



Full wwPDB EM Validation Report ⓘ

Mar 6, 2026 – 10:01 PM UTC

PDB ID : 8Q3R / pdb_00008q3r
EMDB ID : EMD-18134
Title : Cryo-EM structure of the DNA polymerase holoenzyme E9-A20-D4 of vaccinia virus
Authors : Burmeister, W.P.; Ballandras-Colas, A.; Boettcher, B.; Grimm, C.
Deposited on : 2023-08-04
Resolution : 3.80 Å(reported)
Based on initial models : 4od8, 8hg1, .

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

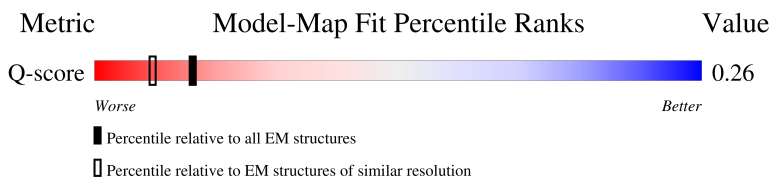
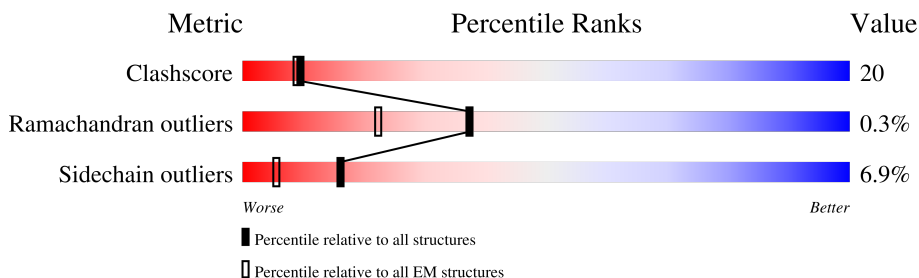
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.80 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	10198 (3.30 - 4.30)

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	D	242	
2	A	426	
3	E	1033	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 11921 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 Uracil-DNA glycosylase.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	D	218	Total	C	N	O	S	0	0
			1771	1147	293	325	6		

There are 25 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	-23	MET	-	initiating methionine	UNP P20536
D	-22	ALA	-	expression tag	UNP P20536
D	-21	SER	-	expression tag	UNP P20536
D	-20	TRP	-	expression tag	UNP P20536
D	-19	SER	-	expression tag	UNP P20536
D	-18	HIS	-	expression tag	UNP P20536
D	-17	PRO	-	expression tag	UNP P20536
D	-16	GLN	-	expression tag	UNP P20536
D	-15	PHE	-	expression tag	UNP P20536
D	-14	GLU	-	expression tag	UNP P20536
D	-13	LYS	-	expression tag	UNP P20536
D	-12	SER	-	expression tag	UNP P20536
D	-11	GLY	-	expression tag	UNP P20536
D	-10	GLY	-	expression tag	UNP P20536
D	-9	GLY	-	expression tag	UNP P20536
D	-8	GLY	-	expression tag	UNP P20536
D	-7	GLY	-	expression tag	UNP P20536
D	-6	LEU	-	expression tag	UNP P20536
D	-5	VAL	-	expression tag	UNP P20536
D	-4	PRO	-	expression tag	UNP P20536
D	-3	ARG	-	expression tag	UNP P20536
D	-2	GLY	-	expression tag	UNP P20536
D	-1	SER	-	expression tag	UNP P20536
D	0	ALA	-	expression tag	UNP P20536
D	208	ALA	VAL	conflict	UNP P20536

- Molecule 2 is a protein called DNA polymerase processivity factor component OPG148.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	A	420	Total	C	N	O	S	0	0
			3430	2209	560	651	10		

- Molecule 3 is a protein called DNA polymerase.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	E	827	Total	C	N	O	S	0	0
			6720	4298	1119	1261	42		

