	198	HIS	CD2-NE2	6.96	1.45	1.37
3	H	65	LYS	CA-C	6.88	1.61	1.52
4	L	147	GLN	CA-C	6.75	1.60	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	H	161	VAL	CA-C	6.73	1.59	1.53
3	H	203	TYR	CA-C	-6.62	1.44	1.52
4	L	98	PHE	CA-C	-6.56	1.44	1.52
2	E	109	ALA	CA-C	-6.53	1.44	1.52
4	L	151	ASP	CA-C	6.40	1.61	1.53
3	H	63	LYS	CA-C	6.23	1.61	1.52
3	H	69	THR	CA-C	6.23	1.59	1.52
4	L	208	SER	C-O	-6.04	1.16	1.23
4	L	132	VAL	CA-C	-6.02	1.45	1.52
4	L	177	SER	CA-C	-6.02	1.45	1.52
4	L	69	THR	CA-C	6.02	1.61	1.52
1	B	116	THR	CA-C	-5.93	1.45	1.52
1	B	51	CYS	CA-C	-5.92	1.45	1.52
2	E	53	HIS	CE1-NE2	-5.87	1.26	1.32
3	H	209	HIS	CD2-NE2	5.85	1.44	1.37
4	L	13	ALA	C-O	-5.82	1.16	1.23
2	E	43	MET	SD-CE	5.78	1.94	1.79
3	H	54	ASN	CA-C	5.77	1.60	1.52
3	H	111	TYR	CA-C	-5.68	1.45	1.52
3	H	204	ILE	C-O	-5.64	1.18	1.24
2	E	112	ARG	C-O	-5.62	1.17	1.24
1	B	133	TYR	CA-C	-5.60	1.46	1.52
3	H	60	PHE	C-O	-5.59	1.16	1.23
4	L	48	ILE	CA-C	-5.59	1.46	1.52
1	B	118	GLN	CA-C	5.58	1.59	1.52
4	L	46	LEU	CA-C	5.57	1.60	1.52
4	L	131	SER	CA-C	5.57	1.60	1.53
1	B	86	GLU	CA-C	-5.56	1.45	1.52
3	H	190	VAL	C-O	-5.54	1.18	1.24
4	L	2	ILE	CA-C	5.54	1.59	1.52
3	H	126	LYS	CA-C	5.54	1.59	1.52
1	B	43	VAL	CA-C	-5.54	1.44	1.53
2	E	126	CYS	CA-C	-5.52	1.45	1.52
3	H	28	ILE	CA-C	5.50	1.59	1.52
1	B	139	HIS	CD2-NE2	5.50	1.44	1.37
2	E	89	GLN	C-O	-5.50	1.17	1.23
3	H	207	VAL	C-O	-5.50	1.18	1.24
1	B	143	ALA	C-O	-5.48	1.17	1.23
1	B	49	LEU	CA-C	5.47	1.60	1.52
3	H	100	ASP	C-O	-5.45	1.18	1.23
3	H	150	LEU	CA-C	5.43	1.59	1.52
1	B	126	TYR	C-O	-5.43	1.17	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	L	79	GLN	CA-C	-5.43	1.46	1.52
4	L	209	PHE	CA-C	-5.41	1.46	1.52
3	H	80	TYR	CA-C	5.36	1.59	1.52
3	H	201	GLN	CA-C	5.36	1.59	1.52
3	H	107	PHE	CA-C	5.36	1.59	1.52
3	H	93	VAL	N-CA	5.35	1.52	1.46
3	H	185	TYR	C-O	-5.35	1.17	1.23
4	L	172	THR	N-CA	5.35	1.52	1.45
4	L	144	ALA	C-O	-5.35	1.17	1.23
4	L	122	ASP	N-CA	-5.32	1.39	1.46
2	E	33	VAL	CA-C	-5.29	1.46	1.52
3	H	205	CYS	CA-C	5.28	1.58	1.52
4	L	4	MET	CA-C	-5.28	1.46	1.52
4	L	93	THR	CA-C	-5.25	1.46	1.52
4	L	145	LYS	CA-C	5.25	1.58	1.52
4	L	50	LYS	CA-C	5.23	1.58	1.52
3	H	151	VAL	C-N	5.22	1.40	1.33
2	E	55	PHE	CA-C	-5.20	1.46	1.52
3	H	60	PHE	CA-C	-5.19	1.46	1.52
3	H	100	ASP	CA-C	-5.19	1.46	1.53
3	H	101	ASP	CA-C	-5.18	1.44	1.52
3	H	168	LEU	C-O	5.18	1.29	1.23
3	H	177	ALA	CA-C	-5.17	1.46	1.52
2	E	63	MET	C-O	-5.17	1.17	1.24
1	B	107	SER	CA-C	-5.16	1.46	1.52
3	H	136	SER	CA-C	-5.16	1.46	1.53
4	L	140	TYR	C-O	5.15	1.29	1.24
2	E	57	THR	N-CA	-5.15	1.39	1.46
3	H	27	TYR	C-O	5.14	1.29	1.23
3	H	20	LEU	N-CA	-5.14	1.39	1.46
3	H	141	SER	C-O	-5.14	1.18	1.24
1	B	98	PHE	CA-C	-5.10	1.45	1.53
3	H	204	ILE	C-N	5.09	1.40	1.33
2	E	138	LEU	CA-C	-5.08	1.46	1.52
2	E	83	LYS	CA-C	-5.07	1.46	1.52
4	L	121	SER	C-O	-5.06	1.18	1.23
3	H	143	GLY	C-N	5.05	1.39	1.33
2	E	114	HIS	N-CA	-5.04	1.40	1.46
3	H	37	VAL	C-O	5.04	1.29	1.23
3	H	184	LEU	CA-C	5.03	1.59	1.52
3	H	193	VAL	CA-C	-5.02	1.48	1.52
3	H	139	SER	C-O	-5.01	1.18	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	L	118	PHE	C-O	-5.01	1.17	1.24
4	L	143	GLU	N-CA	-5.01	1.40	1.46
3	H	86	LEU	N-CA	5.00	1.52	1.46
4	L	185	ASP	N-CA	5.00	1.51	1.46

