



Full wwPDB EM Validation Report ⓘ

May 30, 2026 – 01:02 PM EDT

PDB ID : 10EN / pdb_000010en
EMDB ID : EMD-75112
Title : SK3D-Matured in complex with GluN1-GluN2B, full refinement
Authors : Kleeman, S.O.; Furukawa, H.F.
Deposited on : 2026-01-15
Resolution : 3.70 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:

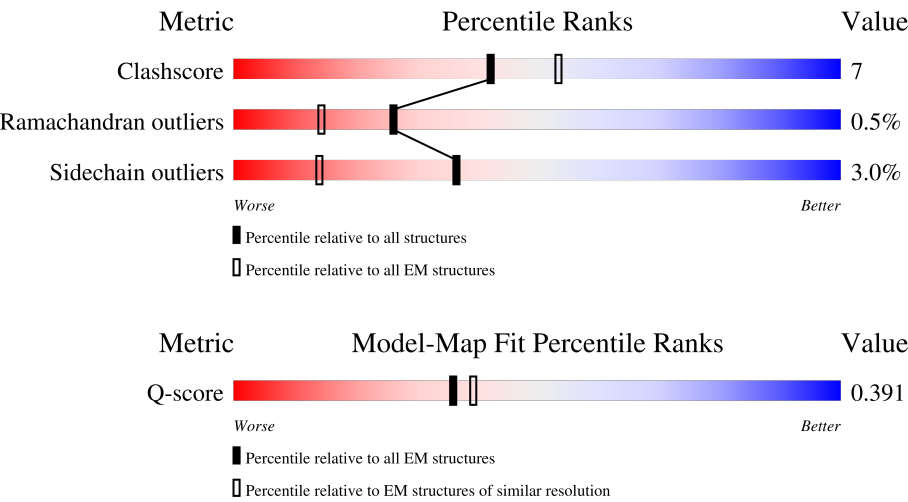
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDb archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh &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

1 Overall quality at a glance i

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

The reported resolution of this entry is 3.70 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	11569 (3.20 - 4.20)

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	817	
1	C	817	
2	B	787	
2	D	787	

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Mol	Chain	Length	Quality of chain
3	E	111	68%30%
3	I	111	68%29%
4	F	120	55%42%
4	J	120	67%30%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 26125 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 4 of Glutamate receptor ionotropic, NMDA 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	795	Total	C	N	O	S	0	0
			5897	3780	1016	1073	28		
1	C	795	Total	C	N	O	S	0	0
			5709	3648	991	1042	28		

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	556	ASN	GLN	conflict	UNP Q5R1P0
A	819	ILE	LEU	conflict	UNP Q5R1P0
A	839	SER	-	expression tag	UNP Q5R1P0
A	840	ARG	-	expression tag	UNP Q5R1P0
A	841	ALA	-	expression tag	UNP Q5R1P0
C	556	ASN	GLN	conflict	UNP Q5R1P0
C	819	ILE	LEU	conflict	UNP Q5R1P0
C	839	SER	-	expression tag	UNP Q5R1P0
C	840	ARG	-	expression tag	UNP Q5R1P0
C	841	ALA	-	expression tag	UNP Q5R1P0

- Molecule 2 is a protein called Glutamate receptor.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	752	Total	C	N	O	S	0	0
			5565	3572	922	1037	34		
2	D	749	Total	C	N	O	S	0	0
			5428	3480	896	1022	30		

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	437	GLU	GLN	conflict	UNP A0A8B7RCM5
B	838	SER	CYS	conflict	UNP A0A8B7RCM5
D	437	GLU	GLN	conflict	UNP A0A8B7RCM5

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Chain	Residue	Modelled	Actual	Comment	Reference
D	838	SER	CYS	conflict	UNP A0A8B7RCM5

- Molecule 3 is a protein called SK3D Light chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	E	111	Total	C	N	O	S	0	0
			844	531	142	168	3		
3	I	111	Total	C	N	O	S	0	0
			844	531	142	168	3		

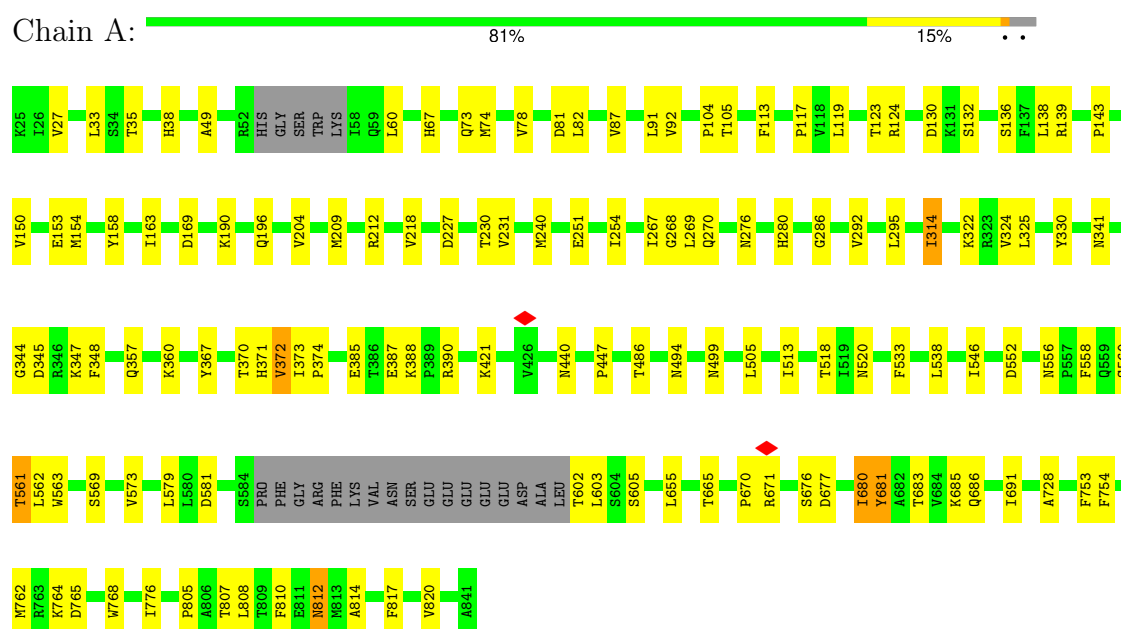
- Molecule 4 is a protein called SK3D Heavy chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	F	120	Total	C	N	O	S	0	0
			919	579	146	188	6		
4	J	120	Total	C	N	O	S	0	0
			919	579	146	188	6		

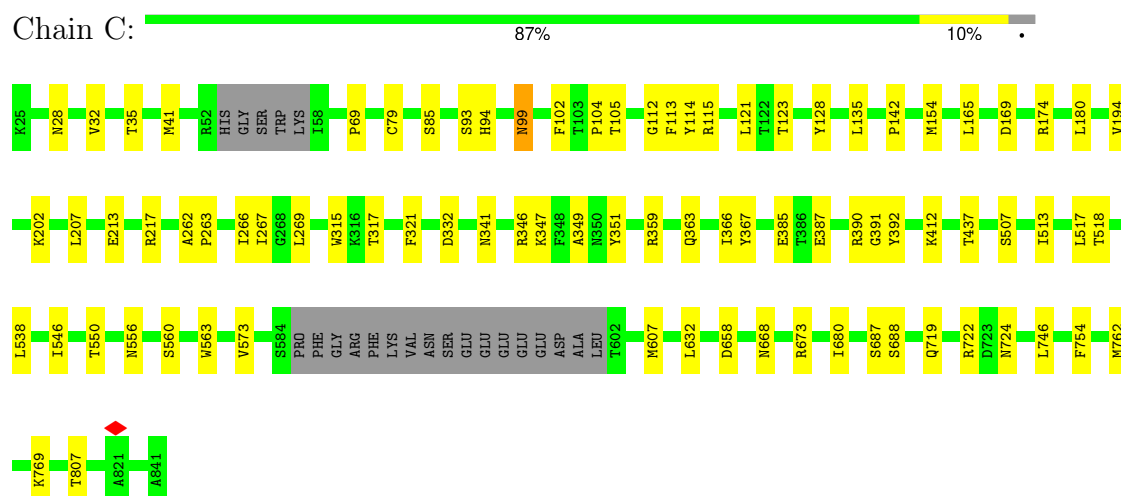
3 Residue-property plots [i](#)

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.

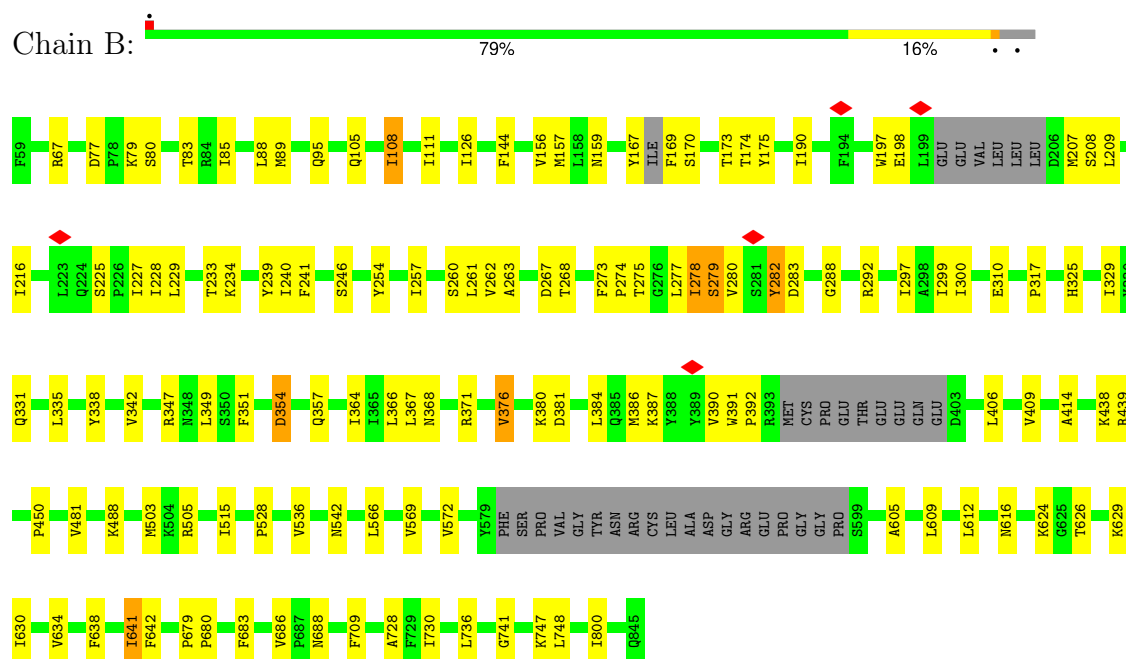
- Molecule 1: Isoform 4 of Glutamate receptor ionotropic, NMDA 1