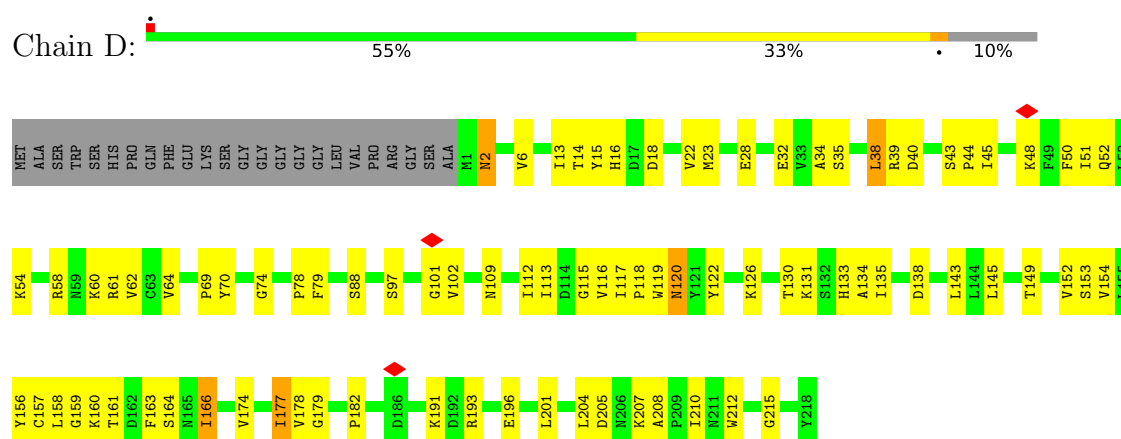
There are 28 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	-26	MET	-	initiating methionine	UNP P20509
E	-25	SER	-	expression tag	UNP P20509
E	-24	TYR	-	expression tag	UNP P20509
E	-23	TYR	-	expression tag	UNP P20509
E	-22	HIS	-	expression tag	UNP P20509
E	-21	HIS	-	expression tag	UNP P20509
E	-20	HIS	-	expression tag	UNP P20509
E	-19	HIS	-	expression tag	UNP P20509
E	-18	HIS	-	expression tag	UNP P20509
E	-17	HIS	-	expression tag	UNP P20509
E	-16	ASP	-	expression tag	UNP P20509
E	-15	TYR	-	expression tag	UNP P20509
E	-14	ASP	-	expression tag	UNP P20509
E	-13	ILE	-	expression tag	UNP P20509
E	-12	PRO	-	expression tag	UNP P20509
E	-11	THR	-	expression tag	UNP P20509
E	-10	THR	-	expression tag	UNP P20509
E	-9	GLU	-	expression tag	UNP P20509
E	-8	ASN	-	expression tag	UNP P20509
E	-7	LEU	-	expression tag	UNP P20509
E	-6	TYR	-	expression tag	UNP P20509
E	-5	PHE	-	expression tag	UNP P20509
E	-4	GLN	-	expression tag	UNP P20509
E	-3	GLY	-	expression tag	UNP P20509
E	-2	ALA	-	expression tag	UNP P20509
E	-1	MET	-	expression tag	UNP P20509
E	0	ASP	-	expression tag	UNP P20509
E	1	PRO	-	expression tag	UNP P20509

