



## Full wwPDB EM Validation Report ⓘ

Nov 20, 2022 – 09:22 PM EST

PDB ID : 7MQS  
EMDB ID : EMD-23951  
Title : The insulin receptor ectodomain in complex with three venom hybrid insulin molecules - asymmetric conformation  
Authors : Blakely, A.D.; Xiong, X.; Kim, J.H.; Menting, J.; Schafer, I.B.; Schubert, H.L.; Agrawal, R.; Gutmann, T.; Delaine, C.; Zhang, Y.; Artik, G.O.; Merriman, A.; Eckert, D.; Lawrence, M.C.; Coskun, U.; Fisher, S.J.; Forbes, B.E.; Safavi-Hemami, H.; Hill, C.P.; Chou, D.H.C.  
Deposited on : 2021-05-06  
Resolution : 4.40 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](mailto: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.dev43  
MolProbity : 4.02b-467  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
MapQ : 1.9.9  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.31.2

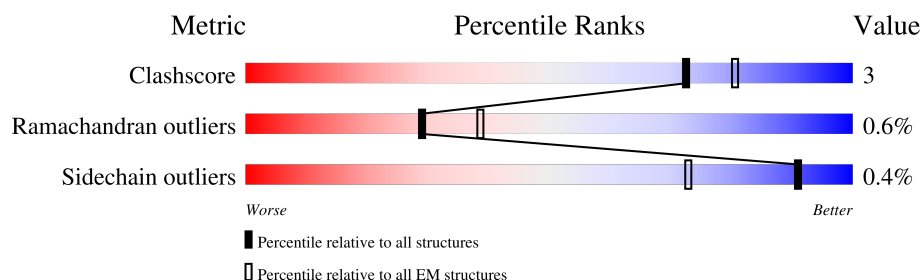
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*ELECTRON MICROSCOPY*

The reported resolution of this entry is 4.40 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)
Clashscore	158937	4297
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826

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  $\geq 3$ , 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	E	916	<div> <div>27%</div> <div>77%</div> <div>11%</div> <div>12%</div> </div>
1	F	916	<div> <div>17%</div> <div>77%</div> <div>11%</div> <div>•</div> <div>11%</div> </div>
2	A	24	<div> <div>17%</div> <div>79%</div> <div>17%</div> <div>•</div> </div>
2	C	24	<div> <div>88%</div> <div>83%</div> <div>17%</div> </div>
2	G	24	<div> <div>8%</div> <div>79%</div> <div>21%</div> </div>
3	B	22	<div> <div>50%</div> <div>23%</div> <div>27%</div> </div>
3	D	22	<div> <div>55%</div> <div>59%</div> <div>9%</div> <div>5%</div> <div>27%</div> </div>
3	H	22	<div> <div>68%</div> <div>14%</div> <div>18%</div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 14028 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 Isoform Short of Insulin receptor.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	E	806	Total	C	N	O	S	0	0
			6521	4139	1124	1210	48		
1	F	811	Total	C	N	O	S	0	0
			6563	4164	1131	1220	48		

- Molecule 2 is a protein called Insulin A chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	A	24	Total	C	N	O	S	0	0
			188	115	30	39	4		
2	C	24	Total	C	N	O	S	0	0
			188	115	30	39	4		
2	G	24	Total	C	N	O	S	0	0
			188	115	30	39	4		

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	21	HIS	ASN	engineered mutation	UNP P01308
A	22	SER	-	insertion	UNP P01308
A	23	LEU	-	insertion	UNP P01308
A	24	GLN	-	insertion	UNP P01308
C	21	HIS	ASN	engineered mutation	UNP P01308
C	22	SER	-	insertion	UNP P01308
C	23	LEU	-	insertion	UNP P01308
C	24	GLN	-	insertion	UNP P01308
G	21	HIS	ASN	engineered mutation	UNP P01308
G	22	SER	-	insertion	UNP P01308
G	23	LEU	-	insertion	UNP P01308
G	24	GLN	-	insertion	UNP P01308

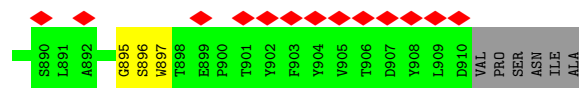
- Molecule 3 is a protein called Insulin B chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	B	16	Total 121	C 79	N 18	O 22	S 2	0	0
3	D	16	Total 121	C 79	N 18	O 22	S 2	0	0
3	H	18	Total 138	C 88	N 22	O 26	S 2	0	0

There are 6 discrepancies between the modelled and reference sequences:

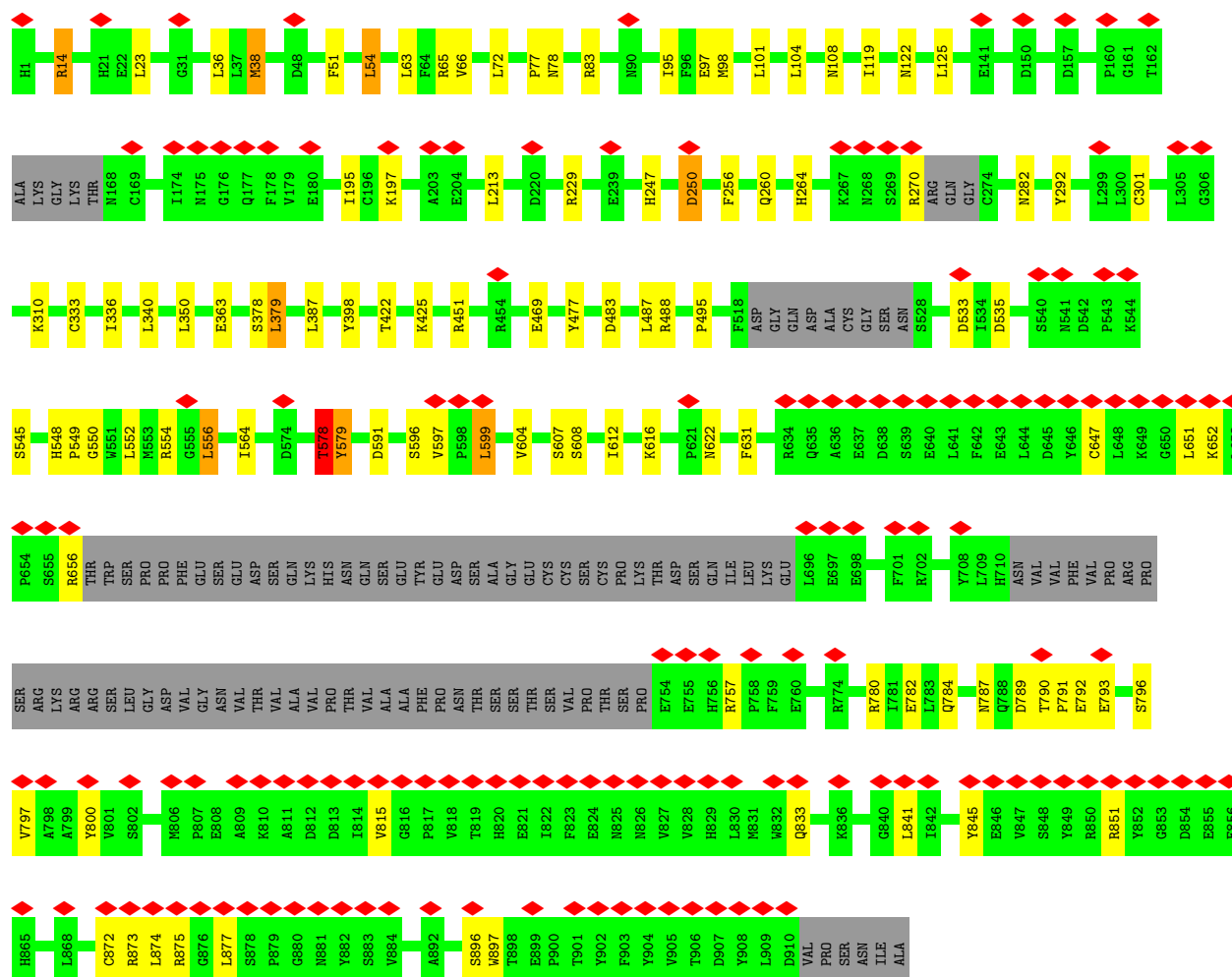
Chain	Residue	Modelled	Actual	Comment	Reference
B	10	GLU	HIS	engineered mutation	UNP P01308
B	20	LEU	GLY	engineered mutation	UNP P01308
D	10	GLU	HIS	engineered mutation	UNP P01308
D	20	LEU	GLY	engineered mutation	UNP P01308
H	10	GLU	HIS	engineered mutation	UNP P01308
H	20	LEU	GLY	engineered mutation	UNP P01308