All (80) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	H	209	HIS	ND1-CE1-NE2	8.43	116.83	108.40
4	L	124	GLN	OE1-CD-NE2	-8.02	114.58	122.60
1	B	32	VAL	N-CA-C	7.82	119.54	108.36
3	H	191	VAL	N-CA-C	7.62	119.45	108.48
4	L	186	TYR	N-CA-C	-7.39	104.11	113.43
4	L	117	ILE	N-CA-C	7.27	117.17	106.85
4	L	100	GLN	OE1-CD-NE2	-7.22	115.38	122.60
4	L	140	TYR	CA-C-O	-7.16	113.75	120.34
4	L	163	VAL	N-CA-C	7.00	117.91	108.11
3	H	209	HIS	CE1-NE2-CD2	-6.89	102.11	109.00
3	H	3	GLN	OE1-CD-NE2	-6.88	115.72	122.60
1	B	129	GLU	N-CA-C	6.79	119.59	108.52
1	B	61	ASP	CA-CB-CG	6.75	119.35	112.60
3	H	164	ASN	CA-CB-CG	6.68	119.28	112.60
2	E	128	PHE	CA-CB-CG	-6.51	107.29	113.80
2	E	86	GLU	N-CA-C	-6.37	105.66	113.50
3	H	217	ASP	CA-CB-CG	6.35	118.95	112.60
4	L	18	ARG	N-CA-C	6.31	119.79	109.76
2	E	116	VAL	CA-C-O	-6.27	115.82	121.97
2	E	115	ASN	N-CA-C	-6.26	102.83	111.28
2	E	76	VAL	CA-C-O	-6.20	114.28	120.85
1	B	46	ASN	CA-CB-CG	-6.19	106.41	112.60
4	L	166	GLN	OE1-CD-NE2	-6.17	116.43	122.60
1	B	133	TYR	N-CA-C	6.06	118.83	109.07
4	L	158	ASN	OD1-CG-ND2	-6.04	116.56	122.60
2	E	114	HIS	CB-CG-CD2	-6.03	123.37	131.20
3	H	106	PRO	CA-C-O	-5.94	114.69	121.11
1	B	71	PHE	CA-C-O	-5.94	114.58	120.70
3	H	33	TYR	N-CA-C	-5.93	102.70	110.53
4	L	3	GLN	N-CA-C	-5.93	101.75	110.52
4	L	152	ASN	CA-CB-CG	-5.89	106.71	112.60
2	E	67	TRP	N-CA-C	5.88	118.25	109.25
4	L	117	ILE	N-CA-CB	-5.86	104.62	111.83
3	H	40	ARG	O-C-N	5.85	128.05	121.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	46	ASN	CA-C-O	-5.83	114.32	121.78
1	B	98	PHE	CA-C-O	-5.82	114.85	121.72
4	L	138	ASN	OD1-CG-ND2	-5.82	116.78	122.60
3	H	193	VAL	N-CA-CB	-5.78	106.76	111.67
3	H	52	ASN	OD1-CG-ND2	5.73	128.33	122.60
2	E	93	TYR	N-CA-C	-5.72	105.04	113.61
3	H	6	GLN	OE1-CD-NE2	-5.71	116.89	122.60
3	H	101	ASP	CA-CB-CG	5.71	118.31	112.60
3	H	18	VAL	CA-C-O	-5.70	116.17	120.96
4	L	15	VAL	N-CA-C	5.68	116.93	109.21
4	L	30	ASN	CA-CB-CG	5.67	118.27	112.60
2	E	98	SER	CA-C-O	-5.66	113.58	120.54
3	H	90	ASP	CA-CB-CG	5.66	118.25	112.60
1	B	141	VAL	CA-C-O	-5.62	114.49	120.39
4	L	54	LEU	N-CA-C	5.61	118.48	110.24
1	B	132	SER	CA-C-O	5.55	126.69	120.92
1	B	105	ARG	N-CA-C	-5.51	107.20	114.31
4	L	37	GLN	OE1-CD-NE2	-5.48	117.12	122.60
4	L	179	LEU	N-CA-C	5.47	118.53	109.46
3	H	81	MET	CA-C-O	5.44	126.12	120.30
3	H	157	GLU	CA-C-O	-5.39	115.08	120.64
4	L	27	GLN	OE1-CD-NE2	-5.37	117.23	122.60
2	E	77	ASN	O-C-N	-5.36	118.15	123.46
3	H	83	LEU	N-CA-C	-5.34	101.87	110.14
3	H	64	PHE	CA-CB-CG	5.33	119.13	113.80
3	H	10	GLU	N-CA-C	5.31	117.48	109.41
4	L	122	ASP	N-CA-C	-5.29	104.88	112.13
4	L	185	ASP	N-CA-C	5.29	118.54	112.57
3	H	70	LEU	N-CA-C	5.24	118.09	109.76
3	H	193	VAL	N-CA-C	5.23	114.73	108.45
3	H	52	ASN	N-CA-C	5.22	116.50	109.24
3	H	213	ASN	OD1-CG-ND2	5.17	127.78	122.60
4	L	49	TYR	CA-C-O	5.16	126.60	120.66
2	E	32	SER	CA-C-N	-5.16	116.44	123.10
2	E	32	SER	C-N-CA	-5.16	116.44	123.10
1	B	48	THR	CA-C-O	5.12	125.66	120.24
4	L	2	ILE	CA-C-O	-5.11	115.03	121.11
1	B	129	GLU	CB-CG-CD	5.10	121.27	112.60
3	H	214	THR	N-CA-C	-5.09	102.54	109.71
1	B	141	VAL	N-CA-C	5.05	115.18	108.11
2	E	66	LYS	N-CA-C	5.03	116.36	107.61
4	L	87	TYR	CA-C-N	-5.01	115.02	122.74

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	L	87	TYR	C-N-CA	-5.01	115.02	122.74
1	B	54	SER	N-CA-C	5.01	119.56	112.75
4	L	118	PHE	CA-C-N	-5.00	115.23	120.38
4	L	118	PHE	C-N-CA	-5.00	115.23	120.38

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	65	ARG	Peptide
2	E	131	GLY	Peptide
3	H	155	PHE	Peptide
4	L	7	SER	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	B	920	0	887	3	0
2	E	889	0	872	0	0
3	H	1710	0	1680	4	0
4	L	1654	0	1602	2	0
All	All	5173	0	5041	7	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 (7) 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:B:33:VAL:HG23	1:B:52:HIS:CE1	2.21	0.75
1:B:33:VAL:CG2	1:B:52:HIS:CE1	2.70	0.75
3:H:173:HIS:NE2	4:L:137:ASN:OD1	2.24	0.69
3:H:173:HIS:HB2	3:H:190:VAL:HG22	1.77	0.66
3:H:173:HIS:CE1	4:L:174:SER:HB2	2.36	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:50:GLU:HG2	3:H:59:LYS:HB2	1.83	0.61
1:B:46:ASN:OD1	1:B:113:HIS:HA	2.20	0.42

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	B	113/534 (21%)	111 (98%)	2 (2%)	0	100	100
2	E	114/340 (34%)	114 (100%)	0	0	100	100
3	H	223/255 (88%)	222 (100%)	1 (0%)	0	100	100
4	L	212/236 (90%)	212 (100%)	0	0	100	100
All	All	662/1365 (48%)	659 (100%)	3 (0%)	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	B	99/472 (21%)	99 (100%)	0	100	100
2	E	95/278 (34%)	95 (100%)	0	100	100
3	H	193/218 (88%)	193 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
4	L	189/207 (91%)	189 (100%)	0	100	100
All	All	576/1175 (49%)	576 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	52	HIS
1	B	139	HIS
3	H	201	GLN
4	L	27	GLN
4	L	92	GLN
4	L	100	GLN
4	L	155	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 ⓘ

There are no chain breaks in this entry.