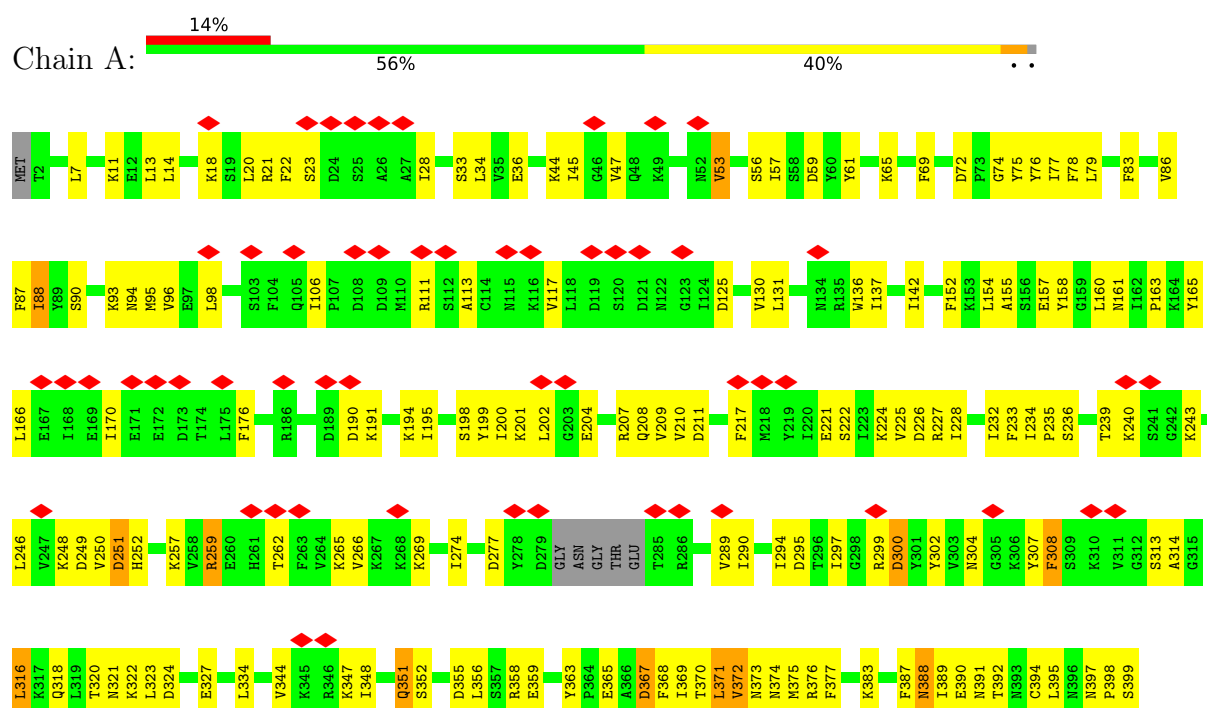
3 Residue-property plots

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: Uracil-DNA glycosylase

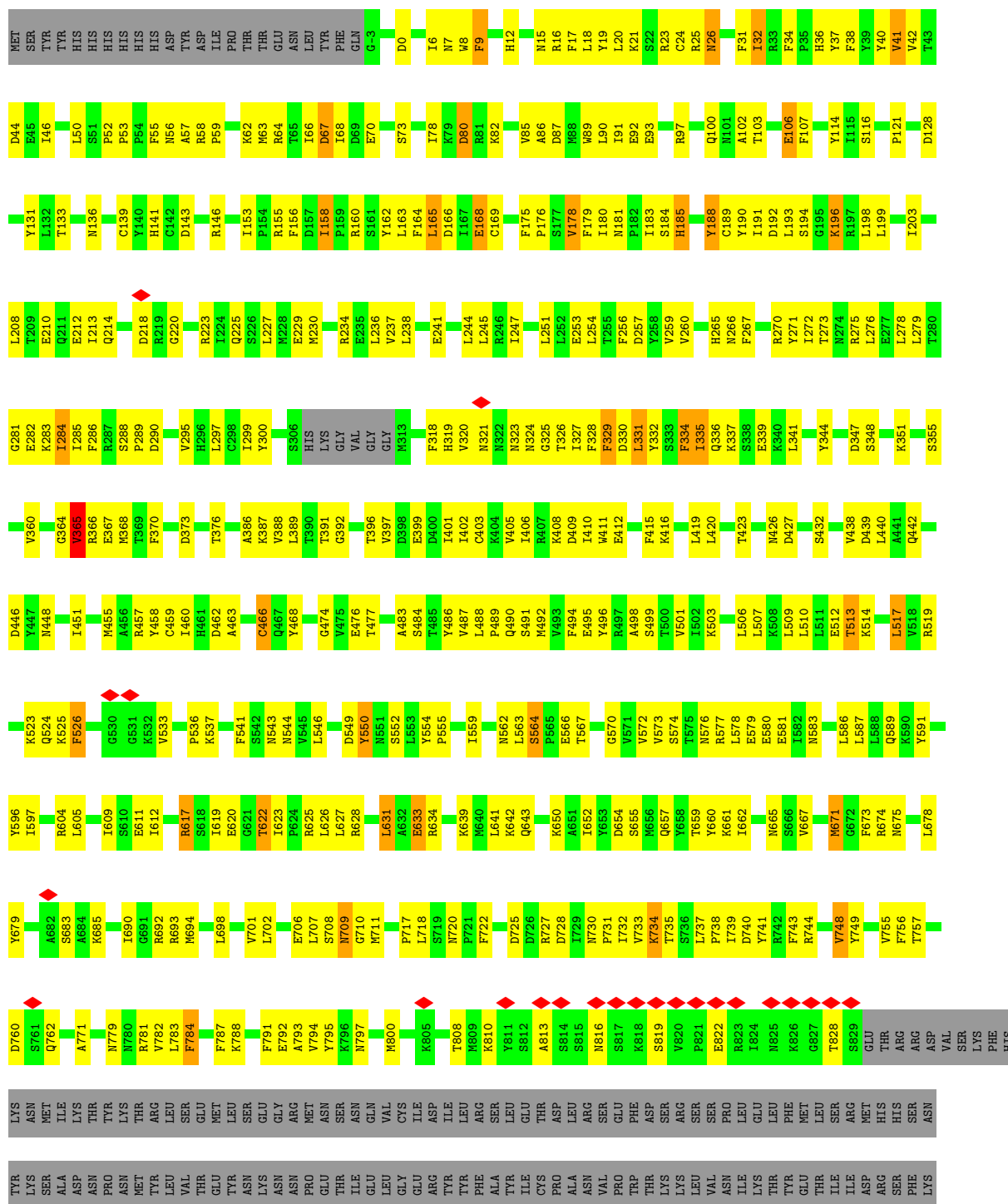


• Molecule 2: DNA polymerase processivity factor component OPG148





• Molecule 3: DNA polymerase



LEU	GLY	SER	ASP	GLN	ARG	ILE	PHE	TYR	GLU	VAL	TYR	PHE	LYS	ARG	LEU	THR	SER	GLU	ILE	VAL	ASN	LEU	LEU	ASP	ASN	LYS	VAL	LEU	CYS	ILE	SER	PHE	PHE	GLU	ARG	MET	PHE	GLY	SER	LYS	PRO	THR	PHE	TYR	GLU	ALA
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4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	104239	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 ($e^-/\text{\AA}^2$)	68	Depositor
Minimum defocus (nm)	1400	Depositor
Maximum defocus (nm)	2400	Depositor
Magnification	42000	Depositor
Image detector	FEI FALCON III (4k x 4k)	Depositor
Maximum map value	3.666	Depositor
Minimum map value	-1.485	Depositor
Average map value	0.004	Depositor
Map value standard deviation	0.061	Depositor
Recommended contour level	0.224	Depositor
Map size (Å)	272.384, 272.384, 272.384	wwPDB
Map dimensions	256, 256, 256	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.064, 1.064, 1.064	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	D	0.18	0/1820	0.34	0/2474
2	A	0.14	0/3496	0.29	0/4713
3	E	0.27	0/6864	0.38	0/9278
All	All	0.23	0/12180	0.35	0/16465

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	1771	0	1766	60	0
2	A	3430	0	3438	131	0
3	E	6720	0	6687	297	0
All	All	11921	0	11891	476	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 20.