• Molecule 1: Isoform Short of Insulin receptor

Chain F: 17% 77% 11% 11%



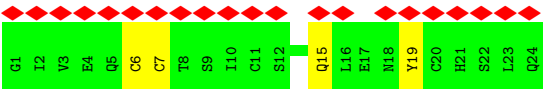
• Molecule 2: Insulin A chain

Chain A: 17% 79% 17%

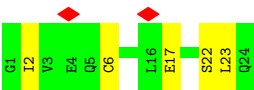
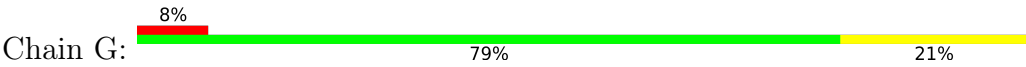


• Molecule 2: Insulin A chain

Chain C: 88% 83% 17%



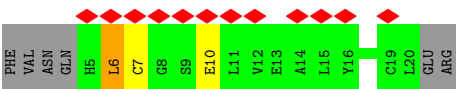
• Molecule 2: Insulin A chain



• Molecule 3: Insulin B chain



• Molecule 3: Insulin B chain



• Molecule 3: Insulin B chain



## 4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	43457	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	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	2.019	Depositor
Minimum map value	-1.588	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.048	Depositor
Recommended contour level	0.455	Depositor
Map size ( $\text{\AA}$ )	349.44, 349.44, 349.44	wwPDB
Map dimensions	256, 256, 256	wwPDB
Map angles ( $^\circ$ )	90.0, 90.0, 90.0	wwPDB
Pixel spacing ( $\text{\AA}$ )	1.365, 1.365, 1.365	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	E	0.90	1/6678 (0.0%)	0.95	20/9050 (0.2%)
1	F	0.90	1/6725 (0.0%)	0.98	28/9118 (0.3%)
2	A	0.88	0/190	1.17	2/255 (0.8%)
2	C	0.88	0/190	0.92	0/255
2	G	1.11	0/190	1.36	3/255 (1.2%)
3	B	0.96	0/122	1.69	4/165 (2.4%)
3	D	0.83	0/122	1.22	1/165 (0.6%)
3	H	0.86	0/139	1.20	0/188
All	All	0.90	2/14356 (0.0%)	0.99	58/19451 (0.3%)

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	E	0	4
1	F	0	4
2	C	0	2
3	B	0	1
All	All	0	11

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	790	THR	C-N	6.29	1.46	1.34
1	F	790	THR	C-N	5.07	1.43	1.34

All (58) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	37	LEU	CA-CB-CG	11.73	142.28	115.30
1	F	250	ASP	CB-CG-OD1	11.34	128.50	118.30

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	533	ASP	CB-CG-OD2	10.88	128.09	118.30
1	F	23	LEU	CA-CB-CG	10.14	138.61	115.30
1	E	330	LEU	CA-CB-CG	9.69	137.58	115.30
1	F	652	LYS	C-N-CA	9.05	144.32	121.70
1	F	72	LEU	CA-CB-CG	8.54	134.95	115.30
1	F	841	LEU	CA-CB-CG	8.17	134.10	115.30
3	B	6	LEU	CB-CG-CD2	8.15	124.85	111.00
1	E	844	LEU	CA-CB-CG	8.12	133.98	115.30
1	E	23	LEU	CA-CB-CG	8.00	133.70	115.30
3	B	15	LEU	CA-CB-CG	7.78	133.19	115.30
1	E	651	LEU	CA-CB-CG	7.70	133.01	115.30
1	E	17	LEU	CA-CB-CG	7.23	131.93	115.30
1	F	38	MET	CA-CB-CG	7.13	125.42	113.30
1	F	591	ASP	CB-CG-OD1	7.04	124.63	118.30
3	B	11	LEU	CA-CB-CG	6.95	131.29	115.30
2	G	23	LEU	CB-CG-CD2	6.85	122.64	111.00
1	E	870	ARG	CA-CB-CG	6.81	128.39	113.40
1	E	843	VAL	CG1-CB-CG2	-6.71	100.17	110.90
1	E	505	LEU	CA-CB-CG	6.47	130.18	115.30
1	F	564	ILE	CG1-CB-CG2	-6.38	97.36	111.40
1	F	36	LEU	CA-CB-CG	6.25	129.68	115.30
1	F	38	MET	CB-CG-SD	6.24	131.11	112.40
2	A	14	TYR	CB-CG-CD1	-6.21	117.27	121.00
1	F	487	LEU	CA-CB-CG	6.20	129.56	115.30
2	G	17	GLU	N-CA-CB	5.97	121.36	110.60
1	F	98	MET	CA-CB-CG	5.91	123.35	113.30
3	B	6	LEU	CA-CB-CG	5.84	128.73	115.30
2	A	16	LEU	CA-CB-CG	5.77	128.58	115.30
1	E	99	VAL	C-N-CA	5.77	136.12	121.70
1	E	37	LEU	CB-CG-CD1	5.72	120.73	111.00
1	E	806	MET	CA-CB-CG	5.67	122.95	113.30
1	E	648	LEU	CA-CB-CG	5.66	128.31	115.30
1	E	250	ASP	CB-CG-OD1	5.62	123.35	118.30
1	E	63	LEU	CA-CB-CG	5.54	128.04	115.30
1	E	841	LEU	CA-CB-CG	5.49	127.92	115.30
1	F	469	GLU	C-N-CA	5.49	135.41	121.70
1	F	791	PRO	C-N-CA	5.46	135.36	121.70
1	F	556	LEU	C-N-CA	5.42	135.25	121.70
1	F	578	THR	C-N-CA	5.38	135.15	121.70
3	D	6	LEU	CA-CB-CG	5.29	127.47	115.30
1	E	36	LEU	CA-CB-CG	5.28	127.45	115.30
1	E	791	PRO	C-N-CA	5.25	134.84	121.70

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	379	LEU	CA-CB-CG	5.20	127.26	115.30
1	F	54	LEU	CA-CB-CG	5.19	127.23	115.30
1	F	14	ARG	C-N-CA	5.17	134.62	121.70
2	G	17	GLU	CA-CB-CG	5.15	124.74	113.40
1	F	552	LEU	CA-CB-CG	5.14	127.13	115.30
1	F	104	LEU	CA-CB-CG	5.13	127.10	115.30
1	F	545	SER	C-N-CA	5.12	134.49	121.70
1	F	599	LEU	C-N-CA	5.12	134.49	121.70
1	F	350	LEU	CA-CB-CG	5.11	127.04	115.30
1	F	101	LEU	CA-CB-CG	5.10	127.03	115.30
1	F	651	LEU	CA-CB-CG	5.09	127.01	115.30
1	E	98	MET	CG-SD-CE	5.06	108.30	100.20
1	E	434	LEU	CA-CB-CG	5.04	126.88	115.30
1	F	250	ASP	CB-CG-OD2	-5.03	113.77	118.30

There are no chirality outliers.

All (11) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	B	7	CYS	Peptide
2	C	6	CYS	Peptide
2	C	7	CYS	Peptide
1	E	111	ASN	Peptide
1	E	152	ASN	Peptide
1	E	276	GLN	Peptide
1	E	895	GLY	Peptide
1	F	495	PRO	Peptide
1	F	554	ARG	Peptide
1	F	578	THR	Peptide
1	F	647	CYS	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	E	6521	0	6339	42	0
1	F	6563	0	6364	45	0

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	188	0	174	3	0
2	C	188	0	174	1	0
2	G	188	0	174	2	0
3	B	121	0	121	1	0
3	D	121	0	121	1	0
3	H	138	0	135	1	0
All	All	14028	0	13602	93	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 3.