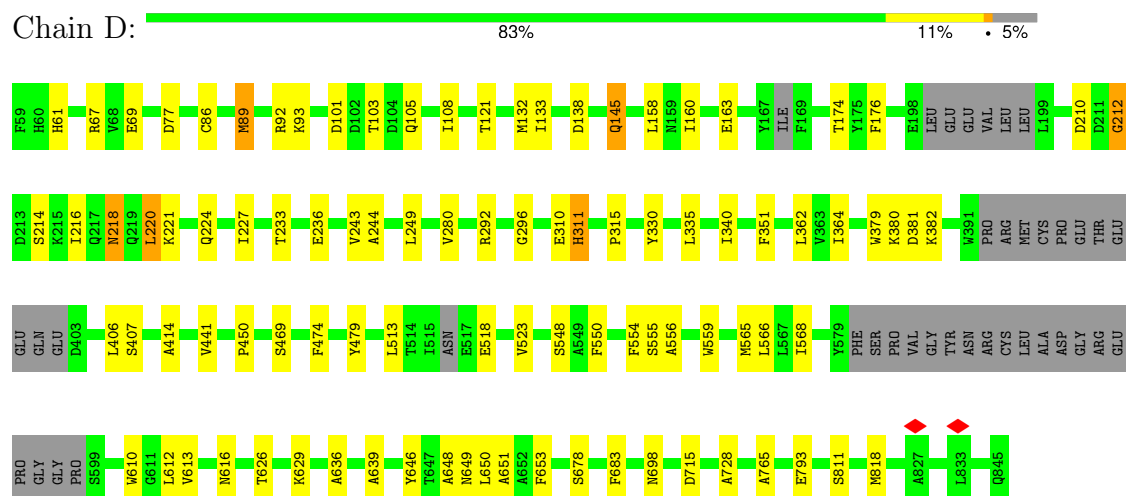
- Molecule 1: Isoform 4 of Glutamate receptor ionotropic, NMDA 1



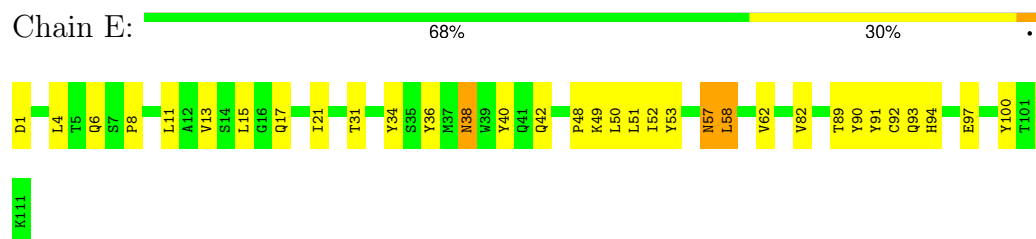
- Molecule 2: Glutamate receptor



- Molecule 2: Glutamate receptor

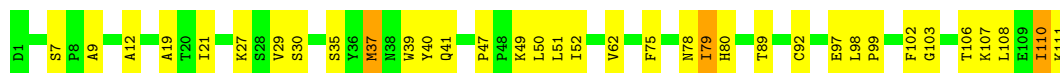


- Molecule 3: SK3D Light chain



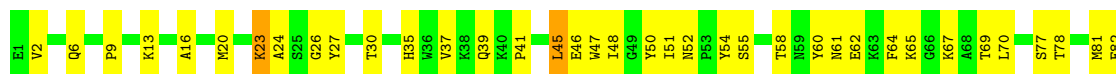
- Molecule 3: SK3D Light chain

Chain I:  68% 29% .



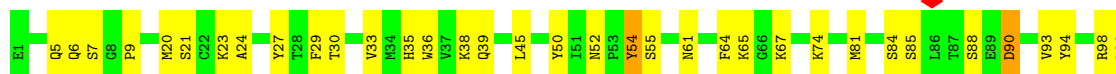
• Molecule 4: SK3D Heavy chain

Chain F:  55% 42% .



• Molecule 4: SK3D Heavy chain

Chain J:  67% 30% .



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	130398	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 ($e^-/\text{\AA}^2$)	71.7	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2200	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOCONTINUUM (6k x 4k)	Depositor
Maximum map value	0.666	Depositor
Minimum map value	-0.343	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.012	Depositor
Recommended contour level	0.036	Depositor
Map size ($\text{\AA}$)	400.0, 400.0, 400.0	wwPDB
Map dimensions	400, 400, 400	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	1.0, 1.0, 1.0	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	0/6030	0.58	0/8224
1	C	0.19	0/5836	0.49	2/7978 (0.0%)
2	B	0.23	0/5684	0.63	1/7731 (0.0%)
2	D	0.21	0/5542	0.58	3/7546 (0.0%)
3	E	0.28	0/863	0.67	0/1171
3	I	0.29	0/863	0.75	3/1171 (0.3%)
4	F	0.28	0/940	0.74	0/1274
4	J	0.31	0/940	0.81	1/1274 (0.1%)
All	All	0.23	0/26698	0.60	10/36369 (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	B	0	3
2	D	0	1
All	All	0	4

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	138	ASP	CA-C-N	6.11	132.14	122.61
2	D	138	ASP	C-N-CA	6.11	132.14	122.61
3	I	79	ILE	N-CA-C	-5.46	107.95	113.47
4	J	54	TYR	CA-CB-CG	5.31	123.47	113.90
2	B	198	GLU	N-CA-C	5.29	117.10	110.91
3	I	27	LYS	CA-CB-CG	5.18	124.46	114.10
3	I	37	MET	CB-CG-SD	5.07	127.90	112.70
2	D	89	MET	CB-CG-SD	5.04	127.83	112.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	315	TRP	CA-C-N	5.03	131.15	121.54
1	C	315	TRP	C-N-CA	5.03	131.15	121.54

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	279	SER	Peptide
2	B	380	LYS	Peptide
2	B	381	ASP	Peptide
2	D	212	GLY	Peptide

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5897	0	5623	79	0
1	C	5709	0	5256	39	0
2	B	5565	0	5202	67	0
2	D	5428	0	4955	56	0
3	E	844	0	822	23	0
3	I	844	0	824	19	0
4	F	919	0	883	36	0
4	J	919	0	883	31	0
All	All	26125	0	24448	339	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 7.