All (476) 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)
2:A:368:PHE:HA	2:A:371:LEU:HD23	1.60	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:8:TRP:HB2	3:E:489:PRO:HB3	1.62	0.82
2:A:21:ARG:HB3	2:A:56:SER:H	1.45	0.82
2:A:78:PHE:HB2	2:A:166:LEU:HB2	1.60	0.81
2:A:371:LEU:HB2	2:A:400:ILE:HG23	1.64	0.79
2:A:74:GLY:HA3	2:A:201:LYS:HE2	1.66	0.77
2:A:348:ILE:HD12	2:A:356:LEU:HD21	1.65	0.77
3:E:70:GLU:OE2	3:E:604:ARG:NH1	2.18	0.76
2:A:404:TYR:HA	2:A:407:PHE:HB2	1.67	0.76
3:E:339:GLU:HB3	3:E:341:LEU:HG	1.68	0.75
3:E:67:ASP:HB3	3:E:82:LYS:HB2	1.69	0.75
3:E:365:VAL:HG13	3:E:367:GLU:H	1.52	0.75
3:E:523:LYS:NZ	3:E:674:ARG:O	2.16	0.73
3:E:396:THR:HB	3:E:432:SER:HB2	1.69	0.73
3:E:265:HIS:NE2	3:E:329:PHE:O	2.21	0.73
3:E:562:ASN:ND2	3:E:619:ILE:O	2.22	0.73
2:A:225:VAL:HA	2:A:235:PRO:HA	1.71	0.72
2:A:419:ASP:OD1	2:A:423:ARG:NH2	2.22	0.72
3:E:604:ARG:HG3	3:E:605:LEU:HG	1.72	0.71
3:E:163:LEU:HD23	3:E:259:VAL:HG22	1.73	0.71
3:E:199:LEU:HD22	3:E:460:ILE:HD11	1.73	0.71
1:D:39:ARG:NH1	3:E:278:LEU:O	2.24	0.70
3:E:64:ARG:HA	3:E:85:VAL:HA	1.72	0.70
3:E:164:PHE:HD2	3:E:466:CYS:HB3	1.56	0.70
3:E:116:SER:OG	3:E:146:ARG:NH2	2.24	0.70
3:E:178:VAL:O	3:E:275:ARG:NH1	2.24	0.70
2:A:376:ARG:NH1	3:E:581:GLU:OE2	2.24	0.70
3:E:816:ASN:OD1	3:E:819:SER:OG	2.11	0.69
3:E:20:LEU:HB3	3:E:32:ILE:HG13	1.72	0.69
2:A:370:THR:O	2:A:374:ASN:ND2	2.27	0.68
3:E:164:PHE:HB2	3:E:189:CYS:HB3	1.76	0.68
3:E:160:ARG:NH1	3:E:476:GLU:OE2	2.27	0.68
2:A:392:THR:HA	2:A:395:LEU:HD23	1.76	0.68
2:A:397:ASN:HB3	2:A:400:ILE:HG12	1.76	0.68
2:A:411:VAL:O	2:A:415:ASN:ND2	2.25	0.68
3:E:56:ASN:ND2	3:E:92:GLU:OE1	2.27	0.68
3:E:265:HIS:HD2	3:E:330:ASP:HB2	1.58	0.67
2:A:321:ASN:OD1	2:A:322:LYS:N	2.26	0.67
3:E:718:LEU:O	3:E:727:ARG:NH2	2.26	0.67
3:E:546:LEU:HB2	3:E:757:THR:HB	1.75	0.67
3:E:58:ARG:NH2	3:E:116:SER:OG	2.28	0.67
2:A:45:ILE:HG22	2:A:47:VAL:H	1.61	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:488:LEU:HD11	3:E:501:VAL:HG12	1.77	0.66
3:E:546:LEU:N	3:E:757:THR:O	2.26	0.66
2:A:314:ALA:HA	2:A:318:GLN:HG3	1.78	0.66
3:E:114:TYR:HD2	3:E:487:VAL:HG12	1.61	0.66
2:A:77:ILE:HB	2:A:200:ILE:HG12	1.78	0.65
3:E:728:ASP:OD2	3:E:781:ARG:NH2	2.29	0.65
2:A:65:LYS:HD3	2:A:210:VAL:HG12	1.78	0.65
3:E:169:CYS:HA	3:E:184:SER:H	1.60	0.65
3:E:537:LYS:HD2	3:E:748:VAL:HA	1.78	0.65
1:D:201:LEU:HD23	1:D:204:LEU:HD12	1.79	0.64
3:E:579:GLU:HG2	3:E:583:ASN:HD21	1.61	0.64
2:A:222:SER:OG	2:A:224:LYS:NZ	2.30	0.64
3:E:783:LEU:HD13	3:E:787:PHE:HB3	1.78	0.64
3:E:734:LYS:NZ	3:E:735:THR:O	2.30	0.64