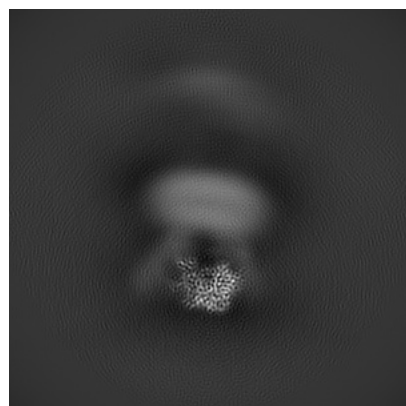
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-62751. These allow visual inspection of the internal detail of the map and identification of artifacts.

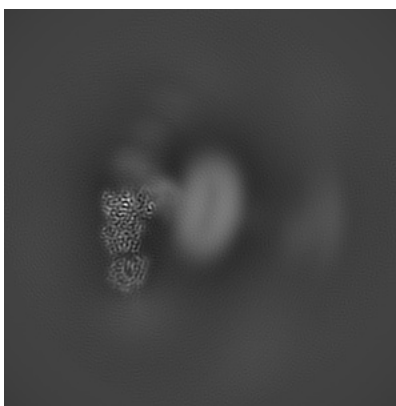
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

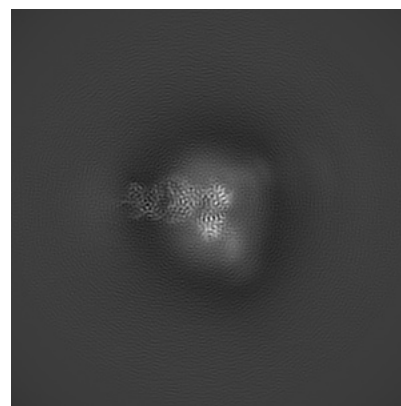
6.1.1 Primary map



X

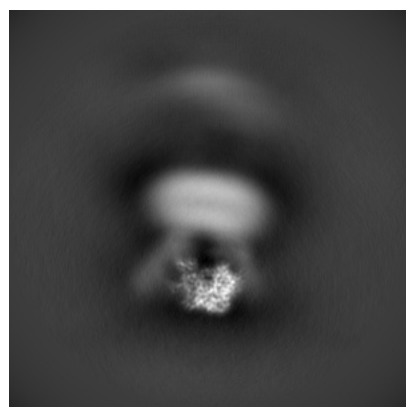


Y

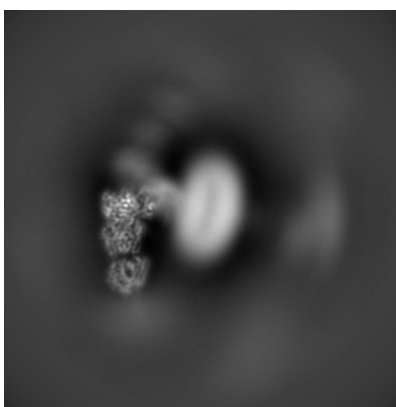


Z

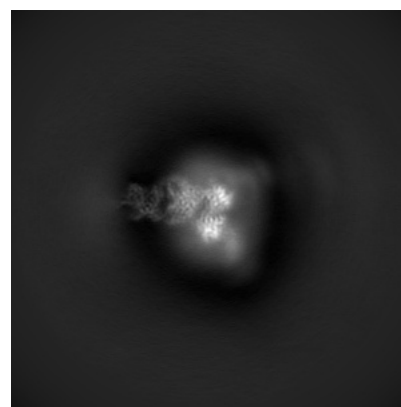
6.1.2 Raw map



X



Y

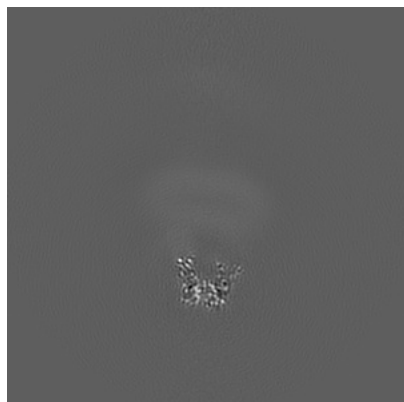


Z

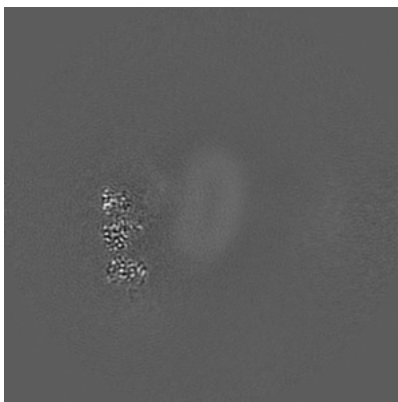
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