All (339) 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)
4:J:94:TYR:HB2	4:J:114:THR:HB	1.70	0.74
4:J:52:ASN:HD21	4:J:55:SER:HB2	1.57	0.69
2:B:275:THR:HG22	2:B:368:ASN:HA	1.79	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:1:ASP:N	3:E:97:GLU:OE1	2.29	0.65
4:J:30:THR:HB	4:J:54:TYR:HB2	1.79	0.64
1:A:685:LYS:HG2	1:A:686:GLN:HG3	1.80	0.64
4:F:69:THR:HB	4:F:82:GLU:HB3	1.79	0.64
1:A:440:ASN:HD21	1:A:447:PRO:HB2	1.62	0.63
3:I:7:SER:HB2	3:I:21:ILE:HD12	1.81	0.62
1:C:104:PRO:HG3	1:C:123:THR:HG21	1.81	0.62
1:C:385:GLU:HG2	1:C:387:GLU:H	1.63	0.62
3:E:38:ASN:HA	3:E:53:TYR:HA	1.81	0.61
3:I:40:TYR:HA	3:I:50:LEU:HA	1.80	0.61
1:A:73:GLN:NE2	4:J:55:SER:OG	2.28	0.60
2:B:169:PHE:N	2:B:225:SER:HG	1.98	0.60
1:A:124:ARG:HB3	1:A:143:PRO:HA	1.84	0.60
2:D:61:HIS:O	2:D:61:HIS:ND1	2.34	0.59
4:F:47:TRP:O	4:F:61:ASN:ND2	2.35	0.59
1:C:262:ALA:O	1:C:359:ARG:NH1	2.34	0.59
1:A:357:GLN:O	1:A:360:LYS:NZ	2.32	0.59
3:E:8:PRO:HB2	3:E:11:LEU:HD22	1.85	0.59
2:D:612:LEU:HA	2:D:616:ASN:HB2	1.84	0.59
4:J:38:LYS:HA	4:J:94:TYR:HA	1.83	0.59
1:A:117:PRO:HA	1:A:136:SER:HB2	1.85	0.58
2:B:299:ILE:HG23	2:B:342:VAL:HG11	1.85	0.58
4:J:20:MET:SD	4:J:21:SER:N	2.76	0.58
3:E:15:LEU:HA	3:E:82:VAL:HG23	1.86	0.58
3:I:19:ALA:HB3	3:I:79:ILE:HD13	1.85	0.58
2:D:566:LEU:HD11	2:D:610:TRP:HB2	1.86	0.58
4:F:46:GLU:OE1	4:F:61:ASN:ND2	2.37	0.58
1:C:390:ARG:NH1	1:C:392:TYR:O	2.37	0.57
4:J:35:HIS:CE1	4:J:99:LEU:HD13	2.38	0.57
1:A:295:LEU:HB2	1:A:324:VAL:HG11	1.84	0.57
4:J:6:GLN:NE2	4:J:94:TYR:O	2.37	0.57
1:A:558:PHE:HD2	1:A:562:LEU:HG	1.69	0.57
4:F:39:GLN:HB3	4:F:93:VAL:HB	1.86	0.57
4:F:67:LYS:HA	4:F:84:SER:HB3	1.87	0.57
3:I:51:LEU:HD22	3:I:62:VAL:HG22	1.88	0.56
1:A:67:HIS:HA	1:A:74:MET:HE1	1.88	0.56
2:B:67:ARG:NH2	2:B:95:GLN:OE1	2.37	0.56
3:I:97:GLU:HG3	3:I:99:PRO:HD2	1.86	0.56
4:F:48:ILE:HG23	4:F:64:PHE:HE2	1.70	0.56
2:B:439:ARG:HD3	2:B:450:PRO:HB3	1.87	0.56
1:C:762:MET:HE1	1:C:769:LYS:HA	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:F:13:LYS:HG3	4:F:16:ALA:HB2	1.86	0.56
4:J:64:PHE:HD1	4:J:67:LYS:HZ3	1.54	0.56
1:A:113:PHE:HE2	2:B:105:GLN:HG2	1.71	0.56
1:A:671:ARG:NH1	2:B:800:ILE:O	2.38	0.56
1:C:687:SER:OG	1:C:688:SER:N	2.39	0.56
1:A:762:MET:HE1	1:A:768:TRP:HB2	1.86	0.55
2:B:364:ILE:H	2:B:376:VAL:HB	1.71	0.55
3:E:40:TYR:OH	4:F:107:MET:O	2.24	0.55
3:I:78:ASN:HD21	3:I:80:HIS:HB3	1.71	0.55
2:D:683:PHE:HB3	2:D:728:ALA:HB3	1.87	0.55
3:I:39:TRP:HB2	3:I:52:ILE:HB	1.88	0.55
1:A:552:ASP:O	1:A:556:ASN:N	2.40	0.55
2:B:569:VAL:HA	2:B:572:VAL:HG12	1.89	0.54
1:A:292:VAL:HA	1:A:295:LEU:HB3	1.89	0.54
2:D:613:VAL:HG23	2:D:639:ALA:HA	1.89	0.54
2:D:649:ASN:O	2:D:653:PHE:N	2.32	0.54
1:A:807:THR:HG21	2:D:555:SER:HB3	1.90	0.54
1:A:33:LEU:HD23	1:A:38:HIS:HB3	1.90	0.54
2:D:315:PRO:HG2	2:D:335:LEU:HD22	1.89	0.54
1:A:169:ASP:OD1	1:A:196:GLN:NE2	2.41	0.54
1:A:665:THR:H	1:A:670:PRO:HD3	1.72	0.54
2:B:503:MET:SD	2:B:505:ARG:NH2	2.80	0.54
2:D:216:ILE:HG23	2:D:243:VAL:HB	1.90	0.54
4:F:64:PHE:HB2	4:F:67:LYS:HB3	1.89	0.54
1:C:112:GLY:O	1:C:115:ARG:NH1	2.42	0.53
4:F:9:PRO:HB3	4:F:115:SER:H	1.74	0.53
1:C:154:MET:HE1	1:C:267:ILE:HG21	1.89	0.53
3:E:42:GLN:O	3:E:89:THR:N	2.40	0.53
4:F:39:GLN:N	4:F:93:VAL:O	2.40	0.53
4:J:39:GLN:N	4:J:93:VAL:O	2.36	0.53
1:A:373:ILE:HD12	1:A:374:PRO:HD2	1.91	0.53
2:B:263:ALA:H	2:B:279:SER:HB2	1.74	0.53
2:D:441:VAL:HA	2:D:450:PRO:HA	1.91	0.53
2:D:554:PHE:O	2:D:559:TRP:NE1	2.42	0.53
1:A:817:PHE:HA	1:A:820:VAL:HG12	1.91	0.52
2:B:366:LEU:HD21	2:B:391:TRP:HE1	1.75	0.52
1:C:538:LEU:HD13	1:C:754:PHE:HB3	1.90	0.52
1:A:505:LEU:HB2	1:A:513:ILE:HD11	1.90	0.52
2:D:67:ARG:HH22	2:D:92:ARG:HH22	1.57	0.52
2:B:260:SER:HA	2:B:278:ILE:HB	1.92	0.52
4:J:98:ARG:NH2	4:J:108:ASP:OD1	2.43	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:154:MET:HE1	1:A:267:ILE:HG21	1.91	0.51
2:D:233:THR:HG22	2:D:236:GLU:HG3	1.91	0.51
3:E:90:TYR:HB2	3:E:106:THR:HB	1.90	0.51
1:A:251:GLU:OE2	1:A:270:GLN:NE2	2.42	0.51
1:A:655:LEU:HD21	1:A:805:PRO:HG2	1.91	0.51
1:A:676:SER:O	1:A:680:ILE:N	2.38	0.51
4:F:52:ASN:OD1	4:F:55:SER:N	2.42	0.51
2:B:566:LEU:HD11	2:B:609:LEU:HD12	1.93	0.51
4:F:51:ILE:HB	4:F:70:LEU:HD13	1.92	0.51
2:B:208:SER:OG	2:B:209:LEU:N	2.43	0.51
2:B:747:LYS:HG3	2:B:748:LEU:HD22	1.93	0.51
2:D:210:ASP:OD1	2:D:210:ASP:N	2.39	0.51
2:D:362:LEU:HB2	2:D:379:TRP:HD1	1.76	0.51
3:I:37:MET:HG2	3:I:75:PHE:CD2	2.46	0.51
1:A:367:TYR:HD1	1:A:372:VAL:HB	1.75	0.50
2:B:89:MET:SD	2:B:89:MET:N	2.84	0.50
1:C:573:VAL:HB	1:C:632:LEU:HD21	1.91	0.50
2:B:280:VAL:HG12	2:B:364:ILE:HD13	1.93	0.50
2:B:207:MET:SD	2:B:207:MET:N	2.83	0.50
1:A:486:THR:HG22	1:A:691:ILE:HD11	1.92	0.50
3:I:92:CYS:H	3:I:103:GLY:HA3	1.76	0.50
2:D:550:PHE:O	2:D:646:TYR:OH	2.26	0.50
1:A:681:TYR:HB3	1:A:728:ALA:HB3	1.94	0.50
1:A:808:LEU:HD13	2:D:649:ASN:HD21	1.77	0.50
4:F:41:PRO:HD3	4:F:92:ALA:HA	1.94	0.49
4:J:24:ALA:HB1	4:J:27:TYR:HE1	1.77	0.49
1:A:385:GLU:HG3	1:A:387:GLU:H	1.77	0.49
1:A:602:THR:HG23	1:A:605:SER:H	1.77	0.49
1:A:520:ASN:OD1	1:A:520:ASN:N	2.45	0.49
2:B:170:SER:H	2:B:228:ILE:HG22	1.77	0.49
1:C:28:ASN:ND2	1:C:85:SER:O	2.45	0.49
4:F:30:THR:HB	4:F:54:TYR:HB2	1.94	0.49
2:B:260:SER:OG	2:B:261:LEU:N	2.45	0.49
2:B:310:GLU:OE1	2:B:338:TYR:OH	2.30	0.49
2:B:630:ILE:O	2:B:634:VAL:HG23	2.13	0.49
4:F:62:GLU:HA	4:F:65:LYS:HG2	1.94	0.49
4:J:33:VAL:HG11	4:J:35:HIS:HE1	1.77	0.49
2:B:282:TYR:OH	2:B:357:GLN:OE1	2.26	0.49
3:E:52:ILE:H	3:E:52:ILE:HD12	1.78	0.49
3:I:47:PRO:O	3:I:49:LYS:NZ	2.45	0.49
2:B:85:ILE:HA	2:B:88:LEU:HD12	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:126:ILE:HD11	2:B:292:ARG:HB3	1.95	0.49
4:J:9:PRO:HA	4:J:116:VAL:HA	1.95	0.49
1:A:370:THR:OG1	1:A:371:HIS:N	2.45	0.48
2:B:262:VAL:HG21	2:B:277:LEU:HD22	1.95	0.48
2:B:273:PHE:O	2:B:371:ARG:NH1	2.35	0.48
1:C:363:GLN:HE22	1:C:366:ILE:HG13	1.78	0.48
1:C:213:GLU:OE2	1:C:213:GLU:N	2.41	0.48
4:F:6:GLN:HE22	4:F:112:GLN:HB3	1.77	0.48
1:A:139:ARG:HG3	1:A:345:ASP:HA	1.95	0.48
2:B:156:VAL:HG11	2:B:364:ILE:HG13	1.94	0.48
4:F:35:HIS:HA	4:F:50:TYR:HA	1.96	0.48
1:A:49:ALA:HB1	1:A:60:LEU:HD22	1.95	0.48
2:B:229:LEU:HB3	2:B:257:ILE:HG13	1.95	0.48
3:I:9:ALA:HA	3:I:106:THR:HA	1.94	0.48
1:A:158:TYR:HD1	1:A:388:LYS:HE2	1.79	0.48
4:F:2:VAL:HA	4:F:26:GLY:HA3	1.95	0.48
1:C:341:ASN:HD22	1:C:347:LYS:HG2	1.78	0.48
2:D:636:ALA:O	2:D:639:ALA:HB3	2.14	0.48
4:J:39:GLN:HB3	4:J:93:VAL:HB	1.95	0.48
2:B:190:ILE:HG12	2:B:197:TRP:HB3	1.95	0.47
1:A:138:LEU:HD11	1:A:325:LEU:HD13	1.96	0.47
2:B:488:LYS:NZ	2:B:688:ASN:OD1	2.41	0.47
2:B:741:GLY:O	2:B:800:ILE:N	2.45	0.47
3:E:92:CYS:O	3:E:103:GLY:N	2.47	0.47
4:F:98:ARG:HG3	4:F:108:ASP:HB3	1.96	0.47
1:A:494:ASN:OD1	1:A:494:ASN:N	2.48	0.47
2:B:626:THR:HA	2:B:629:LYS:HE2	1.97	0.47
1:C:99:ASN:OD1	1:C:99:ASN:N	2.46	0.47
4:F:20:MET:HE2	4:F:81:MET:HG3	1.96	0.47
2:B:679:PRO:HA	2:B:680:PRO:HD3	1.74	0.47
1:C:93:SER:HB3	1:C:121:LEU:HD12	1.97	0.47
1:C:437:THR:O	1:C:437:THR:OG1	2.33	0.47
2:D:548:SER:HB2	2:D:818:MET:HA	1.97	0.47
3:I:12:ALA:HB1	3:I:111:LYS:HE3	1.97	0.47
4:J:7:SER:HB2	4:J:20:MET:HE1	1.97	0.47
1:A:254:ILE:HG21	1:A:268:GLY:HA3	1.96	0.46
1:A:35:THR:HG23	1:A:38:HIS:H	1.80	0.46
1:A:602:THR:OG1	1:A:603:LEU:N	2.48	0.46
2:B:79:LYS:O	2:B:83:THR:OG1	2.28	0.46
1:C:722:ARG:HG2	1:C:746:LEU:HD11	1.98	0.46
2:D:381:ASP:HA	2:D:382:LYS:HA	1.55	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:523:VAL:HG11	2:D:765:ALA:HB1	1.96	0.46
2:D:612:LEU:HD21	2:D:636:ALA:HB2	1.98	0.46
2:D:793:GLU:OE2	2:D:793:GLU:N	2.42	0.46
1:A:209:MET:HA	1:A:212:ARG:HB2	1.98	0.46
4:F:55:SER:O	4:F:55:SER:OG	2.33	0.46
4:J:84:SER:OG	4:J:85:SER:N	2.49	0.46
1:A:286:GLY:HA3	1:A:330:TYR:CZ	2.50	0.46
2:D:362:LEU:HB2	2:D:379:TRP:CD1	2.51	0.46