All (93) 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:E:351:ALA:O	1:E:355:GLU:HB2	1.99	0.62
1:E:256:PHE:O	1:E:260:GLN:HB2	2.01	0.61
2:G:2:ILE:O	2:G:6:CYS:HB3	2.07	0.54
1:E:261:ASP:O	1:E:265:LYS:HB2	2.08	0.54
1:F:780:ARG:HD2	1:F:800:TYR:HB3	1.92	0.52
1:F:63:LEU:HD12	1:F:95:ILE:HG12	1.92	0.51
1:E:476:SER:H	1:E:489:TRP:HA	1.74	0.51
1:E:407:ASN:HA	1:E:433:LYS:HD2	1.93	0.50
1:F:256:PHE:O	1:F:260:GLN:HB2	2.11	0.50
1:E:477:TYR:HB3	1:E:488:ARG:HB2	1.94	0.48
1:F:398:TYR:HA	1:F:425:LYS:H	1.78	0.48
1:F:872:CYS:SG	1:F:873:ARG:N	2.87	0.48
1:F:292:TYR:HB3	1:F:301:CYS:HB3	1.96	0.48
1:E:599:LEU:HD12	1:E:616:LYS:HD2	1.97	0.47
1:F:310:LYS:NZ	1:F:333:CYS:SG	2.88	0.47
1:E:310:LYS:NZ	1:E:312:CYS:SG	2.87	0.46
1:F:38:MET:HB2	1:F:66:VAL:HA	1.96	0.46
1:F:604:VAL:HB	1:F:612:ILE:HB	1.98	0.46
1:E:646:TYR:HB2	1:E:844:LEU:HD21	1.97	0.46
1:F:874:LEU:HB3	1:F:877:LEU:HD11	1.98	0.46
2:A:16:LEU:HA	2:A:19:TYR:HB2	1.98	0.46
1:E:535:ASP:OD1	1:E:535:ASP:N	2.49	0.46
1:F:78:ASN:HA	1:F:108:ASN:HD22	1.81	0.46
1:E:631:PHE:HB2	1:E:782:GLU:HB3	1.99	0.45
1:F:247:HIS:ND1	1:F:282:ASN:O	2.50	0.45
1:E:48:ASP:N	1:E:48:ASP:OD1	2.50	0.45
1:F:83:ARG:NH2	1:F:250:ASP:O	2.50	0.45

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:535:ASP:OD1	1:F:535:ASP:N	2.50	0.45
1:F:845:TYR:HB2	1:F:861:VAL:HB	1.98	0.45
1:E:119:ILE:HB	1:E:145:ILE:HA	1.98	0.44
1:E:72:LEU:HA	1:E:75:LEU:HB2	2.00	0.44
1:F:607:SER:OG	1:F:608:SER:N	2.50	0.44
1:F:363:GLU:HG3	1:F:387:LEU:HB3	2.00	0.44
1:F:422:THR:O	1:F:451:ARG:NH2	2.50	0.44
1:E:87:LEU:HD21	1:E:114:ARG:HG2	1.98	0.44
1:E:775:HIS:NE2	1:E:837:GLU:OE1	2.50	0.44
1:E:177:GLN:O	1:E:181:ARG:NH2	2.51	0.44
1:E:594:ASN:O	1:E:622:ASN:ND2	2.50	0.44
1:E:851:ARG:HA	1:E:882:TYR:HA	1.99	0.44
1:F:578:THR:OG1	1:F:579:TYR:N	2.50	0.44
1:E:596:SER:OG	1:E:597:VAL:N	2.50	0.44
1:F:14:ARG:NH1	2:G:22:SER:O	2.50	0.44
1:F:77:PRO:O	1:F:108:ASN:ND2	2.50	0.44
1:F:195:ILE:O	1:F:197:LYS:NZ	2.50	0.44
1:E:247:HIS:ND1	1:E:282:ASN:O	2.51	0.44
1:E:654:PRO:HG2	1:E:656:ARG:HH21	1.83	0.44
1:F:596:SER:OG	1:F:597:VAL:N	2.51	0.44
1:F:784:GLN:NE2	1:F:796:SER:O	2.51	0.44
1:F:488:ARG:NH1	1:F:548:HIS:O	2.51	0.43
1:E:821:GLU:OE2	1:E:829:HIS:ND1	2.52	0.43
1:F:796:SER:OG	1:F:797:VAL:N	2.51	0.43
1:E:323:SER:O	1:E:327:ALA:HB2	2.18	0.43
1:F:789:ASP:HA	1:F:793:GLU:HA	2.00	0.43
1:E:851:ARG:NH1	1:E:879:PRO:O	2.51	0.43
1:E:641:LEU:HA	1:E:644:LEU:HG	2.01	0.43
1:F:622:ASN:O	1:F:787:ASN:ND2	2.50	0.43
1:E:458:ALA:O	1:E:462:ASN:ND2	2.52	0.43
1:E:607:SER:OG	1:E:608:SER:N	2.52	0.43
1:F:599:LEU:HB2	1:F:616:LYS:HD3	2.00	0.43
1:E:476:SER:OG	1:E:477:TYR:N	2.52	0.43
1:F:95:ILE:HB	1:F:119:ILE:HG23	2.00	0.43
1:E:32:HIS:HB3	1:E:59:ASP:HB2	2.01	0.42
1:E:896:SER:OG	1:E:897:TRP:N	2.51	0.42
1:F:336:ILE:HD13	1:F:340:LEU:HD21	2.01	0.42
2:A:9:SER:O	3:B:5:HIS:N	2.53	0.42
2:C:15:GLN:O	2:C:19:TYR:HB2	2.19	0.42
1:E:192:CYS:SG	1:E:201:CYS:N	2.86	0.42
1:F:378:SER:OG	1:F:379:LEU:N	2.51	0.42

*Continued on next page...*

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:815:VAL:O	1:F:833:GLN:NE2	2.51	0.42
1:F:873:ARG:O	1:F:875:ARG:NH2	2.50	0.42
1:E:436:LEU:O	1:E:440:HIS:ND1	2.53	0.42
1:E:494:PRO:HG2	1:E:497:PHE:HA	2.00	0.42
1:F:213:LEU:HD22	1:F:229:ARG:HA	2.01	0.42
1:F:483:ASP:HB3	1:F:556:LEU:H	1.84	0.42
3:D:6:LEU:HD12	3:D:10:GLU:HB3	2.01	0.42
1:E:467:SER:O	1:E:582:LYS:NZ	2.52	0.41
1:F:122:ASN:HD22	1:F:125:LEU:HD22	1.84	0.41
1:E:292:TYR:HB3	1:E:301:CYS:HB3	2.02	0.41
1:E:385:LEU:O	1:E:417:HIS:NE2	2.53	0.41
1:F:488:ARG:HD3	1:F:550:GLY:HA3	2.02	0.41
1:E:626:THR:H	1:E:787:ASN:HA	1.85	0.41
1:E:469:GLU:HB2	1:E:582:LYS:HE2	2.01	0.41
1:F:896:SER:OG	1:F:897:TRP:N	2.53	0.41
1:F:631:PHE:HB2	1:F:782:GLU:HB3	2.02	0.41
1:E:133:TRP:HB3	1:E:137:LEU:HD13	2.01	0.41
1:E:266:CYS:HB3	1:E:274:CYS:HB3	1.76	0.41
1:F:51:PHE:HB3	1:F:54:LEU:HD13	2.03	0.40
1:F:477:TYR:HB3	1:F:488:ARG:HB2	2.03	0.40
3:H:10:GLU:HA	3:H:13:GLU:HB2	2.03	0.40
1:F:65:ARG:NH2	1:F:97:GLU:OE2	2.55	0.40
1:F:260:GLN:O	1:F:264:HIS:ND1	2.50	0.40
1:E:488:ARG:HH11	1:E:548:HIS:HB2	1.86	0.40
1:F:488:ARG:NH2	2:A:17:GLU:OE1	2.53	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	E	792/916 (86%)	723 (91%)	65 (8%)	4 (0%)	29 68

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	F	799/916 (87%)	708 (89%)	87 (11%)	4 (0%)	29	68
2	A	22/24 (92%)	21 (96%)	1 (4%)	0	100	100
2	C	22/24 (92%)	19 (86%)	3 (14%)	0	100	100
2	G	22/24 (92%)	20 (91%)	2 (9%)	0	100	100
3	B	14/22 (64%)	12 (86%)	2 (14%)	0	100	100
3	D	14/22 (64%)	12 (86%)	1 (7%)	1 (7%)	1	16
3	H	16/22 (73%)	12 (75%)	3 (19%)	1 (6%)	1	19
All	All	1701/1970 (86%)	1527 (90%)	164 (10%)	10 (1%)	29	65

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	F	792	GLU
3	D	7	CYS
1	E	540	SER
1	F	549	PRO
1	F	579	TYR
1	E	15	ASN
1	E	792	GLU
3	H	7	CYS
1	F	578	THR
1	E	543	PRO

### 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	E	735/830 (89%)	733 (100%)	2 (0%)	92	95
1	F	739/830 (89%)	735 (100%)	4 (0%)	88	93
2	A	23/23 (100%)	23 (100%)	0	100	100
2	C	23/23 (100%)	23 (100%)	0	100	100
2	G	23/23 (100%)	23 (100%)	0	100	100

Continued on next page...