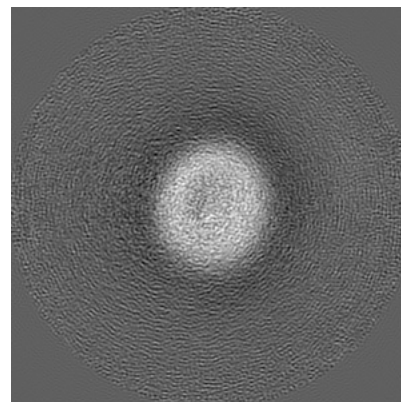
6.2.1 Primary map



X Index: 180

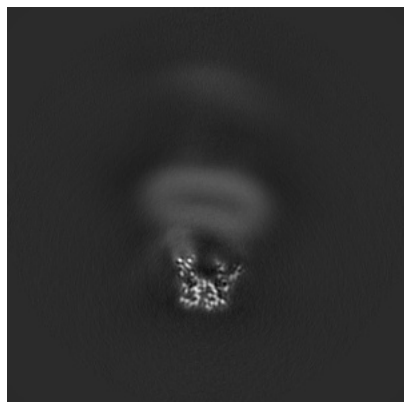


Y Index: 180

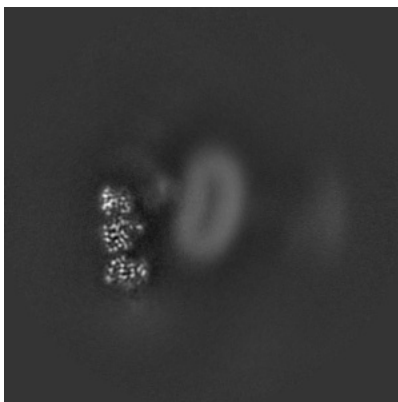


Z Index: 180

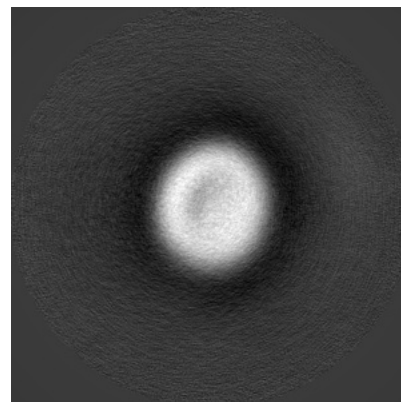
6.2.2 Raw map



X Index: 180



Y Index: 180

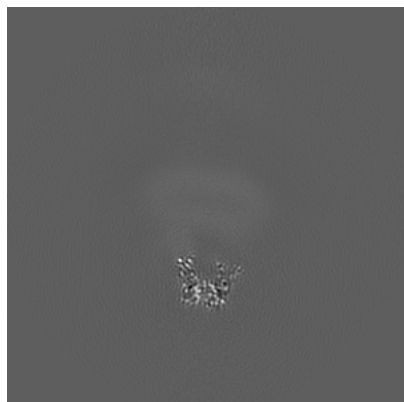


Z Index: 180

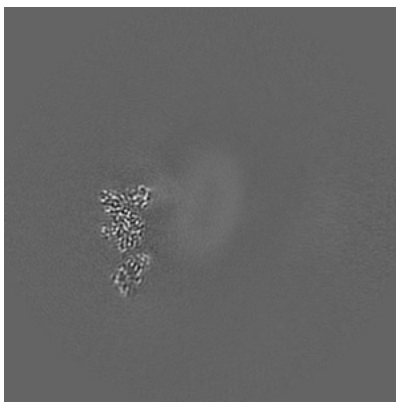
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

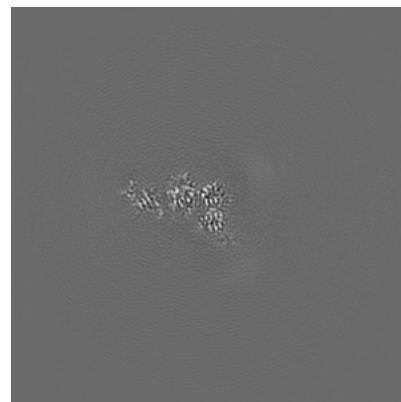
6.3.1 Primary map



X Index: 180

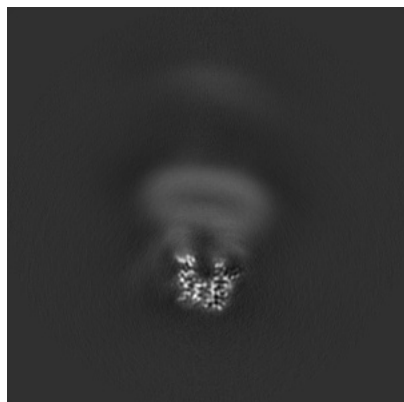


Y Index: 192

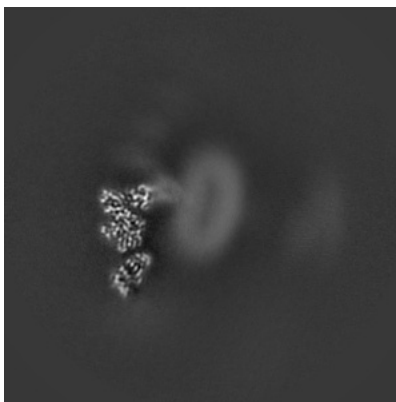


Z Index: 108

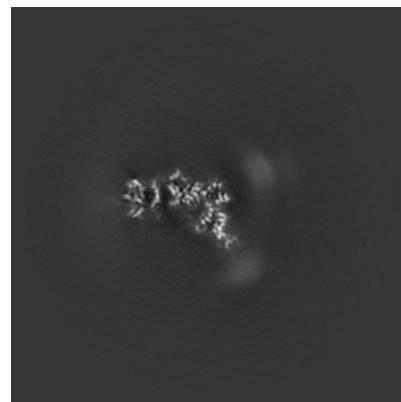
6.3.2 Raw map



X Index: 184



Y Index: 192

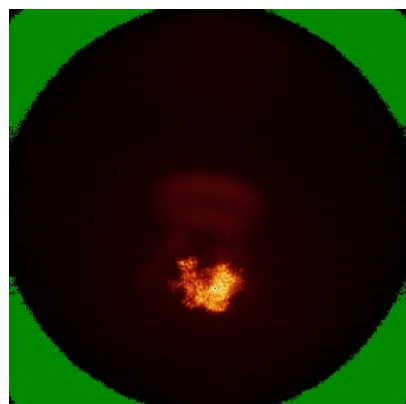


Z Index: 112

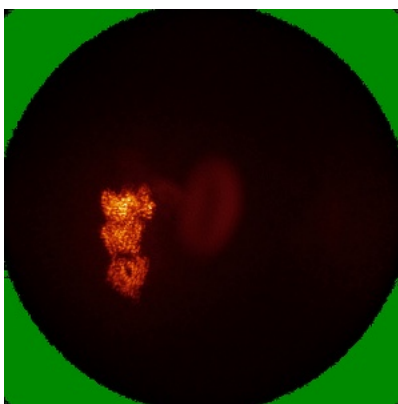
The images above show the largest variance slices of the map in three orthogonal directions.

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

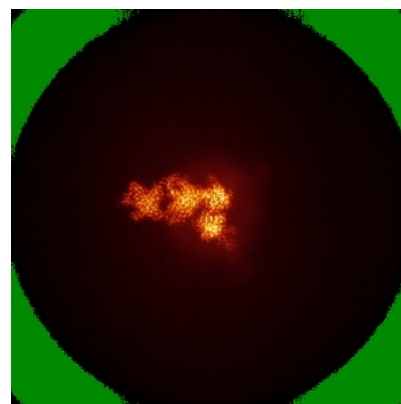
6.4.1 Primary map



X

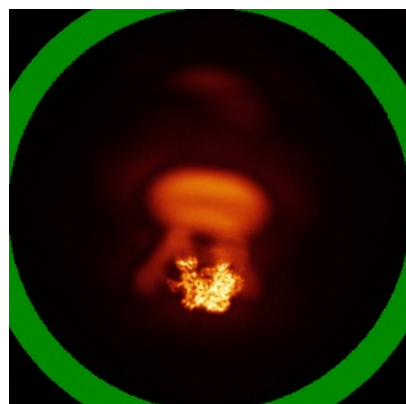


Y

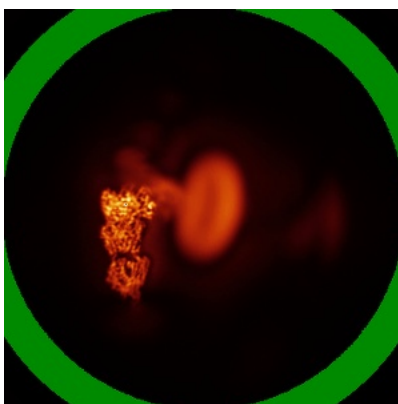


Z

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 0.02. 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.