3:E:34:TYR:HB3	3:E:36:TYR:CE1	2.51	0.46
1:A:499:ASN:OD1	1:A:499:ASN:N	2.46	0.46
2:D:244:ALA:HB1	2:D:249:LEU:HB2	1.97	0.46
3:I:98:LEU:HD11	4:J:50:TYR:HE2	1.81	0.46
2:B:536:VAL:HG12	2:B:730:ILE:HG12	1.98	0.46
1:A:561:THR:HG23	1:A:562:LEU:HD22	1.98	0.45
1:A:677:ASP:OD1	1:A:677:ASP:N	2.48	0.45
2:D:469:SER:OG	2:D:474:PHE:O	2.33	0.45
3:E:91:TYR:HE1	4:F:45:LEU:H	1.65	0.45
1:A:808:LEU:HD23	2:D:554:PHE:CD1	2.51	0.45
2:B:240:ILE:HG22	2:B:241:PHE:HD2	1.81	0.45
1:C:263:PRO:HD2	1:C:266:ILE:HD11	1.97	0.45
3:E:94:HIS:H	3:E:94:HIS:HD1	1.64	0.45
1:A:560:SER:HA	1:A:563:TRP:HD1	1.80	0.45
1:C:169:ASP:O	1:C:174:ARG:NH1	2.46	0.45
2:D:67:ARG:NH2	2:D:69:GLU:OE2	2.50	0.45
4:J:29:PHE:HE1	4:J:74:LYS:HA	1.82	0.45
1:C:69:PRO:HA	1:C:102:PHE:HE1	1.81	0.45
1:C:332:ASP:OD1	1:C:332:ASP:N	2.50	0.45
2:D:224:GLN:N	2:D:224:GLN:OE1	2.49	0.45
2:D:650:LEU:HA	2:D:653:PHE:HB2	1.99	0.45
3:I:51:LEU:HD13	3:I:62:VAL:HG13	1.99	0.45
4:J:6:GLN:NE2	4:J:114:THR:OG1	2.50	0.45
1:C:668:ASN:HA	1:C:673:ARG:HH21	1.82	0.45
1:C:658:ASP:OD1	1:C:658:ASP:N	2.49	0.45
4:J:45:LEU:HD12	4:J:45:LEU:H	1.82	0.45
1:A:163:ILE:HG22	1:A:218:VAL:HB	1.98	0.45
2:B:638:PHE:HA	2:B:641:ILE:HG22	1.99	0.45
1:C:724:ASN:OD1	1:C:724:ASN:N	2.50	0.45
4:J:33:VAL:CG1	4:J:35:HIS:CE1	3.00	0.45
4:J:61:ASN:O	4:J:65:LYS:N	2.50	0.45
2:B:354:ASP:OD1	2:B:354:ASP:N	2.50	0.45
2:B:515:ILE:HD12	2:B:528:PRO:HG3	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:49:LYS:HB3	3:E:49:LYS:HE3	1.71	0.45
1:A:808:LEU:HB3	1:A:812:ASN:HD21	1.82	0.44
2:D:89:MET:HE2	2:D:89:MET:HA	1.97	0.44
2:D:556:ALA:HA	2:D:559:TRP:HD1	1.81	0.44
3:E:52:ILE:HA	3:E:58:LEU:HA	1.99	0.44
2:B:282:TYR:HD1	2:B:283:ASP:H	1.64	0.44
2:D:174:THR:HG23	2:D:176:PHE:H	1.82	0.44
3:E:51:LEU:HA	3:E:62:VAL:HG21	1.99	0.44
3:I:37:MET:HE2	3:I:37:MET:HA	1.97	0.44
1:A:579:LEU:O	1:A:581:ASP:N	2.50	0.44
3:E:8:PRO:HG3	3:E:21:ILE:HA	1.99	0.44
4:F:48:ILE:O	4:F:61:ASN:N	2.51	0.44
1:C:142:PRO:HG2	1:C:349:ALA:HB2	1.99	0.44
2:D:132:MET:O	2:D:145:GLN:NE2	2.51	0.44
4:J:36:TRP:NE1	4:J:81:MET:HB2	2.33	0.44
1:A:808:LEU:HD13	2:D:649:ASN:ND2	2.33	0.44
1:A:533:PHE:HB3	1:A:776:ILE:HD11	1.98	0.44
1:A:538:LEU:HD13	1:A:754:PHE:HB3	1.99	0.44
2:B:683:PHE:HB3	2:B:728:ALA:HB3	1.98	0.44
1:C:550:THR:O	1:C:556:ASN:ND2	2.50	0.44
2:D:218:ASN:H	2:D:221:LYS:HE3	1.82	0.44
4:F:58:THR:HG21	4:F:60:TYR:CZ	2.53	0.44
2:D:160:ILE:HD13	2:D:160:ILE:HA	1.85	0.44
3:I:41:GLN:O	3:I:49:LYS:N	2.48	0.44
2:B:159:ASN:HB3	2:B:384:LEU:HD21	2.00	0.43
2:B:616:ASN:OD1	2:B:616:ASN:N	2.49	0.43
1:C:513:ILE:HG21	1:C:517:LEU:HD22	1.99	0.43
4:F:87:THR:HA	4:F:118:VAL:HG11	1.98	0.43
1:A:74:MET:O	1:A:78:VAL:HG23	2.18	0.43
2:D:310:GLU:OE1	2:D:311:HIS:ND1	2.51	0.43
3:I:89:THR:HB	3:I:107:LYS:HG3	2.00	0.43
2:D:715:ASP:N	2:D:715:ASP:OD1	2.51	0.43
4:J:90:ASP:N	4:J:90:ASP:OD1	2.51	0.43
4:F:9:PRO:HB3	4:F:115:SER:N	2.34	0.43
4:F:37:VAL:HG11	4:F:110:TRP:HZ3	1.82	0.43
1:A:153:GLU:OE2	1:A:153:GLU:N	2.52	0.43
2:B:297:ILE:HA	2:B:300:ILE:HD12	2.01	0.43
2:D:626:THR:HA	2:D:629:LYS:HE2	2.01	0.43
3:E:6:GLN:HA	3:E:6:GLN:NE2	2.33	0.43
4:F:35:HIS:HB3	4:F:47:TRP:HE1	1.83	0.43
2:B:108:ILE:H	2:B:108:ILE:HG13	1.63	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:648:ALA:O	2:D:651:ALA:HB3	2.19	0.43
4:F:2:VAL:HG22	4:F:27:TYR:HB3	2.01	0.43
4:F:51:ILE:HD12	4:F:51:ILE:HA	1.87	0.43
1:A:680:ILE:HG22	1:A:681:TYR:H	1.84	0.42
4:J:6:GLN:HB3	4:J:112:GLN:HB2	2.01	0.42
4:J:88:SER:HA	4:J:118:VAL:HB	2.01	0.42
1:A:212:ARG:HG2	1:A:240:MET:HE3	2.01	0.42
2:D:105:GLN:HB2	2:D:108:ILE:HG13	2.01	0.42
4:F:24:ALA:HB3	4:F:77:SER:HB3	2.01	0.42
1:A:138:LEU:HD22	1:A:344:GLY:HA3	2.01	0.42
1:A:158:TYR:O	1:A:388:LYS:NZ	2.46	0.42
1:A:276:ASN:O	1:A:280:HIS:ND1	2.52	0.42
1:A:810:PHE:HD2	1:A:814:ALA:HB2	1.83	0.42
2:B:233:THR:OG1	2:B:234:LYS:N	2.51	0.42
2:B:542:ASN:OD1	2:B:542:ASN:N	2.53	0.42
3:E:13:VAL:HG13	3:E:17:GLN:HB2	2.01	0.42
2:B:241:PHE:HE1	2:B:274:PRO:HD3	1.84	0.42
2:D:280:VAL:HG12	2:D:364:ILE:HG12	2.02	0.42
4:J:33:VAL:HG12	4:J:99:LEU:HD22	2.01	0.42
2:B:157:MET:H	2:B:157:MET:HG2	1.65	0.42
2:B:347:ARG:O	2:B:349:LEU:N	2.52	0.42
2:D:77:ASP:OD1	2:D:77:ASP:N	2.51	0.42
1:C:165:LEU:HD11	1:C:180:LEU:HD23	2.01	0.42
2:D:220:LEU:HD23	2:D:220:LEU:HA	1.89	0.42
1:A:764:LYS:HE2	1:A:764:LYS:HB3	1.83	0.42
2:D:565:MET:HA	2:D:568:ILE:HG12	2.01	0.42
2:B:144:PHE:HB3	2:B:351:PHE:CD2	2.55	0.42
2:B:144:PHE:HB3	2:B:351:PHE:CE2	2.55	0.42
3:I:30:SER:HB3	3:I:35:SER:HB3	2.02	0.42
1:C:94:HIS:HB2	1:C:123:THR:HG22	2.02	0.42
2:D:133:ILE:HG22	2:D:145:GLN:HG3	2.02	0.42
3:E:93:GLN:HE22	3:E:100:TYR:HB3	1.85	0.42
4:J:33:VAL:HG11	4:J:35:HIS:CE1	2.55	0.42
2:B:167:TYR:HE2	2:B:254:TYR:HB3	1.85	0.42
1:C:202:LYS:HA	1:C:202:LYS:HD3	1.92	0.42
1:A:227:ASP:O	1:A:230:THR:OG1	2.33	0.41
1:A:314:ILE:HD12	1:A:314:ILE:HA	1.89	0.41
2:B:77:ASP:O	2:B:80:SER:OG	2.33	0.41
2:B:174:THR:HB	2:B:175:TYR:H	1.71	0.41
1:C:351:TYR:HD1	1:C:367:TYR:HD2	1.68	0.41
2:D:380:LYS:HE3	2:D:380:LYS:HB3	1.62	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:J:23:LYS:HB3	4:J:23:LYS:HE3	1.73	0.41
1:A:91:LEU:HB2	1:A:119:LEU:HB3	2.02	0.41
1:A:254:ILE:HD13	1:A:268:GLY:HA3	2.02	0.41
1:A:388:LYS:O	1:A:390:ARG:NH1	2.54	0.41
1:C:412:LYS:HA	1:C:412:LYS:HD3	1.88	0.41
2:D:212:GLY:O	2:D:214:SER:N	2.45	0.41
3:E:57:ASN:OD1	3:E:57:ASN:N	2.52	0.41
1:A:136:SER:HA	1:A:322:LYS:NZ	2.35	0.41
1:A:204:VAL:HG21	1:A:231:VAL:HA	2.02	0.41
1:C:105:THR:HG22	1:C:128:TYR:HE1	1.86	0.41
1:C:142:PRO:HD3	1:C:346:ARG:HD2	2.01	0.41
2:B:641:ILE:HD12	2:B:641:ILE:HA	1.80	0.41
1:A:104:PRO:HD3	1:A:123:THR:HG21	2.03	0.41
1:A:130:ASP:OD2	1:A:132:SER:OG	2.37	0.41
1:A:560:SER:HA	1:A:563:TRP:CD1	2.56	0.41
1:A:81:ASP:OD2	1:A:82:LEU:N	2.53	0.41
1:A:341:ASN:HB3	1:A:347:LYS:HB2	2.02	0.41
2:B:95:GLN:HE21	2:B:317:PRO:HD3	1.86	0.41
2:B:605:ALA:O	2:B:609:LEU:HB2	2.21	0.41
2:D:407:SER:HB3	2:D:479:TYR:HE1	1.86	0.41
2:D:513:LEU:HD23	2:D:513:LEU:HA	1.94	0.41
4:F:23:LYS:HA	4:F:78:THR:HA	2.02	0.41
1:A:190:LYS:HA	1:A:190:LYS:HD2	1.78	0.41
1:A:569:SER:O	1:A:573:VAL:HG22	2.21	0.41
2:B:297:ILE:H	2:B:297:ILE:HG12	1.70	0.41
1:C:560:SER:HA	1:C:563:TRP:HD1	1.86	0.41
2:D:132:MET:H	2:D:132:MET:HG2	1.65	0.41
4:F:67:LYS:HE3	4:F:83:LEU:HD13	2.01	0.41
4:F:97:ALA:HB1	4:F:107:MET:SD	2.60	0.41
2:B:156:VAL:HG21	2:B:364:ILE:HG13	2.03	0.41
2:B:624:LYS:HA	2:B:624:LYS:HD2	1.97	0.41
2:B:288:GLY:O	2:B:292:ARG:HG3	2.21	0.40
1:C:217:ARG:HB2	1:C:391:GLY:HA2	2.03	0.40
3:E:93:GLN:NE2	3:E:100:TYR:HB3	2.36	0.40
2:B:331:GLN:O	2:B:335:LEU:HD12	2.22	0.40
2:D:698:ASN:OD1	2:D:698:ASN:N	2.53	0.40
3:E:48:PRO:HD3	4:F:95:TYR:CE1	2.56	0.40
3:E:50:LEU:H	3:E:50:LEU:HD23	1.85	0.40
4:J:5:GLN:HB3	4:J:23:LYS:HB3	2.03	0.40
1:A:150:VAL:HG11	1:A:269:LEU:HG	2.03	0.40
1:A:810:PHE:CD2	1:A:814:ALA:HB2	2.57	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:41:MET:HE3	1:C:41:MET:HB2	1.87	0.40
1:C:317:THR:O	1:C:321:PHE:N	2.47	0.40
2:D:340:ILE:O	2:D:351:PHE:N	2.54	0.40
2:B:277:LEU:HD13	2:B:367:LEU:HD23	2.04	0.40
2:B:387:LYS:HA	2:B:392:PRO:HD3	2.03	0.40
2:D:93:LYS:HA	2:D:93:LYS:HD3	1.90	0.40
2:D:292:ARG:O	2:D:296:GLY:N	2.52	0.40
3:I:108:LEU:HD11	3:I:110:ILE:HB	2.04	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	789/817 (97%)	735 (93%)	50 (6%)	4 (0%)	24	56
1	C	789/817 (97%)	749 (95%)	39 (5%)	1 (0%)	48	78
2	B	742/787 (94%)	653 (88%)	84 (11%)	5 (1%)	18	50
2	D	737/787 (94%)	659 (89%)	71 (10%)	7 (1%)	14	45
3	E	109/111 (98%)	87 (80%)	22 (20%)	0	100	100
3	I	109/111 (98%)	99 (91%)	10 (9%)	0	100	100
4	F	118/120 (98%)	101 (86%)	16 (14%)	1 (1%)	16	48
4	J	118/120 (98%)	106 (90%)	11 (9%)	1 (1%)	16	48
All	All	3511/3670 (96%)	3189 (91%)	303 (9%)	19 (0%)	26	56