*Continued from previous page...*

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	B	14/20 (70%)	14 (100%)	0	100	100
3	D	14/20 (70%)	14 (100%)	0	100	100
3	H	16/20 (80%)	16 (100%)	0	100	100
All	All	1587/1789 (89%)	1581 (100%)	6 (0%)	91	94

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	40	LYS
1	E	870	ARG
1	F	270	ARG
1	F	656	ARG
1	F	757	ARG
1	F	851	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 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 monosaccharides 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.

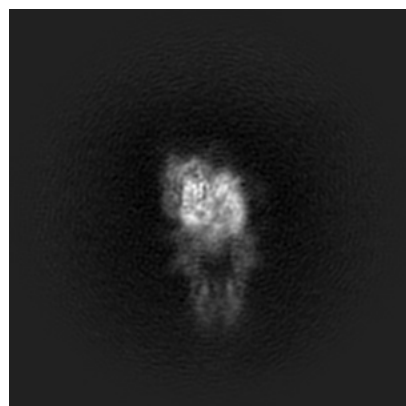
## 6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-23951. These allow visual inspection of the internal detail of the map and identification of artifacts.

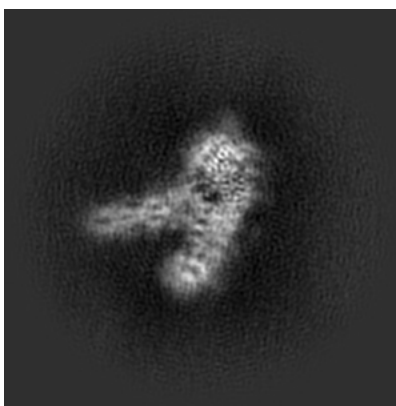
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

### 6.1 Orthogonal projections [i](#)

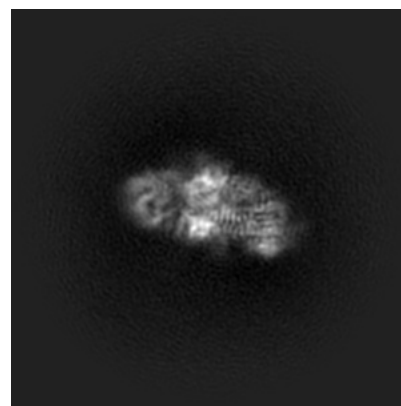
#### 6.1.1 Primary map



X

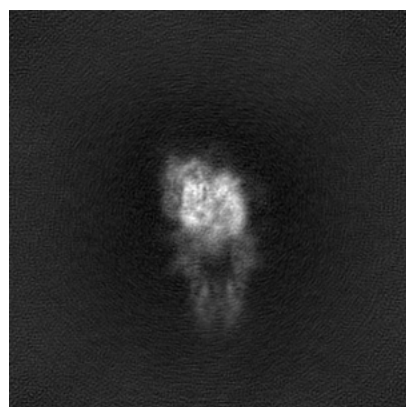


Y

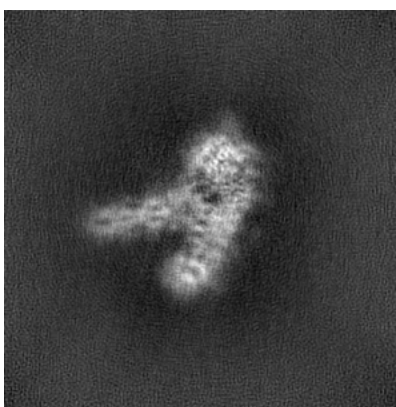


Z

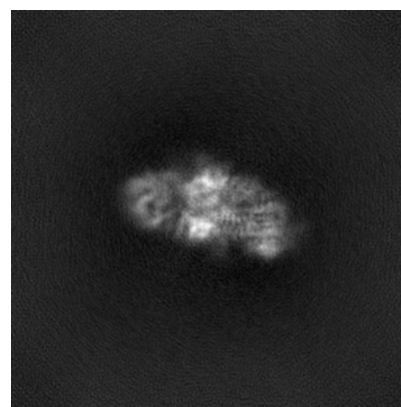
#### 6.1.2 Raw map



X



Y

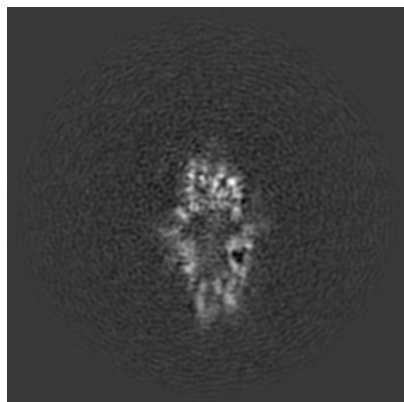


Z

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

## 6.2 Central slices [i](#)

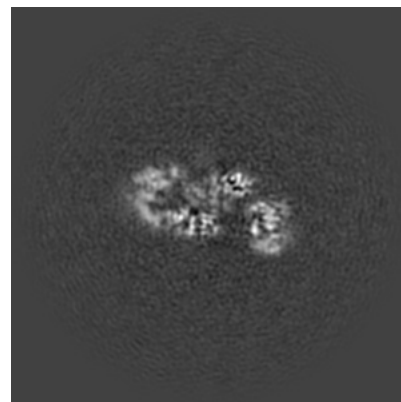
### 6.2.1 Primary map



X Index: 128

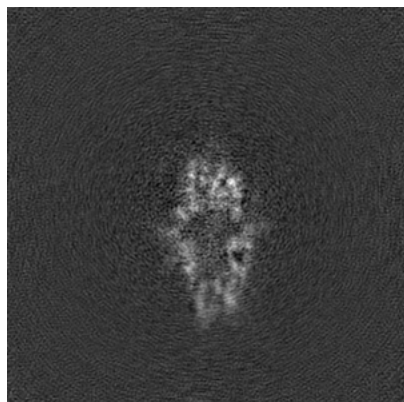


Y Index: 128

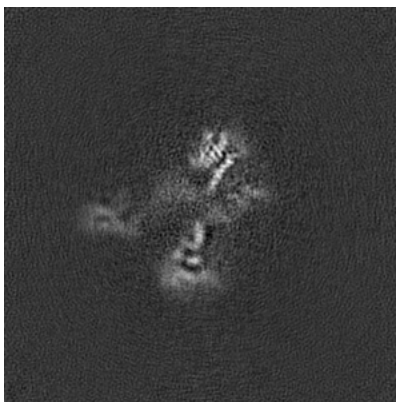


Z Index: 128

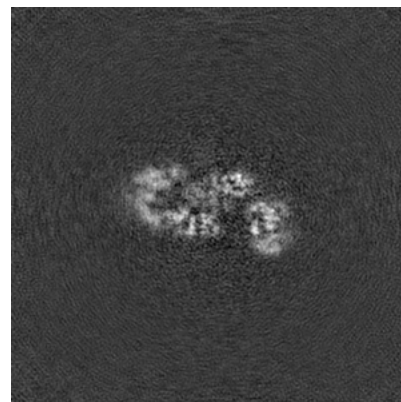
### 6.2.2 Raw map



X Index: 128



Y Index: 128

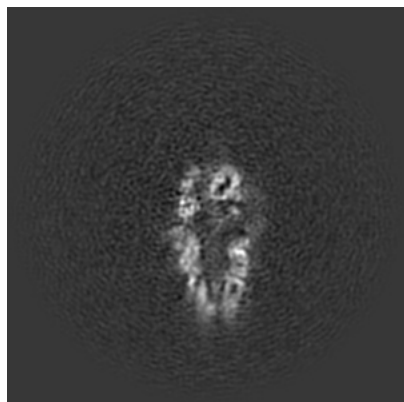


Z Index: 128

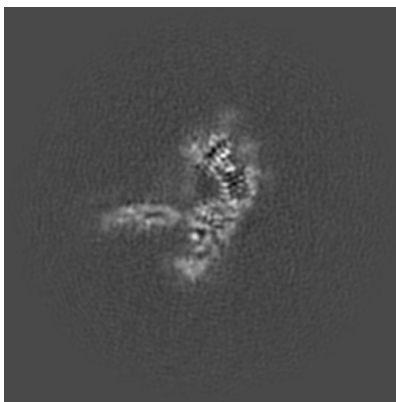
The images above show central slices of the map in three orthogonal directions.