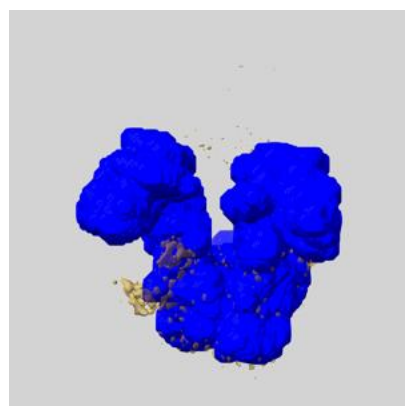
6.6 Mask visualisation [i](#)

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

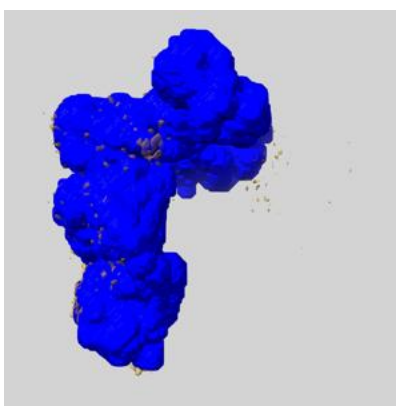
A mask typically either:

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

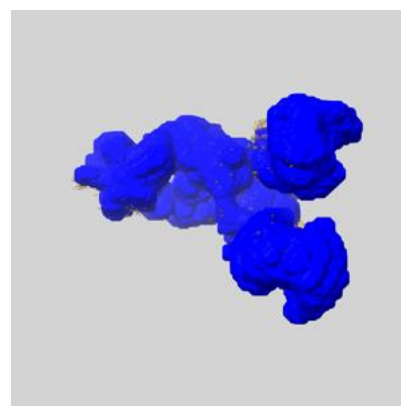
6.6.1 emd_62751_msk_1.map [i](#)