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	680	ILE
1	C	680	ILE

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Mol	Chain	Res	Type
1	A	683	THR
2	B	246	SER
2	D	678	SER
2	B	414	ALA
2	D	227	ILE
2	D	414	ALA
1	A	348	PHE
2	D	163	GLU
2	B	227	ILE
2	B	386	MET
2	B	390	VAL
2	D	218	ASN
2	D	518	GLU
2	D	811	SER
4	F	116	VAL
4	J	116	VAL
1	A	27	VAL

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	586/704 (83%)	573 (98%)	13 (2%)	45	63
1	C	535/704 (76%)	519 (97%)	16 (3%)	36	57
2	B	550/689 (80%)	527 (96%)	23 (4%)	26	51
2	D	518/689 (75%)	508 (98%)	10 (2%)	50	65
3	E	92/92 (100%)	86 (94%)	6 (6%)	15	43
3	I	92/92 (100%)	89 (97%)	3 (3%)	33	55
4	F	102/102 (100%)	97 (95%)	5 (5%)	22	48
4	J	102/102 (100%)	100 (98%)	2 (2%)	48	64
All	All	2577/3174 (81%)	2499 (97%)	78 (3%)	37	57

All (78) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	87	VAL
1	A	92	VAL
1	A	105	THR
1	A	314	ILE
1	A	372	VAL
1	A	421	LYS
1	A	518	THR
1	A	546	ILE
1	A	561	THR
1	A	681	TYR
1	A	753	PHE
1	A	765	ASP
1	A	812	ASN
2	B	108	ILE
2	B	111	ILE
2	B	173	THR
2	B	216	ILE
2	B	239	TYR
2	B	267	ASP
2	B	268	THR
2	B	278	ILE
2	B	282	TYR
2	B	325	HIS
2	B	329	ILE
2	B	354	ASP
2	B	376	VAL
2	B	406	LEU
2	B	409	VAL
2	B	438	LYS
2	B	481	VAL
2	B	612	LEU
2	B	641	ILE
2	B	642	PHE
2	B	686	VAL
2	B	709	PHE
2	B	736	LEU
1	C	32	VAL
1	C	35	THR
1	C	79	CYS
1	C	99	ASN
1	C	113	PHE
1	C	114	TYR
1	C	135	LEU

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Mol	Chain	Res	Type
1	C	194	VAL
1	C	207	LEU
1	C	269	LEU
1	C	507	SER
1	C	518	THR
1	C	546	ILE
1	C	607	MET
1	C	719	GLN
1	C	807	THR
2	D	86	CYS
2	D	101	ASP
2	D	103	THR
2	D	121	THR
2	D	145	GLN
2	D	158	LEU
2	D	220	LEU
2	D	311	HIS
2	D	330	TYR
2	D	406	LEU
3	E	4	LEU
3	E	31	THR
3	E	38	ASN
3	E	57	ASN
3	E	58	LEU
3	E	102	PHE
4	F	23	LYS
4	F	45	LEU
4	F	103	VAL
4	F	110	TRP
4	F	114	THR
3	I	29	VAL
3	I	102	PHE
3	I	110	ILE
4	J	90	ASP
4	J	118	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	368	ASN
1	A	371	HIS
1	A	616	ASN

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Continued from previous page...

Mol	Chain	Res	Type
1	A	742	GLN
2	B	224	GLN
2	B	812	GLN
1	C	28	ASN
1	C	70	ASN
1	C	171	HIS
1	C	355	ASN
1	C	371	HIS
1	C	405	GLN
1	C	440	ASN
1	C	453	GLN
2	D	336	ASN
2	D	649	ASN
2	D	817	ASN
3	E	80	HIS
4	F	59	ASN
4	F	61	ASN
4	F	112	GLN
3	I	46	GLN
4	J	3	GLN
4	J	35	HIS

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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
2	D	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	D	199:LEU	C	206:ASP	N	20.18

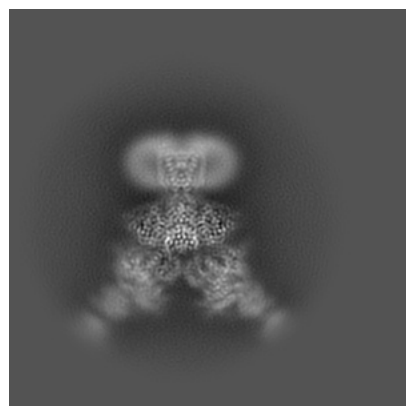
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-75112. These allow visual inspection of the internal detail of the map and identification of artifacts.

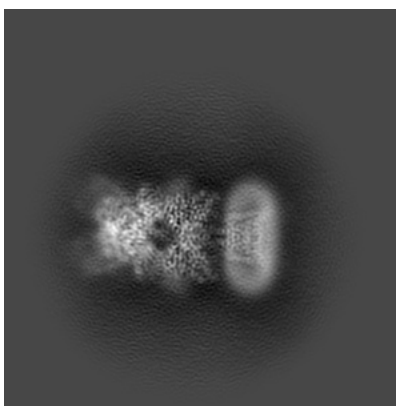
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

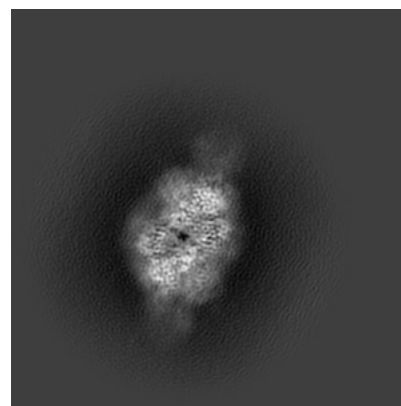
6.1.1 Primary map



X

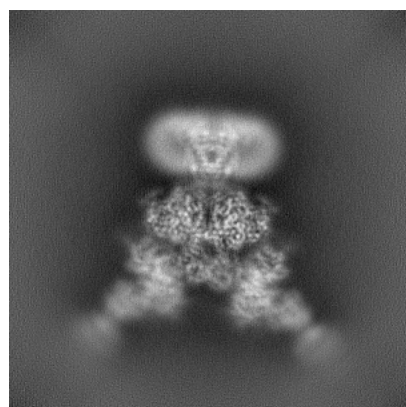


Y

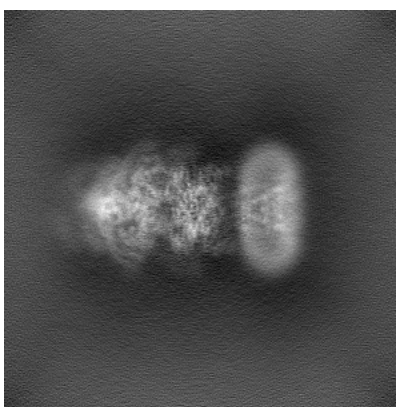


Z

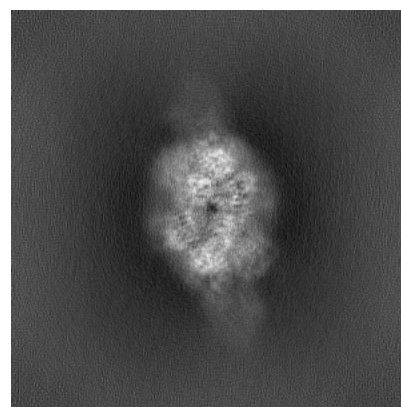
6.1.2 Raw map



X



Y

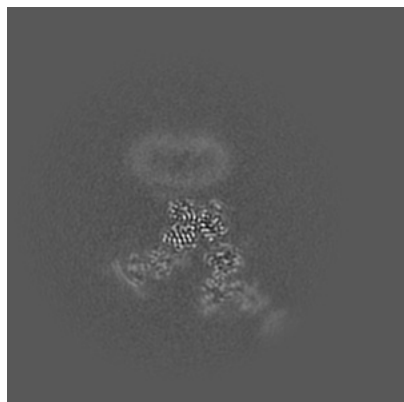


Z

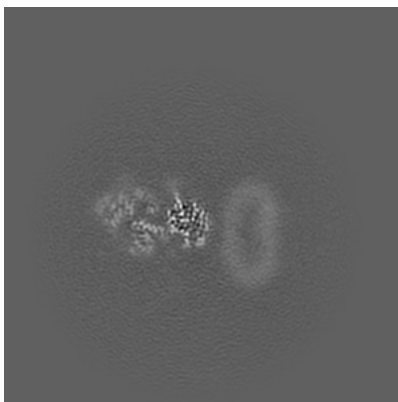
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

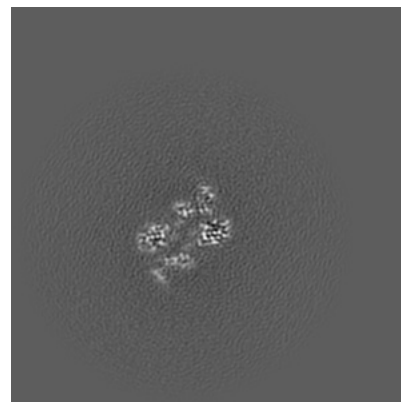
6.2.1 Primary map



X Index: 200

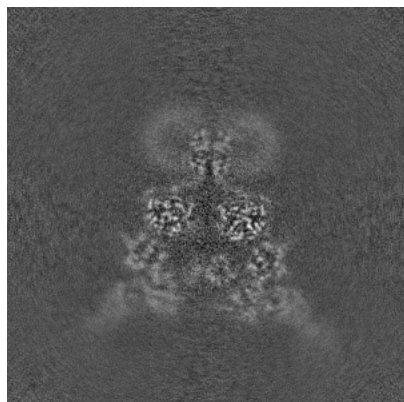


Y Index: 200

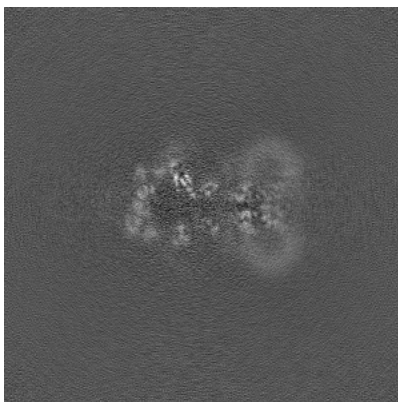


Z Index: 200

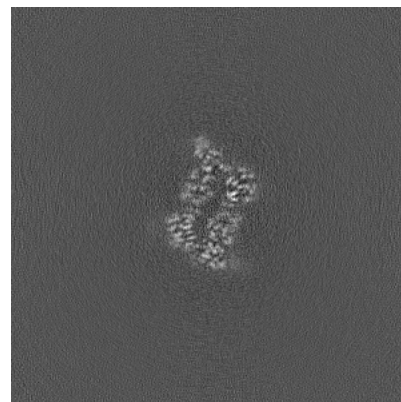
6.2.2 Raw map



X Index: 200



Y Index: 200

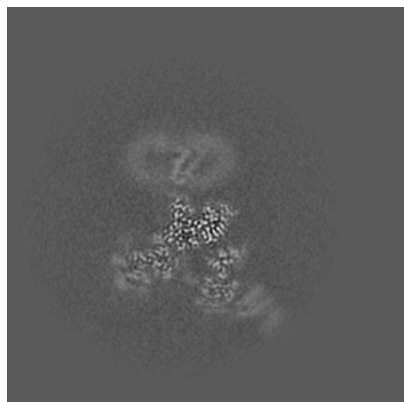


Z Index: 200

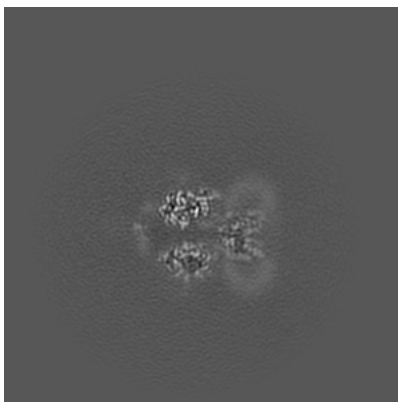
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