## 6.3 Largest variance slices [i](#)

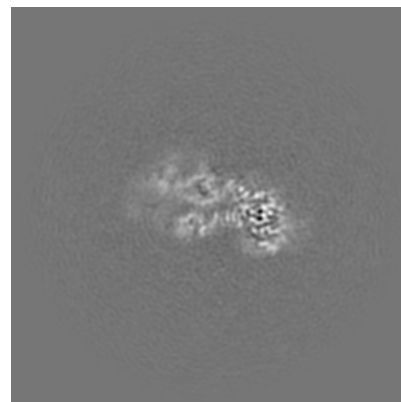
### 6.3.1 Primary map



X Index: 125

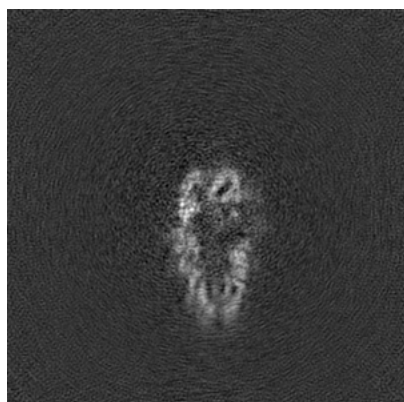


Y Index: 118

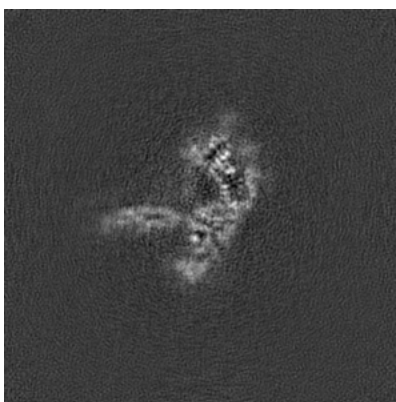


Z Index: 138

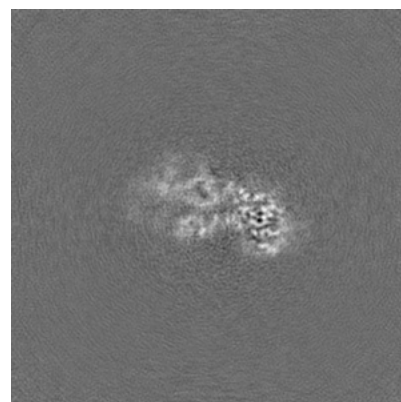
### 6.3.2 Raw map



X Index: 124



Y Index: 118

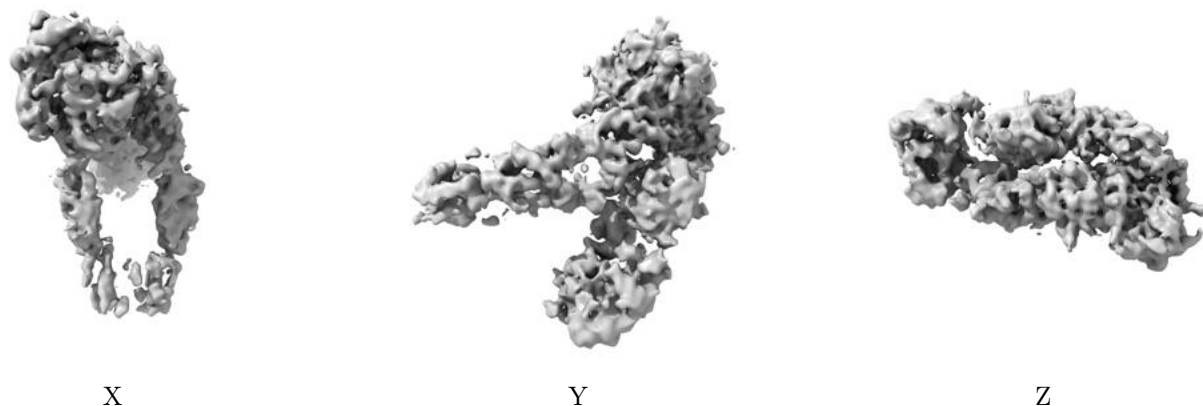


Z Index: 138

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

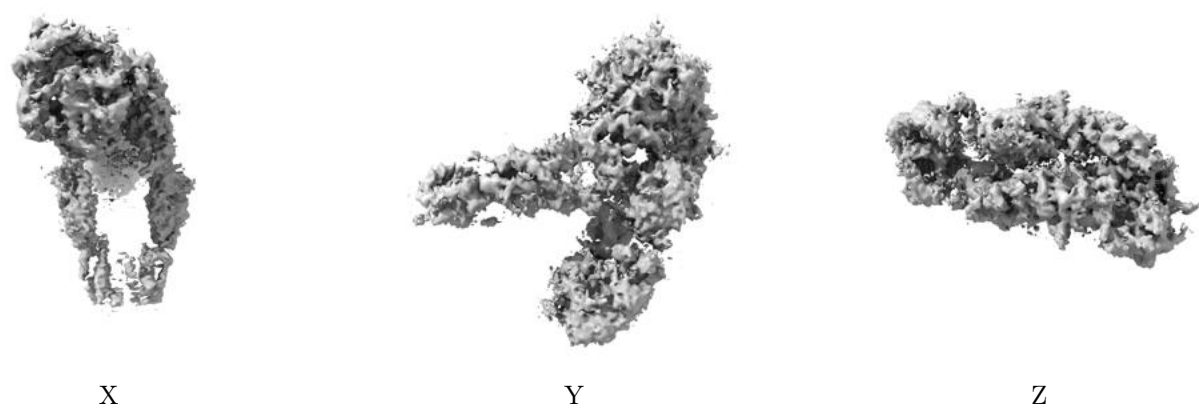
## 6.4 Orthogonal surface views [i](#)

### 6.4.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.455. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

### 6.4.2 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.

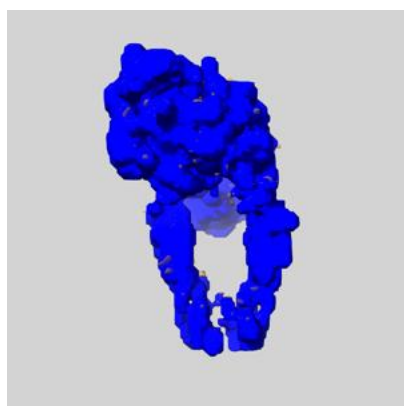
## 6.5 Mask visualisation [i](#)

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

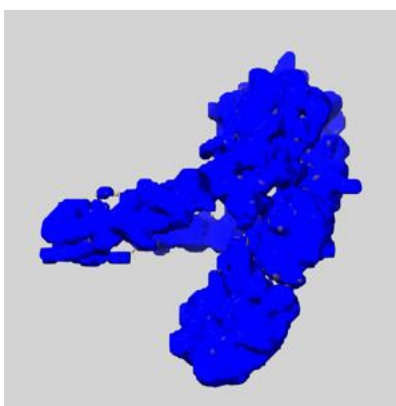
A mask typically either:

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

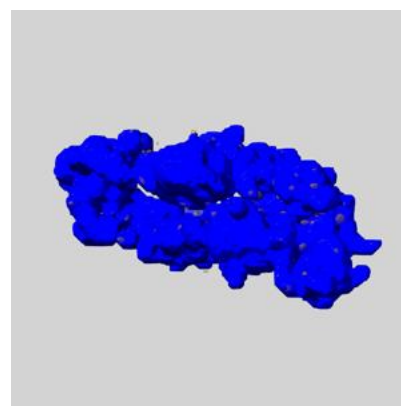
### 6.5.1 emd\_23951\_msk\_1.map [i](#)