X



Y

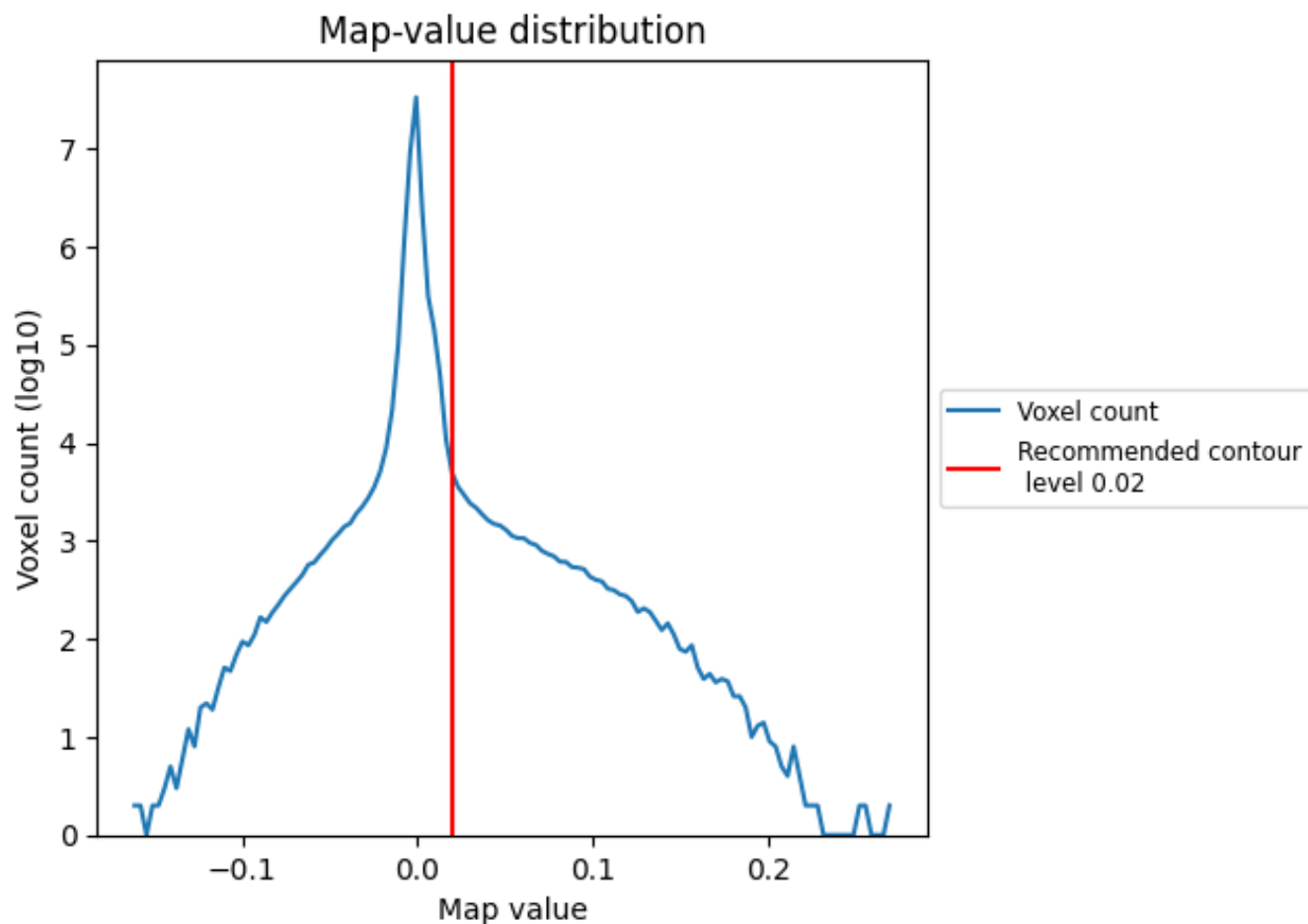


Z

7 Map analysis [i](#)

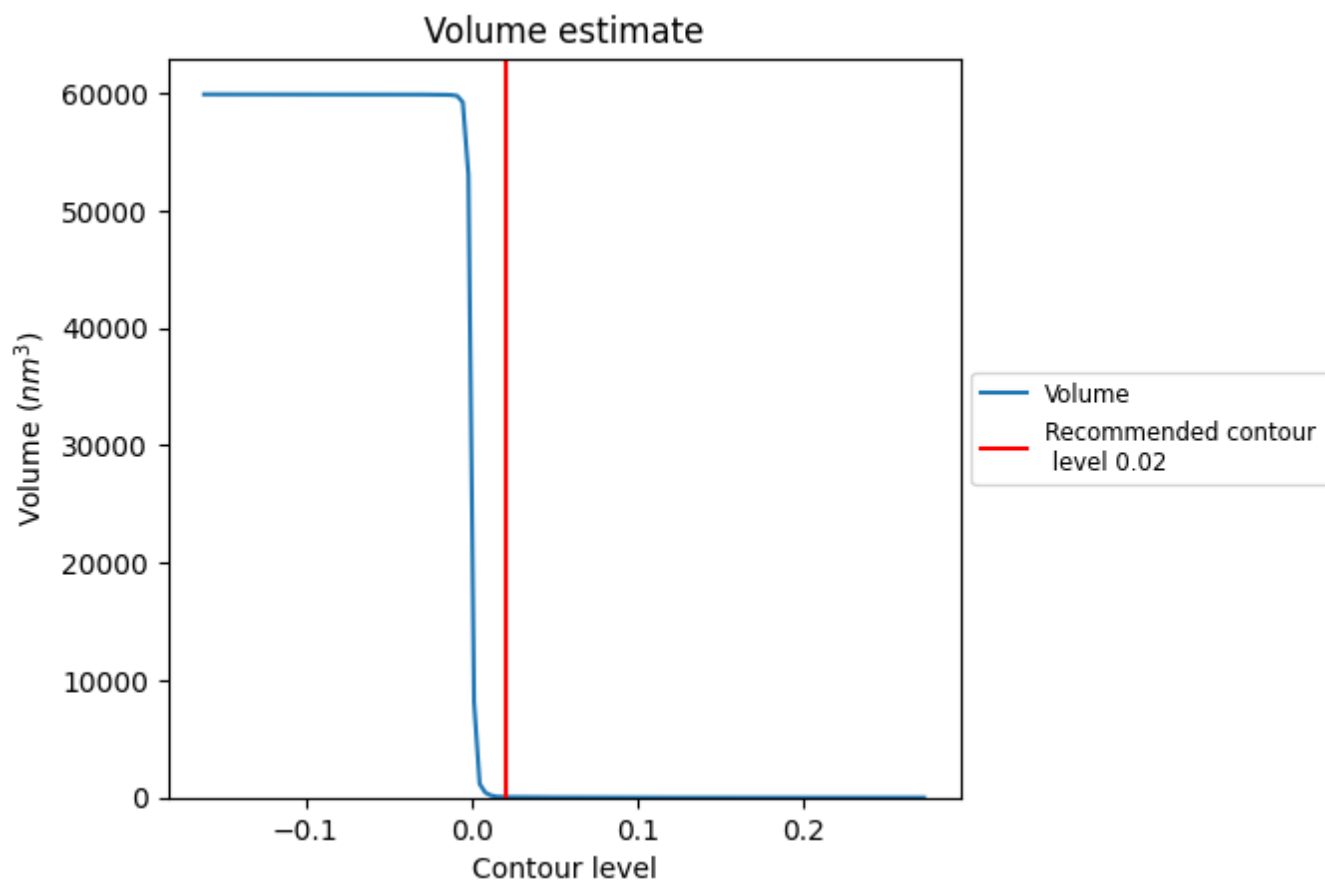
This section contains the results of statistical analysis of the map.

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.

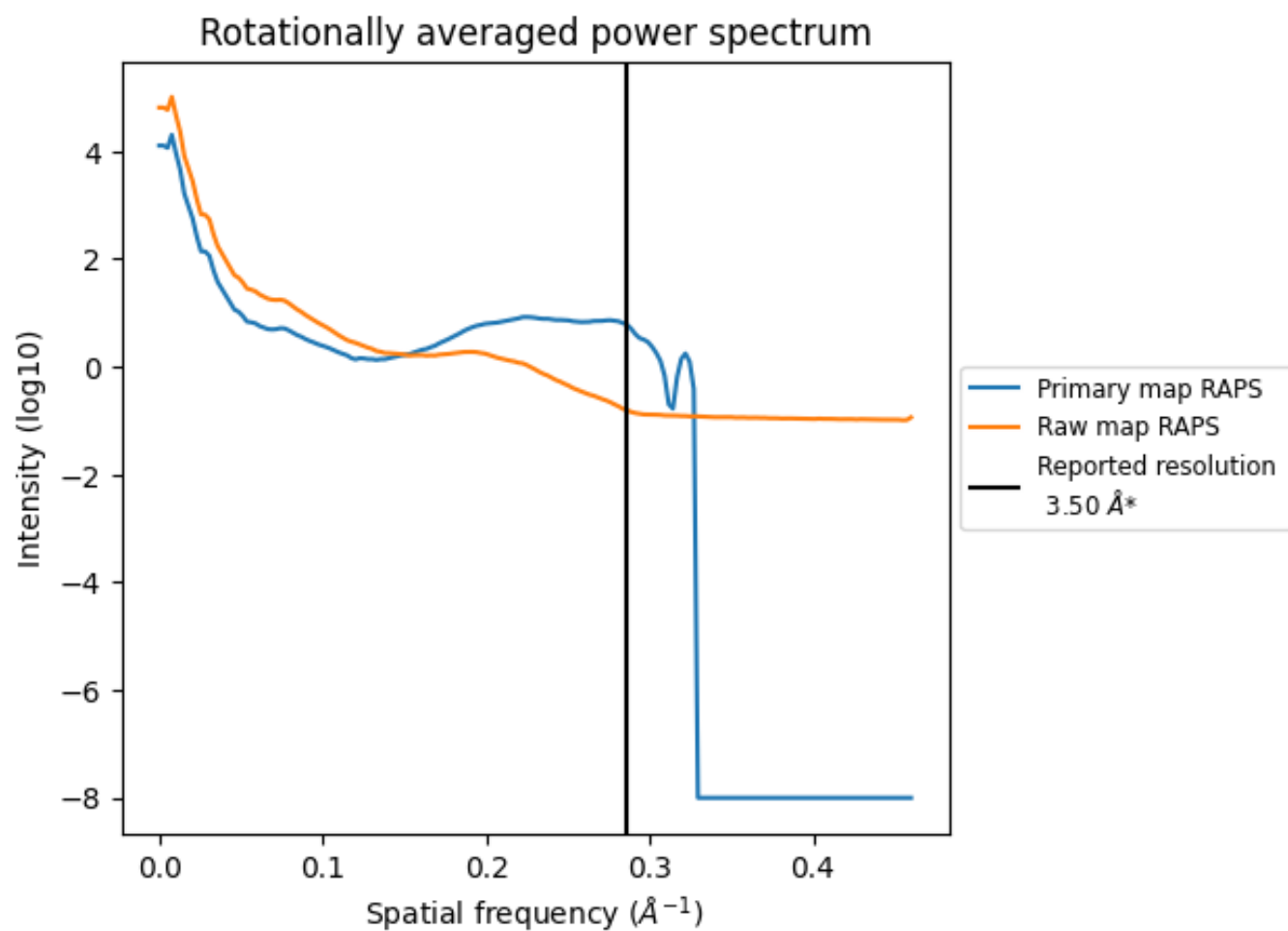
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 49 nm^3 ; this corresponds to an approximate mass of 44 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 ⓘ