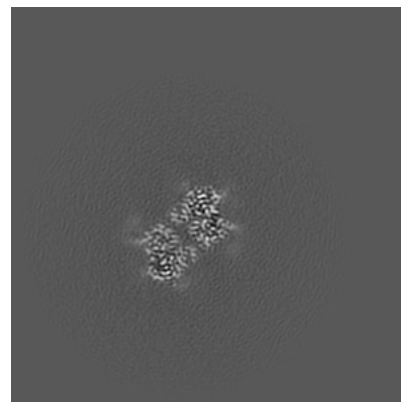
6.3.1 Primary map



X Index: 193

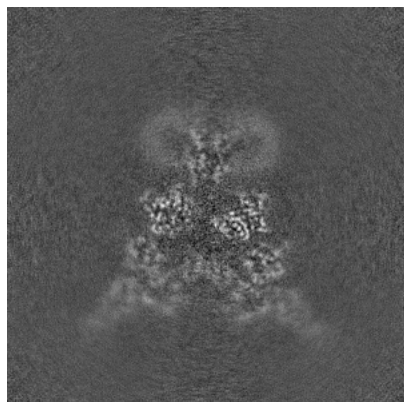


Y Index: 174

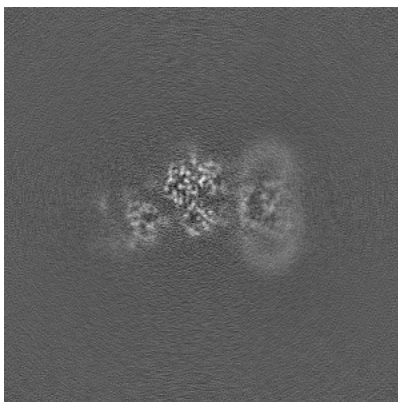


Z Index: 177

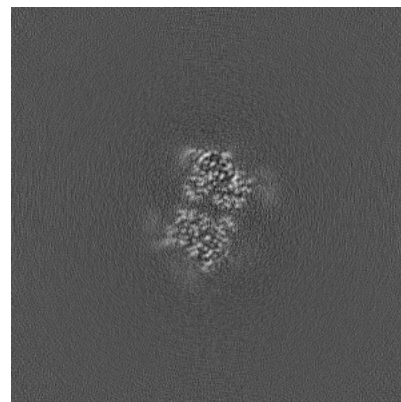
6.3.2 Raw map



X Index: 204



Y Index: 219

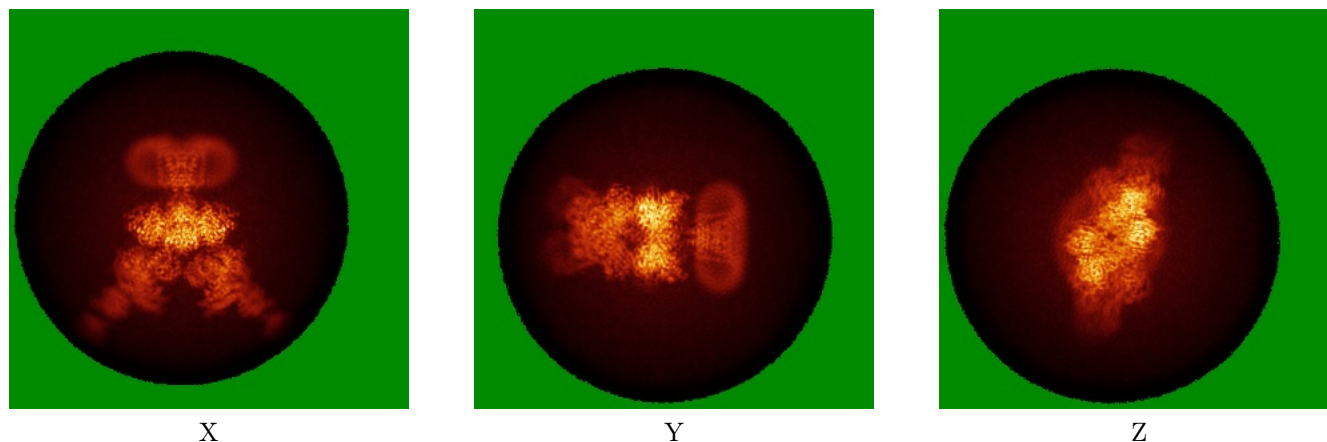


Z Index: 185

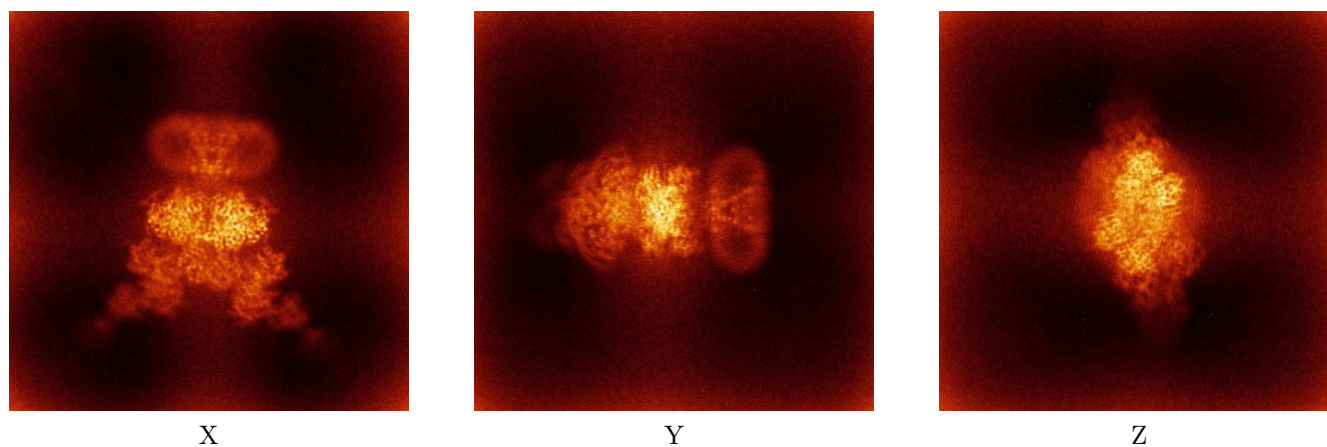
The images above show the largest variance slices of the map in three orthogonal directions.

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

6.4.1 Primary map



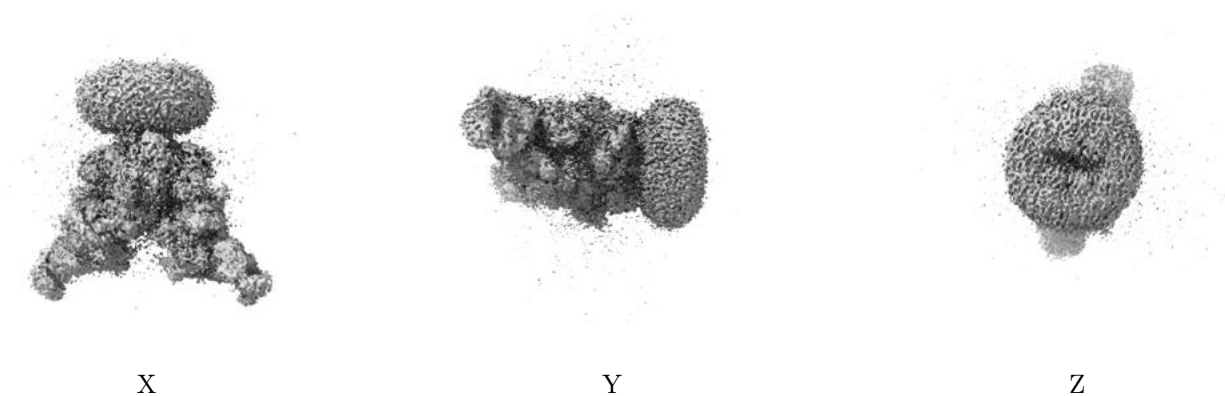
6.4.2 Raw map



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

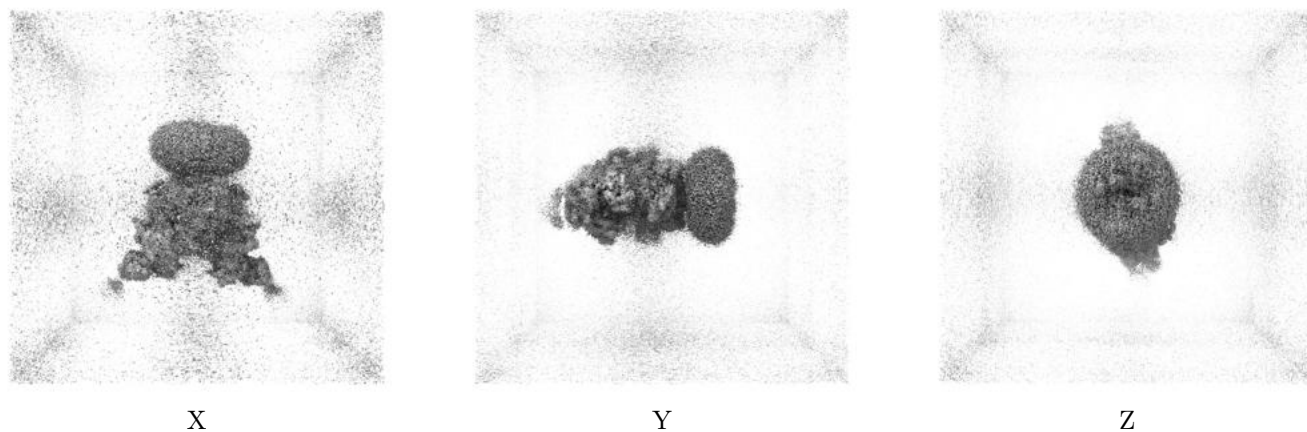
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.036. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



These images show the 3D surface of the raw map. The raw map's contour level was selected so that its surface encloses the same volume as the primary map does at its recommended contour level.