X



Y

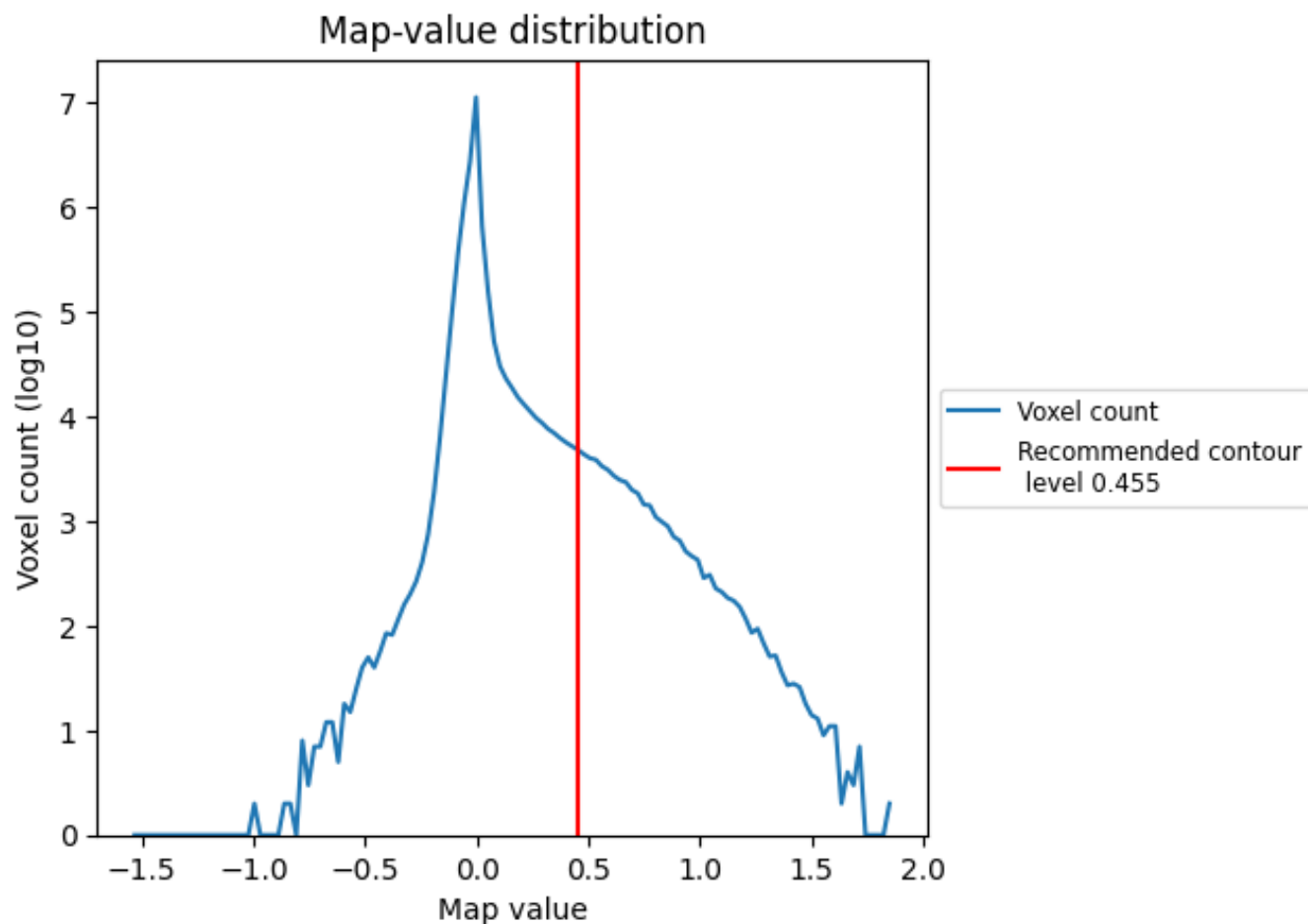


Z

## 7 Map analysis [i](#)

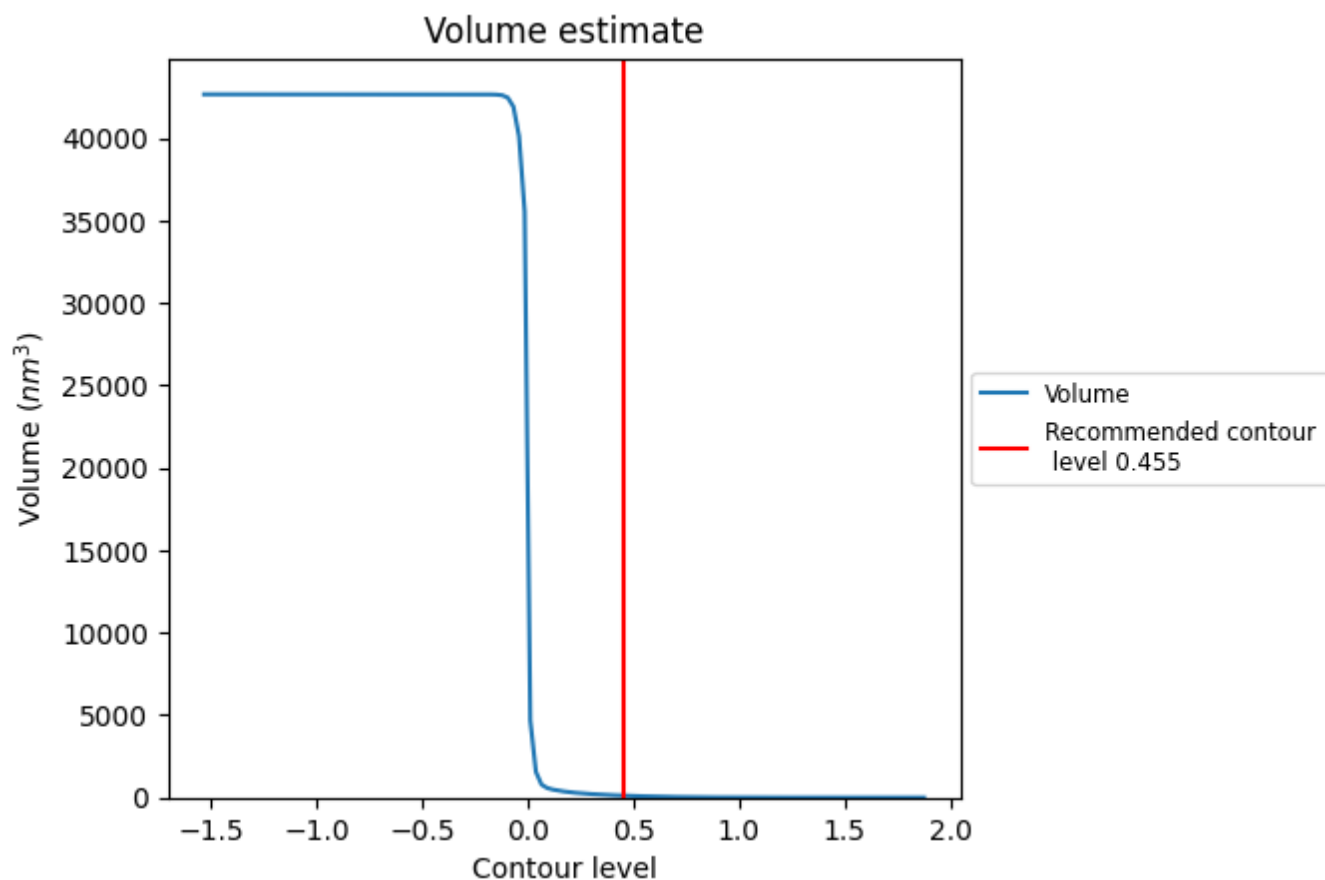
This section contains the results of statistical analysis of the map.