3:E:164:PHE:CD2	3:E:466:CYS:HB3	2.33	0.63
1:D:160:LYS:NZ	1:D:178:VAL:O	2.23	0.63
3:E:552:SER:HB3	3:E:555:PRO:HB2	1.79	0.63
3:E:26:ASN:OD1	3:E:26:ASN:N	2.31	0.63
2:A:376:ARG:N	2:A:388:ASN:O	2.28	0.62
2:A:224:LYS:O	2:A:236:SER:N	2.32	0.62
2:A:294:ILE:HD12	2:A:398:PRO:HB2	1.81	0.62
2:A:249:ASP:OD1	2:A:250:VAL:N	2.32	0.62
1:D:61:ARG:NE	1:D:208:ALA:O	2.31	0.62
3:E:718:LEU:HD22	3:E:782:VAL:HG13	1.82	0.61
2:A:375:MET:HA	2:A:390:GLU:H	1.64	0.61
3:E:389:LEU:HB3	3:E:410:ILE:HD11	1.81	0.61
2:A:195:ILE:O	2:A:269:LYS:NZ	2.33	0.60
2:A:300:ASP:OD1	2:A:300:ASP:N	2.26	0.60
3:E:483:ALA:HB2	3:E:490:GLN:HG3	1.83	0.60
3:E:59:PRO:HA	3:E:89:TRP:HA	1.83	0.60
2:A:211:ASP:OD1	2:A:269:LYS:NZ	2.34	0.60
3:E:9:PHE:HE1	3:E:19:TYR:HB2	1.66	0.60
3:E:324:ASN:OD1	3:E:325:GLY:N	2.34	0.60
2:A:246:LEU:HB2	2:A:274:ILE:HG22	1.83	0.60
1:D:112:ILE:HG13	1:D:215:GLY:HA2	1.84	0.60
3:E:196:LYS:HB3	3:E:198:LEU:HD13	1.83	0.60
3:E:42:VAL:HG12	3:E:102:ALA:HA	1.83	0.59
3:E:203:ILE:HB	3:E:238:LEU:HD22	1.83	0.59
3:E:347:ASP:OD1	3:E:348:SER:N	2.35	0.59
3:E:9:PHE:HE2	3:E:325:GLY:HA2	1.67	0.59
3:E:411:TRP:CD1	3:E:412:GLU:H	2.19	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:95:MET:HG3	2:A:106:ILE:HG21	1.84	0.59
3:E:220:GLY:O	3:E:234:ARG:NH2	2.35	0.59
2:A:208:GLN:N	2:A:208:GLN:OE1	2.34	0.59
3:E:639:LYS:NZ	3:E:643:GLN:OE1	2.34	0.59
3:E:633:GLU:HB3	3:E:660:TYR:CE2	2.38	0.58
1:D:62:VAL:HG22	1:D:154:VAL:HB	1.85	0.58
3:E:80:ASP:N	3:E:80:ASP:OD1	2.37	0.58
3:E:163:LEU:HG	3:E:164:PHE:N	2.18	0.58
2:A:257:LYS:HG2	2:A:289:VAL:HG21	1.86	0.58
3:E:282:GLU:OE1	3:E:282:GLU:N	2.36	0.58
3:E:18:LEU:HB3	3:E:34:PHE:HB2	1.86	0.58
3:E:573:VAL:HG12	3:E:611:GLU:HA	1.86	0.58
1:D:97:SER:O	1:D:101:GLY:N	2.34	0.58
1:D:97:SER:OG	1:D:102:VAL:O	2.20	0.58
3:E:92:GLU:OE2	3:E:146:ARG:NH2	2.36	0.58
3:E:23:ARG:NH1	3:E:257:ASP:OD2	2.32	0.57
3:E:706:GLU:HA	3:E:744:ARG:HA	1.85	0.57
3:E:725:ASP:OD2	3:E:727:ARG:NE	2.37	0.57
1:D:40:ASP:HB2	1:D:133:HIS:CE1	2.39	0.57
2:A:294:ILE:HG12	2:A:308:PHE:HZ	1.70	0.57
2:A:94:ASN:OD1	2:A:111:ARG:NH1	2.37	0.57
3:E:779:ASN:HB3	3:E:788:LYS:NZ	2.20	0.57
1:D:64:VAL:HG13	1:D:156:TYR:HD2	1.69	0.57
3:E:267:PHE:O	3:E:271:TYR:HB2	2.06	0.56
2:A:373:ASN:HD21	3:E:609:ILE:HD13	1.69	0.56
3:E:270:ARG:HH21	3:E:299:ILE:HG21	1.71	0.56
3:E:509:LEU:HD22	3:E:626:LEU:HB2	1.87	0.56
3:E:720:ASN:HB3	3:E:725:ASP:HB2	1.88	0.56
2:A:313:SER:O	2:A:409:GLN:NE2	2.38	0.56
3:E:327:ILE:O	3:E:329:PHE:N	2.38	0.56
3:E:360:VAL:HG22	3:E:370:PHE:HE2	1.69	0.56