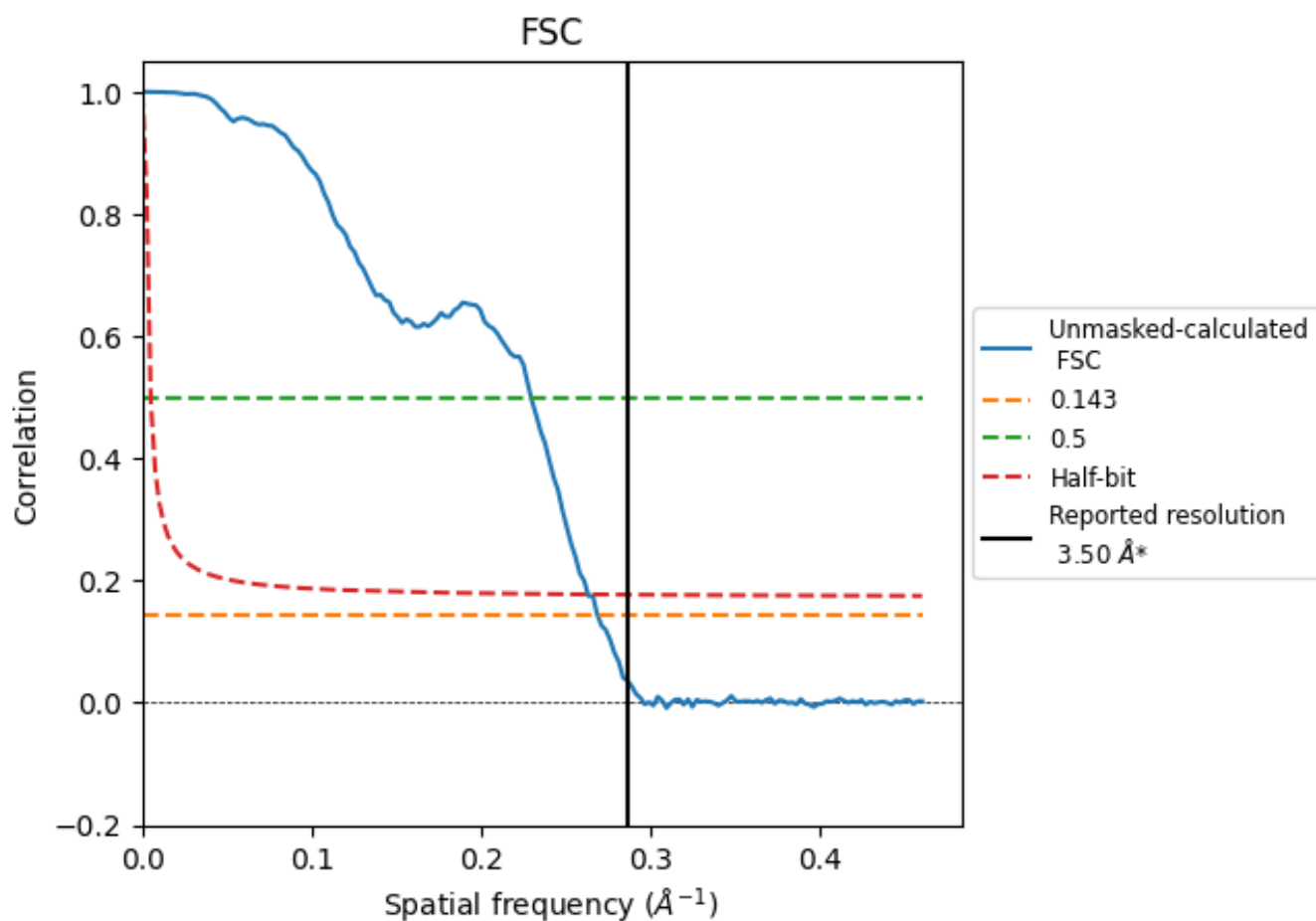


*Reported resolution corresponds to spatial frequency of $0.286 \text{ \AA}^{-1}$

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 0.286 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

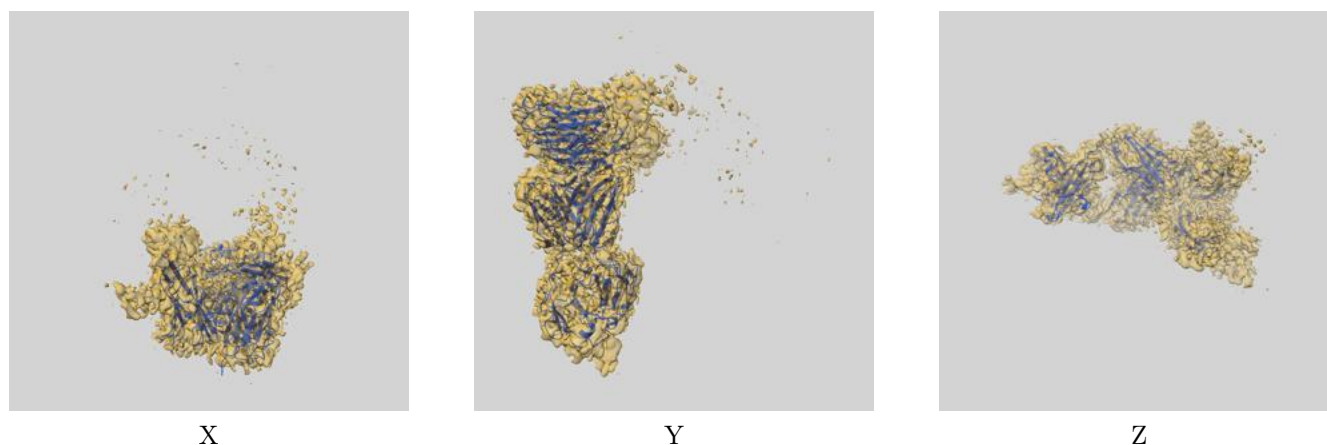
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.50	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	3.72	4.36	3.80

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

9 Map-model fit [i](#)

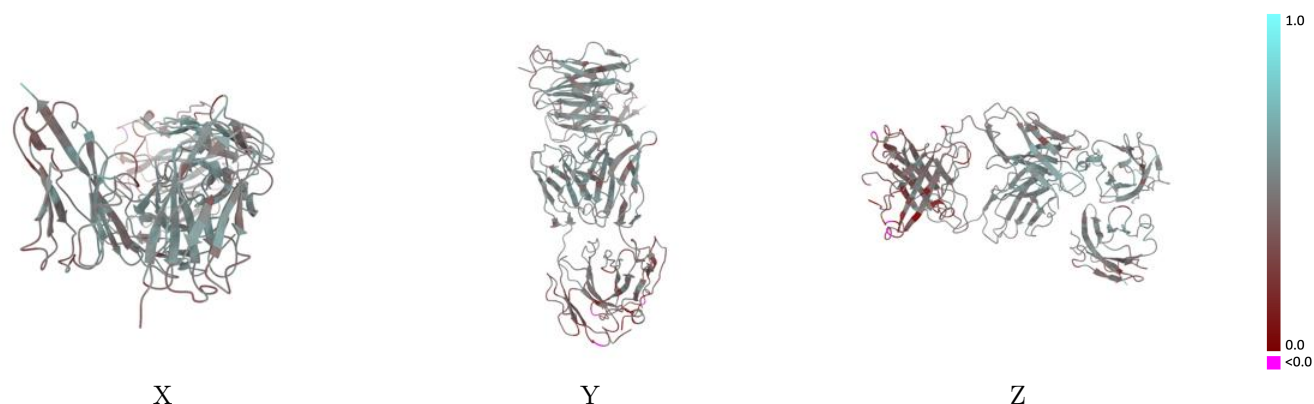
This section contains information regarding the fit between EMDB map EMD-62751 and PDB model 9L1P. Per-residue inclusion information can be found in [section 3](#) on [page 6](#).