### 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](#)

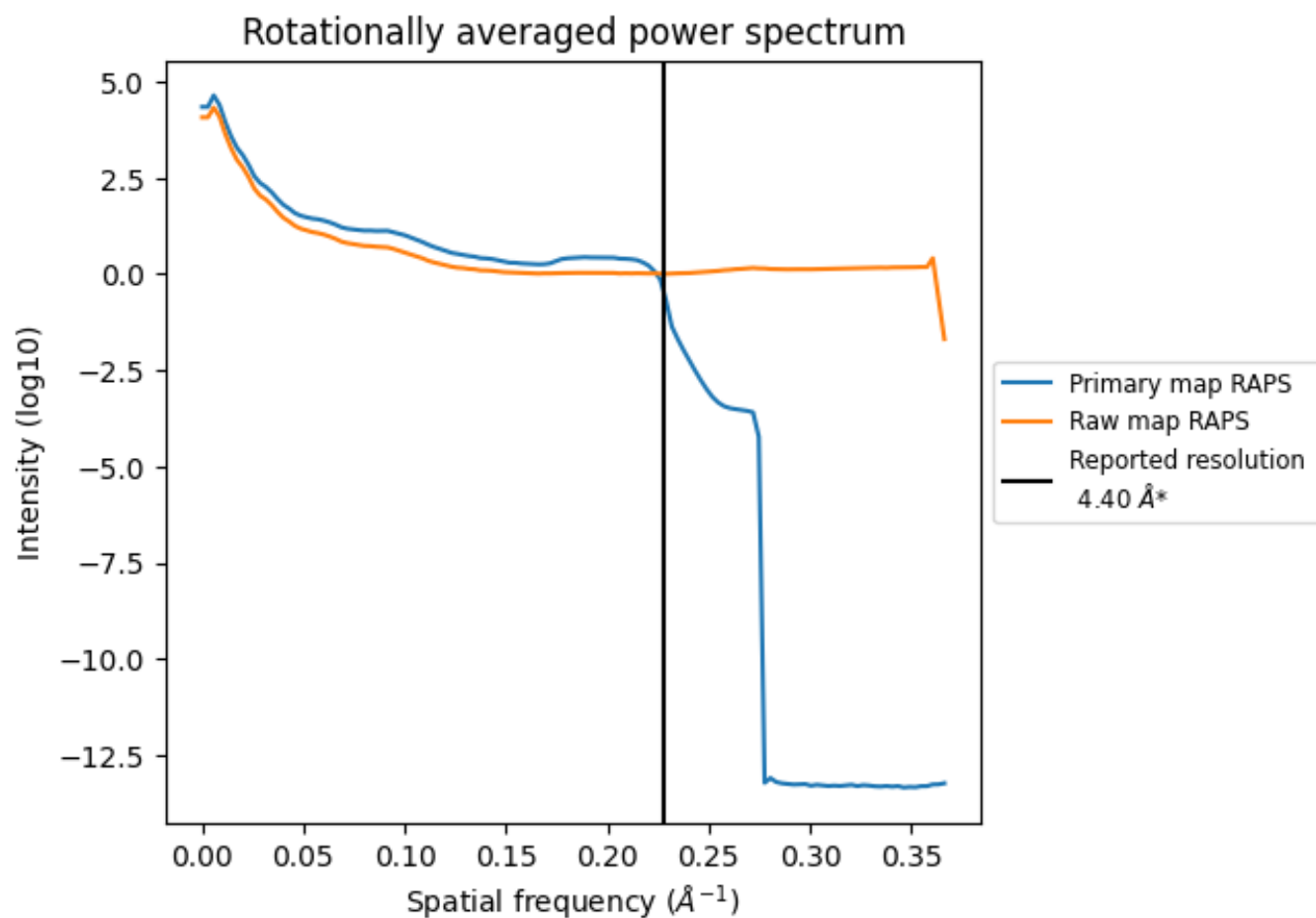


The volume at the recommended contour level is 116 nm<sup>3</sup>; this corresponds to an approximate mass of 105 kDa.

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 ⓘ

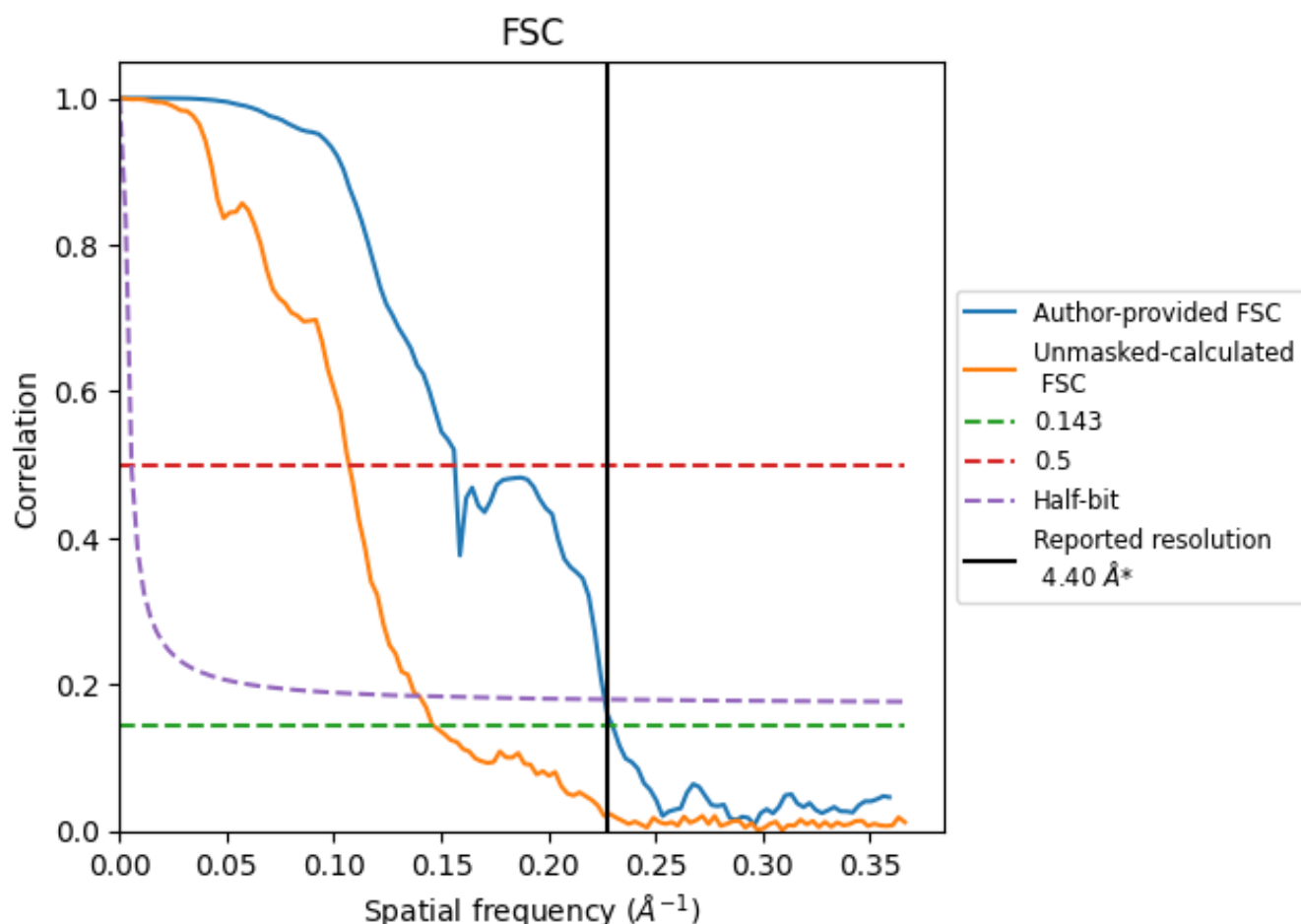


\*Reported resolution corresponds to spatial frequency of 0.227 Å<sup>-1</sup>

## 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 0.227 Å<sup>-1</sup>

## 8.2 Resolution estimates [i](#)

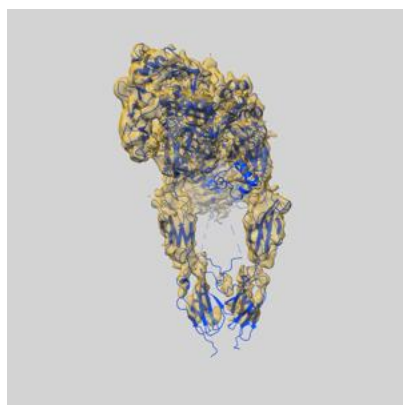
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	4.40	-	-
Author-provided FSC curve	4.35	6.39	4.42
Unmasked-calculated*	6.82	9.34	7.16

\*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 6.82 differs from the reported value 4.4 by more than 10 %

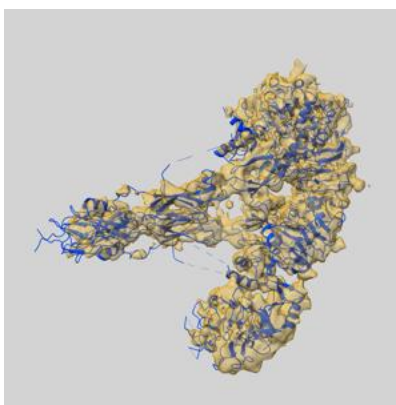
## 9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-23951 and PDB model 7MQS. Per-residue inclusion information can be found in [section 3](#) on [page 5](#).