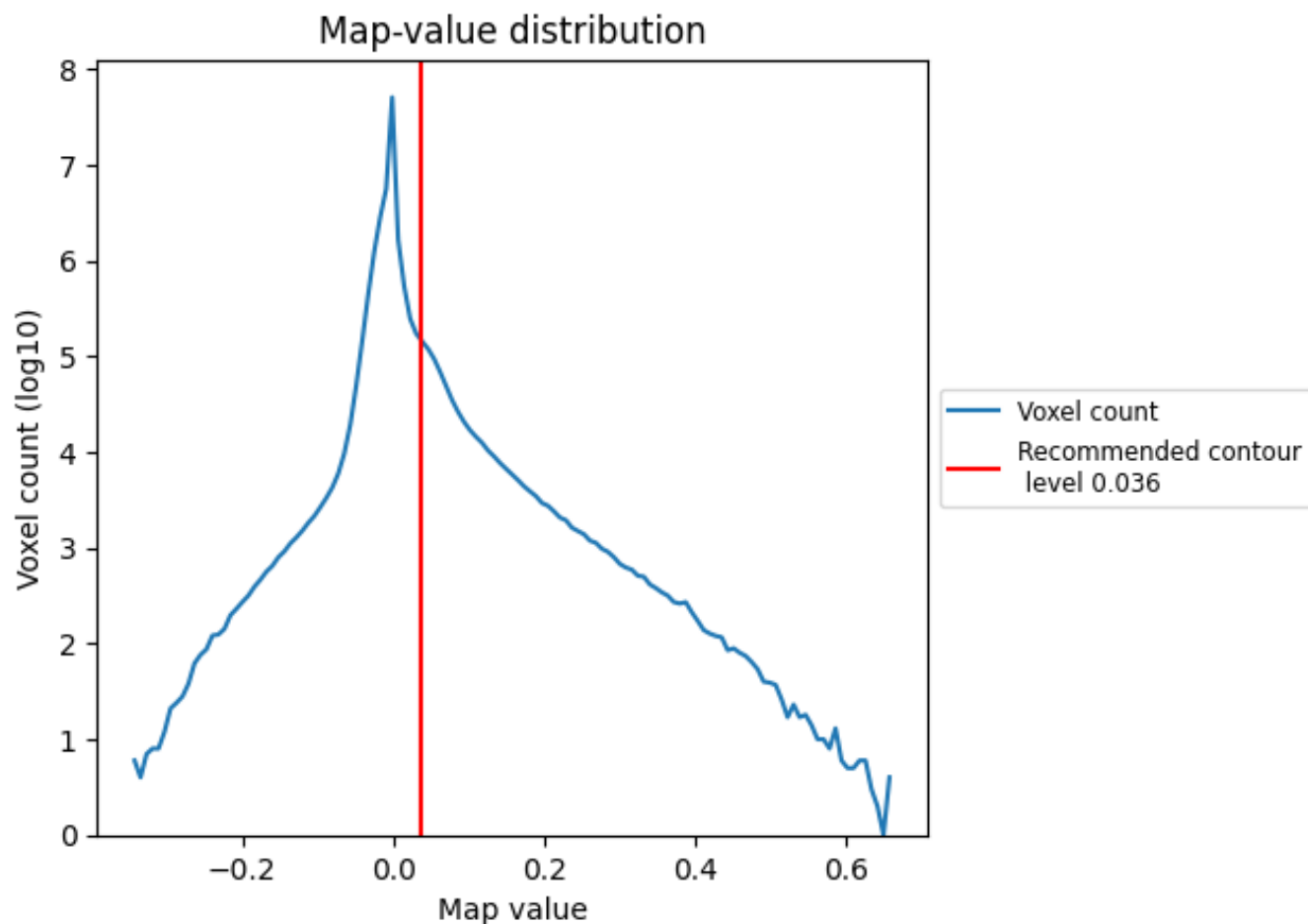
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

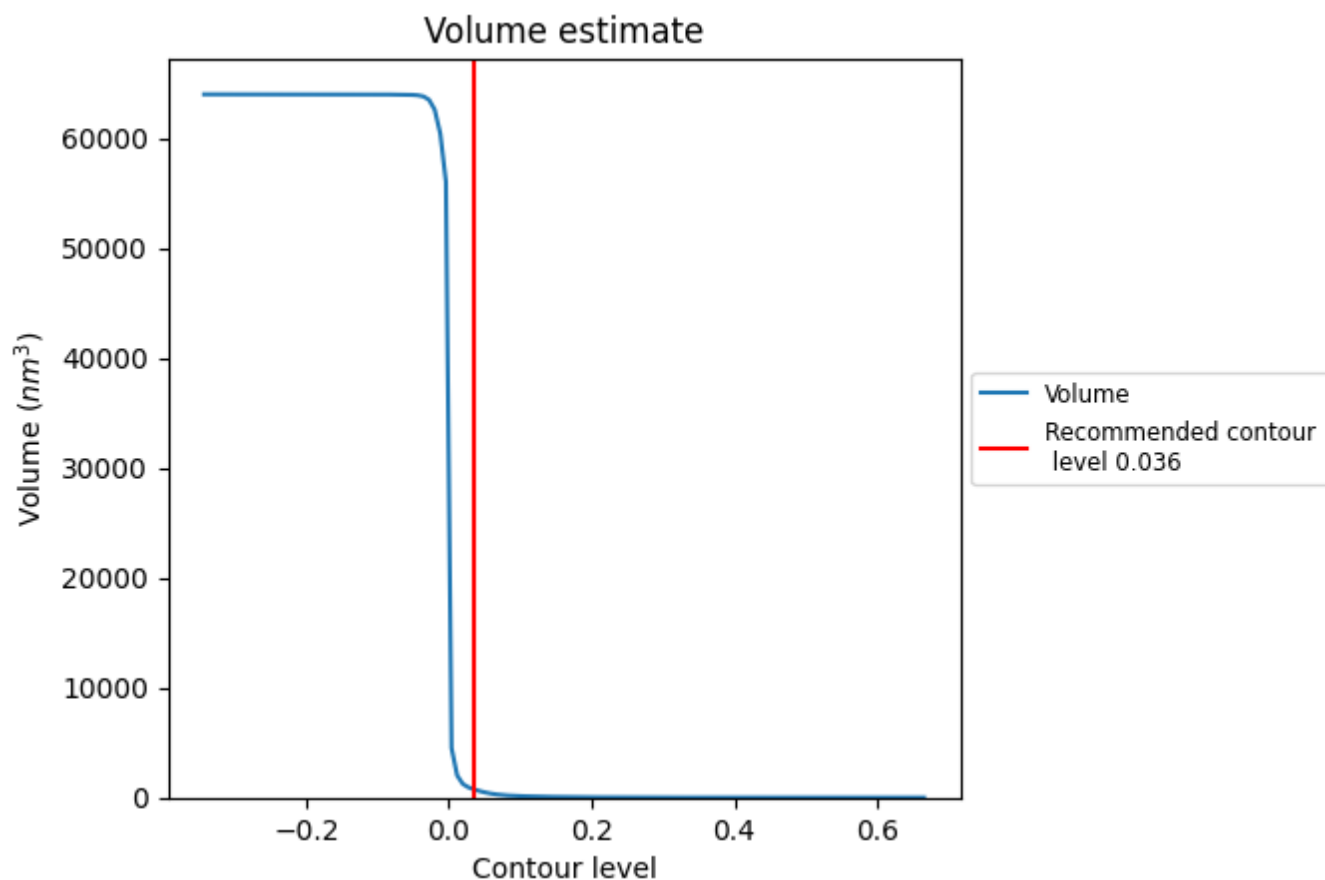
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



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

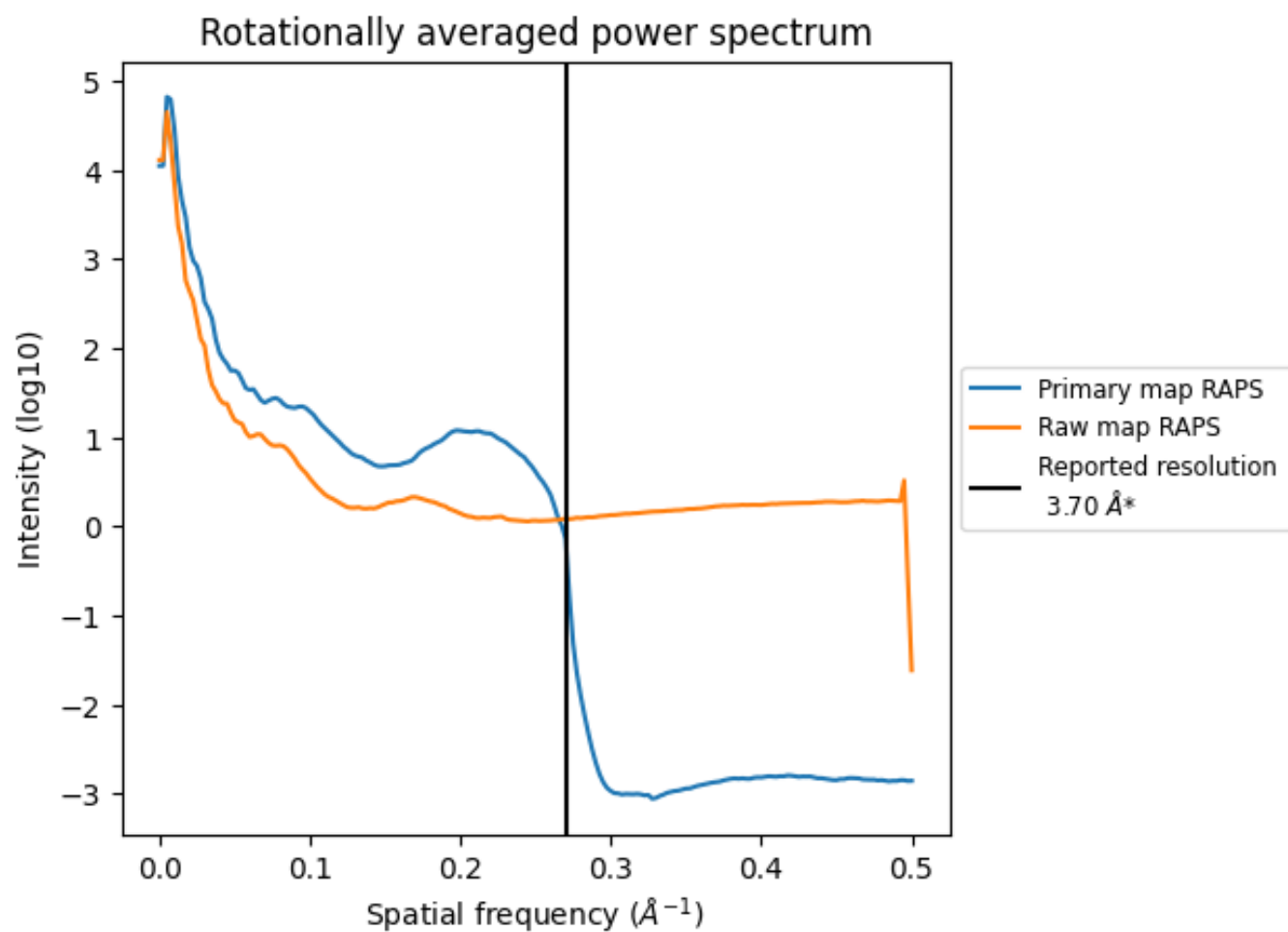
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 727 nm^3 ; this corresponds to an approximate mass of 657 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 ⓘ