9.1 Map-model overlay [i](#)



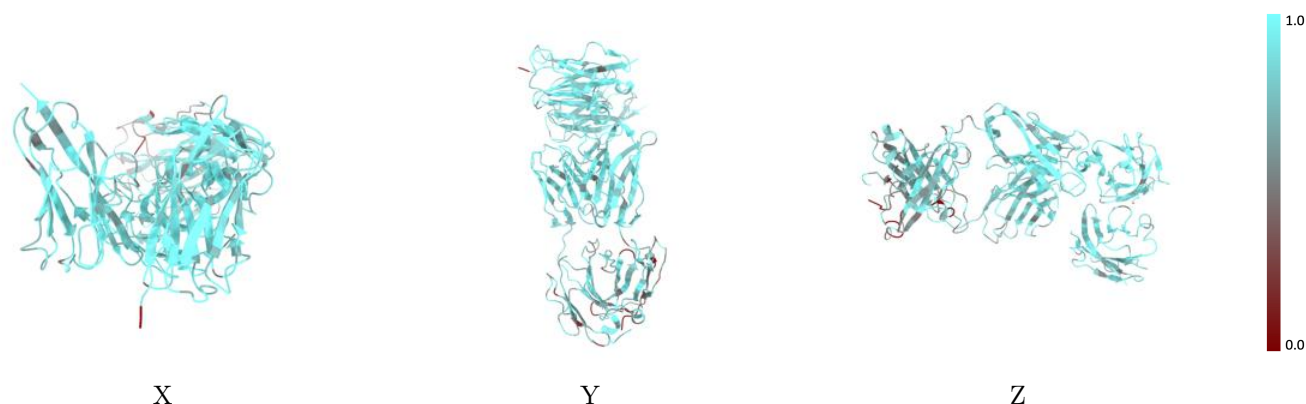
The images above show the 3D surface view of the map at the recommended contour level 0.02 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



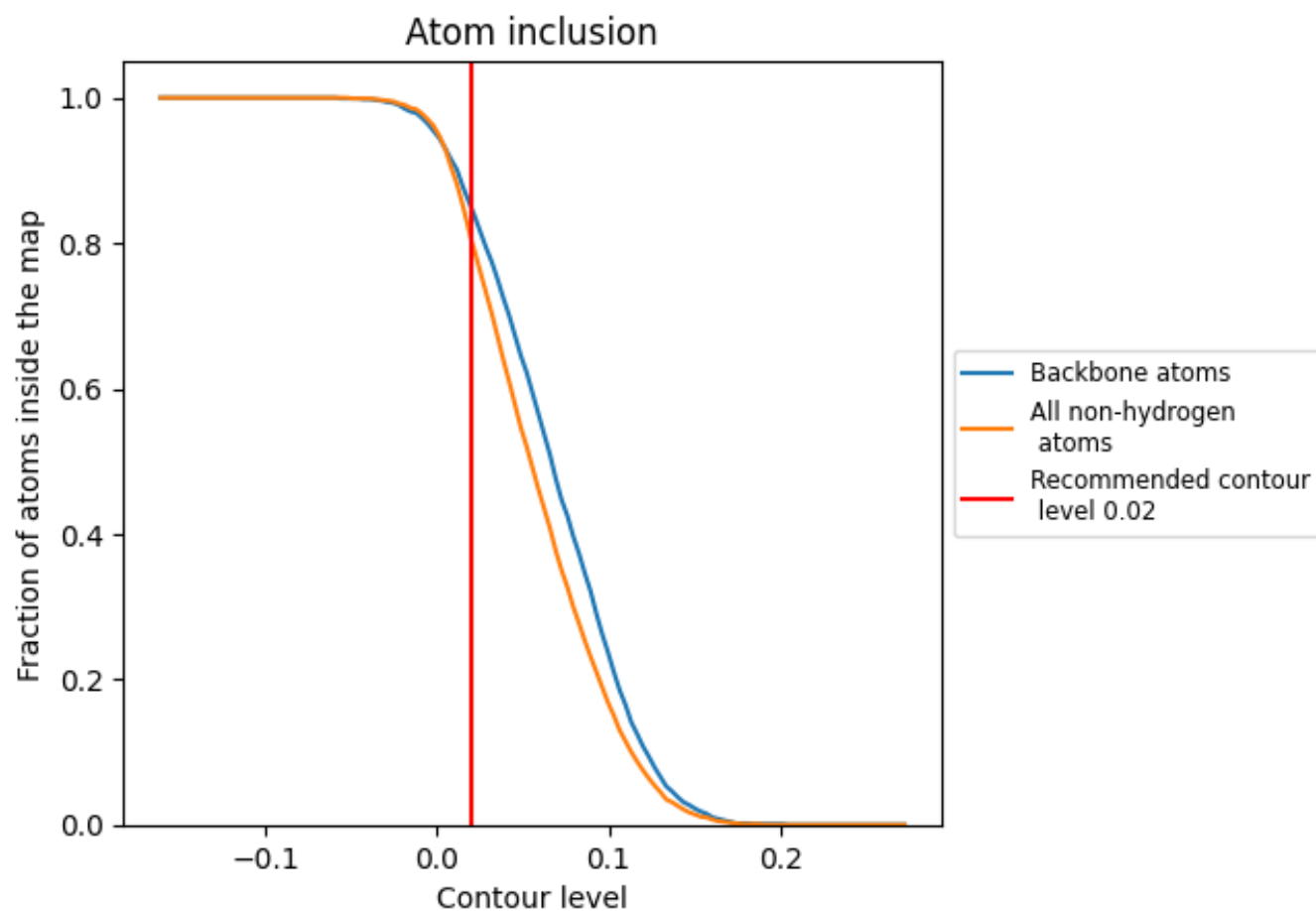
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.02).

9.4 Atom inclusion [i](#)



At the recommended contour level, 85% of all backbone atoms, 80% 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 (0.02) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8030	0.4560
B	0.8400	0.4680
E	0.8810	0.4990
H	0.7560	0.4400
L	0.7900	0.4420