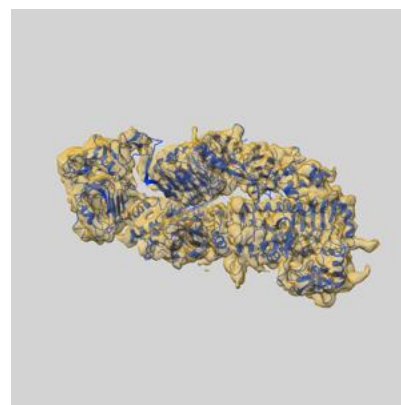
### 9.1 Map-model overlay [i](#)



X



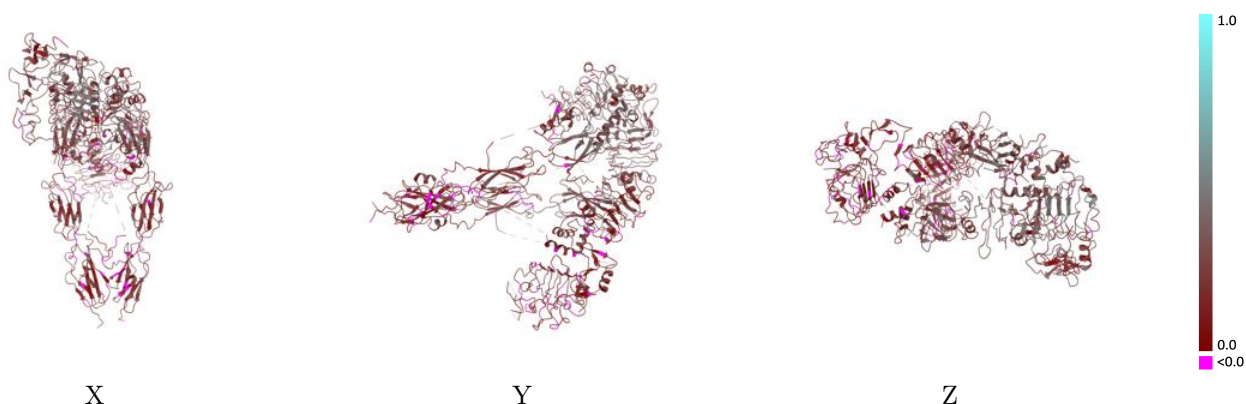
Y



Z

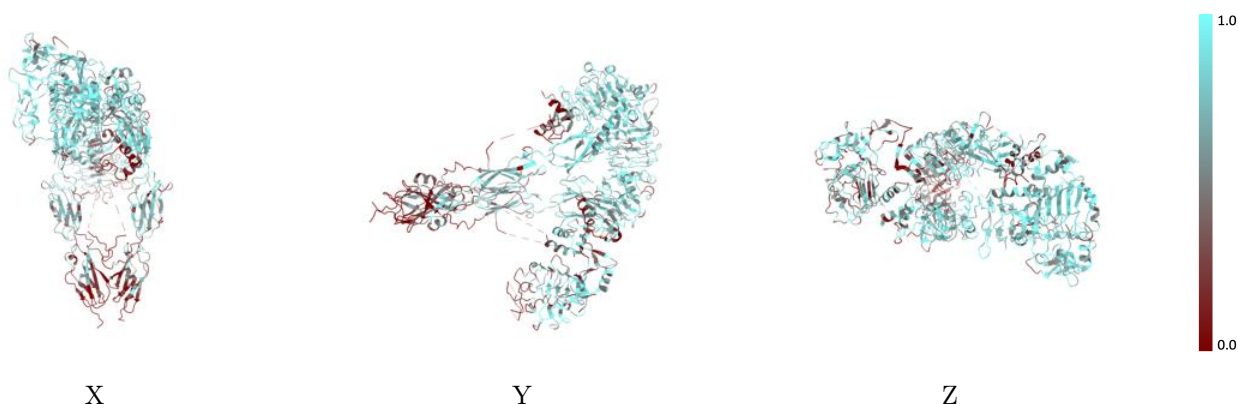
The images above show the 3D surface view of the map at the recommended contour level 0.455 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



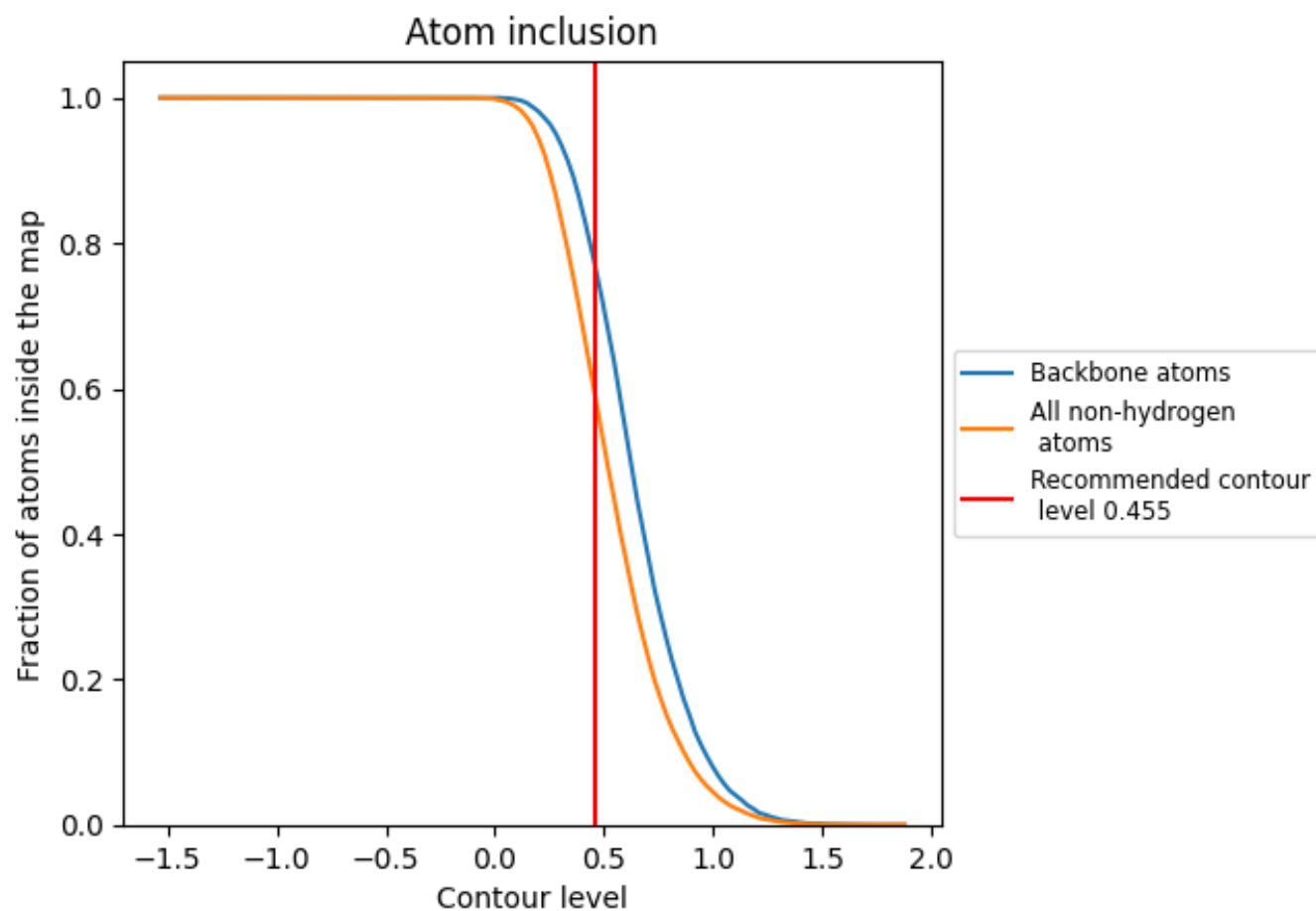
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.455).

## 9.4 Atom inclusion [i](#)



At the recommended contour level, 77% of all backbone atoms, 60% 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 (0.455) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div></div> 0.5967	<div></div> 0.2320
A	<div></div> 0.6667	<div></div> 0.1840
B	<div></div> 0.7000	<div></div> 0.2090
C	<div></div> 0.0914	<div></div> 0.2000
D	<div></div> 0.2333	<div></div> 0.1580
E	<div></div> 0.5637	<div></div> 0.2010
F	<div></div> 0.6399	<div></div> 0.2640
G	<div></div> 0.7151	<div></div> 0.2910
H	<div></div> 0.7664	<div></div> 0.3150

1.0

0.0

<0.0