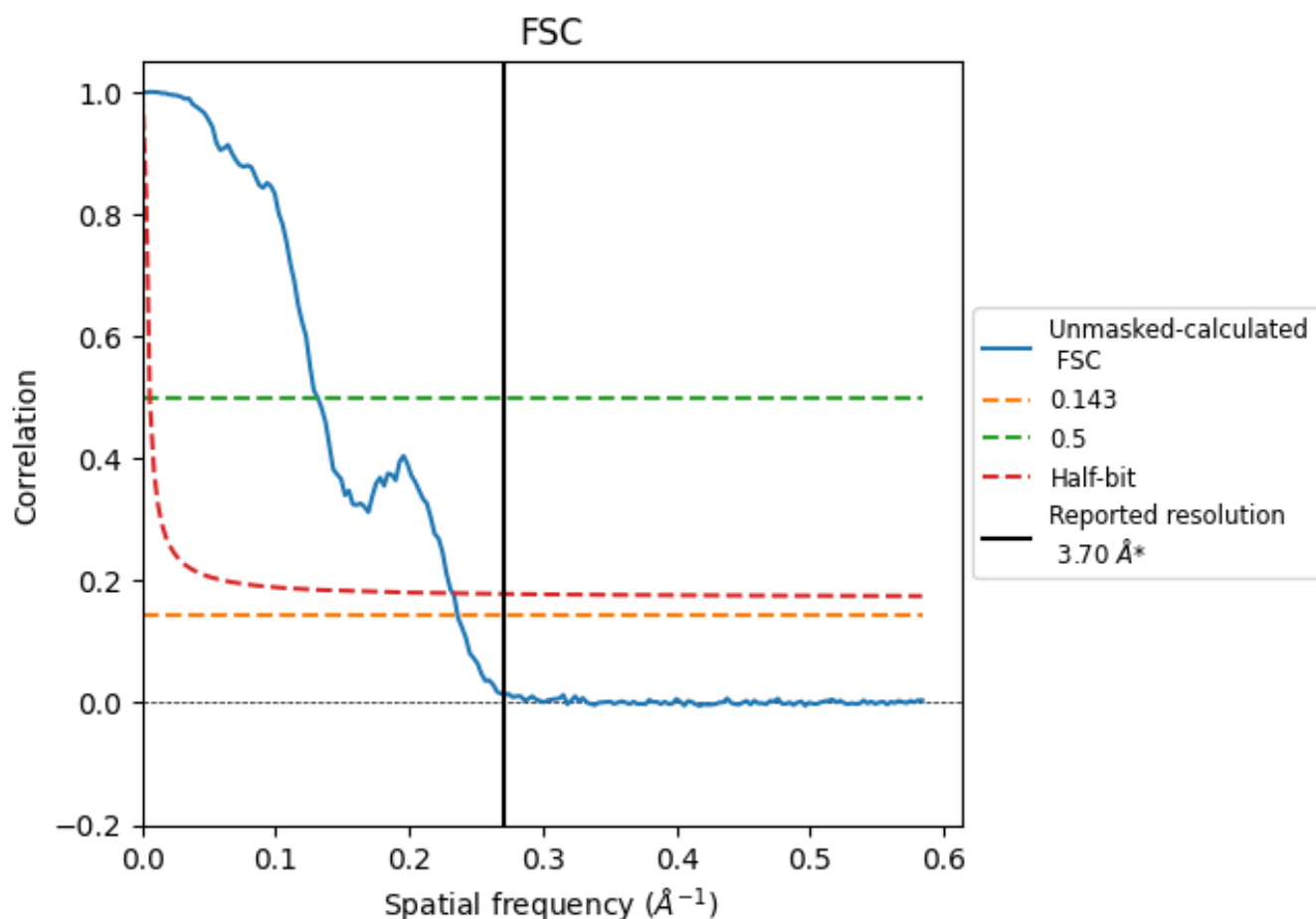


*Reported resolution corresponds to spatial frequency of $0.270 \text{ \AA}^{-1}$

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 $0.270 \text{ \AA}^{-1}$

8.2 Resolution estimates [i](#)

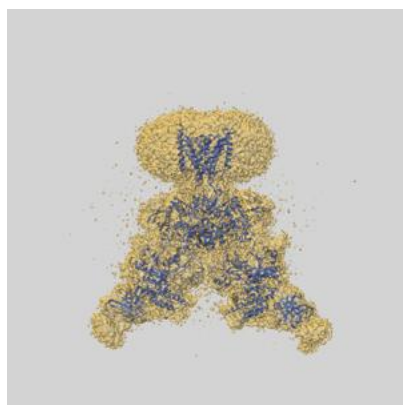
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.70	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	4.23	7.62	4.30

*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 4.23 differs from the reported value 3.7 by more than 10 %

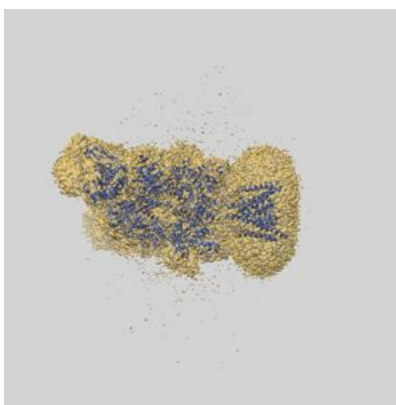
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-75112 and PDB model 10EN. Per-residue inclusion information can be found in section [3](#) on page [6](#).

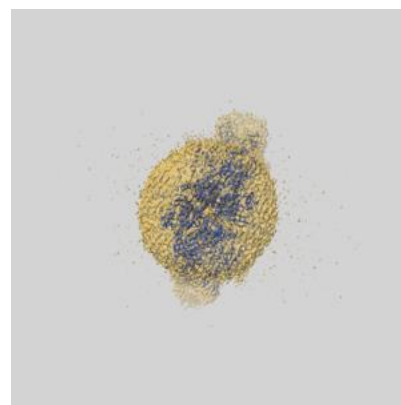
9.1 Map-model overlay [i](#)



X



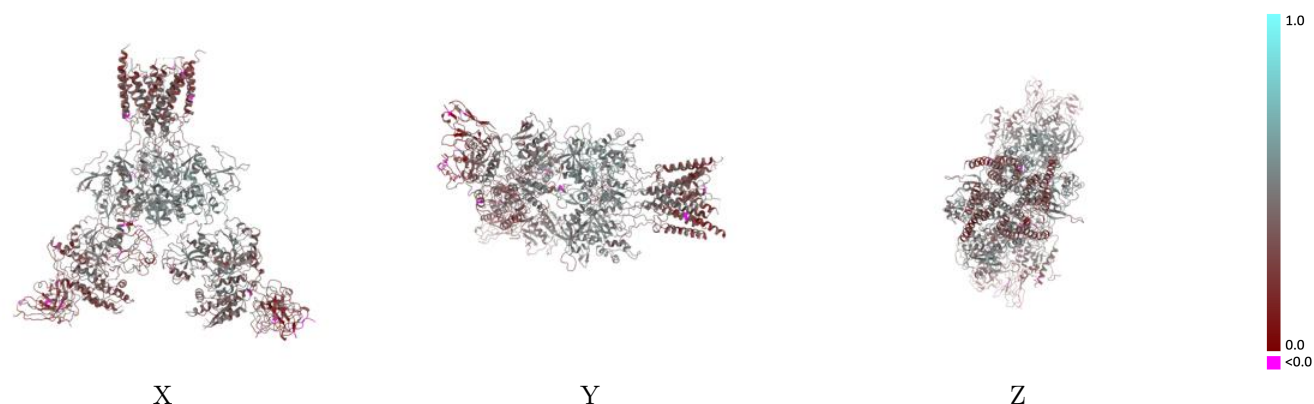
Y



Z

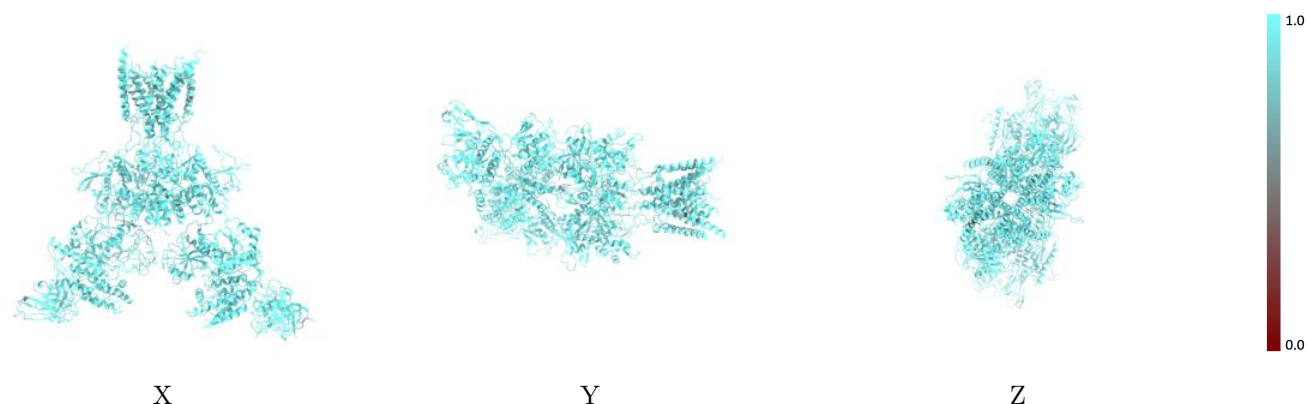
The images above show the 3D surface view of the map at the recommended contour level 0.036 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



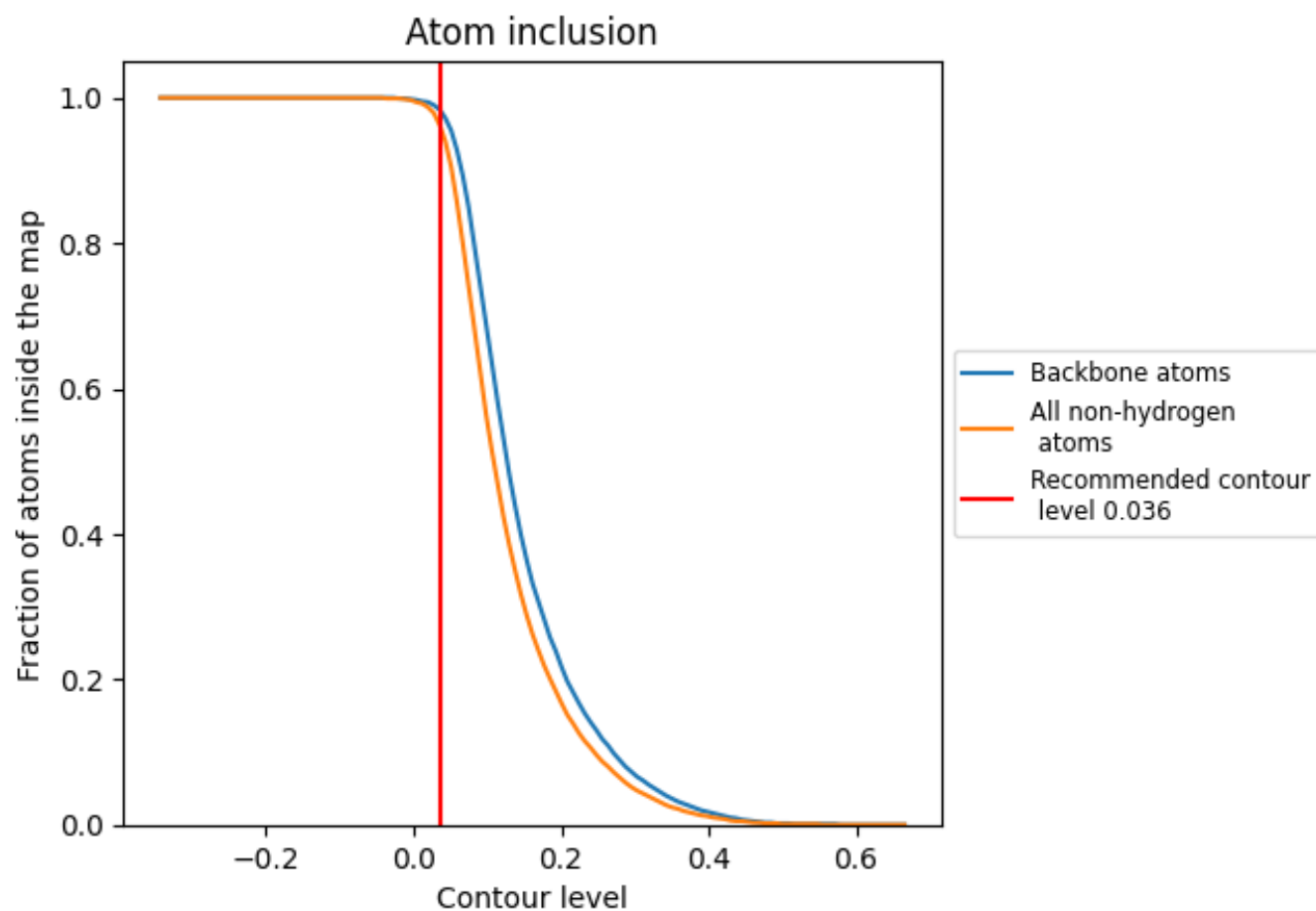
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.036).

9.4 Atom inclusion [i](#)



At the recommended contour level, 98% of all backbone atoms, 96% 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 (0.036) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.9610	0.3910
A	0.9640	0.3950
B	0.9430	0.3980
C	0.9790	0.4460
D	0.9660	0.4080
E	0.9530	0.2690
F	0.9450	0.2760
I	0.9570	0.2430
J	0.9550	0.2480

1.0

0.0

<0.0