3:E:8:TRP:O	3:E:491:SER:OG	2.23	0.56
3:E:63:MET:HE3	3:E:514:LYS:HA	1.87	0.56
3:E:57:ALA:HA	3:E:91:ILE:HA	1.86	0.56
1:D:51:ILE:HA	1:D:54:LYS:HE2	1.87	0.55
2:A:13:LEU:HD13	2:A:34:LEU:HB2	1.87	0.55
2:A:351:GLN:NE2	2:A:355:ASP:OD2	2.39	0.55
3:E:9:PHE:CE2	3:E:325:GLY:HA2	2.41	0.55
3:E:392:GLY:O	3:E:408:LYS:NZ	2.38	0.55
2:A:374:ASN:HB3	2:A:391:ASN:HB3	1.89	0.55
3:E:21:LYS:NZ	3:E:256:PHE:O	2.39	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:417:VAL:O	2:A:421:LYS:HG2	2.06	0.55
3:E:286:PHE:HZ	3:E:297:LEU:HD12	1.71	0.55
3:E:492:MET:SD	3:E:496:TYR:OH	2.58	0.55
2:A:377:PHE:CD2	3:E:578:LEU:HD13	2.42	0.55
3:E:318:PHE:O	3:E:319:HIS:ND1	2.39	0.55
2:A:21:ARG:HD2	2:A:56:SER:HB3	1.87	0.55
3:E:265:HIS:HE1	3:E:328:PHE:HB3	1.71	0.55
3:E:284:ILE:HD12	3:E:297:LEU:HD13	1.88	0.55
3:E:566:GLU:OE2	3:E:622:THR:N	2.40	0.55
1:D:113:ILE:HD13	1:D:212:TRP:HA	1.88	0.54
3:E:190:TYR:HD1	3:E:191:ILE:N	2.05	0.54
3:E:56:ASN:O	3:E:92:GLU:N	2.32	0.54
3:E:58:ARG:N	3:E:90:LEU:O	2.34	0.54
3:E:78:ILE:HD13	3:E:570:GLY:HA3	1.89	0.54
3:E:331:LEU:O	3:E:335:ILE:HB	2.07	0.54
3:E:512:GLU:OE2	3:E:625:ARG:NH2	2.41	0.54
2:A:363:TYR:HE2	2:A:410:PHE:HB2	1.72	0.54
2:A:190:ASP:O	2:A:194:LYS:NZ	2.40	0.54
3:E:116:SER:HG	3:E:146:ARG:HH21	1.52	0.54
2:A:351:GLN:HB3	2:A:355:ASP:HB2	1.90	0.54
3:E:62:LYS:HA	3:E:87:ASP:HA	1.89	0.54
3:E:633:GLU:HB3	3:E:660:TYR:HE2	1.73	0.54
3:E:59:PRO:HB3	3:E:89:TRP:CE2	2.43	0.54
3:E:52:PRO:O	3:E:97:ARG:NH2	2.41	0.54
3:E:351:LYS:O	3:E:355:SER:OG	2.26	0.54
3:E:64:ARG:CZ	3:E:66:ILE:HD11	2.38	0.53
3:E:160:ARG:HB2	3:E:162:TYR:CE1	2.43	0.53
3:E:442:GLN:NE2	3:E:446:ASP:OD2	2.41	0.53
3:E:457:ARG:HA	3:E:460:ILE:HG22	1.89	0.53
3:E:525:LYS:HG2	3:E:674:ARG:HH22	1.72	0.53
3:E:562:ASN:OD1	3:E:620:GLU:HA	2.09	0.53
2:A:228:ILE:HD11	2:A:307:TYR:HE1	1.73	0.53
3:E:68:ILE:HD13	3:E:517:LEU:HB3	1.89	0.53
3:E:203:ILE:HD11	3:E:236:LEU:HD22	1.91	0.53
3:E:265:HIS:CE1	3:E:328:PHE:HB3	2.43	0.53
3:E:330:ASP:OD1	3:E:332:TYR:N	2.42	0.53
2:A:294:ILE:HG12	2:A:308:PHE:CZ	2.43	0.53
3:E:136:ASN:HB2	3:E:139:CYS:HB2	1.91	0.53
3:E:579:GLU:HG2	3:E:583:ASN:ND2	2.24	0.53
2:A:228:ILE:HD11	2:A:307:TYR:CE1	2.43	0.53
3:E:241:GLU:O	3:E:245:LEU:HG	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:181:ASN:O	3:E:271:TYR:OH	2.26	0.53
2:A:316:LEU:HG	2:A:334:LEU:HD11	1.91	0.53
3:E:779:ASN:HB3	3:E:788:LYS:HZ2	1.74	0.53
3:E:52:PRO:HG2	3:E:97:ARG:NH2	2.24	0.52
2:A:388:ASN:OD1	2:A:388:ASN:N	2.41	0.52
3:E:519:ARG:NH2	3:E:679:TYR:O	2.35	0.52
1:D:35:SER:HA	1:D:38:LEU:HD22	1.92	0.52
2:A:28:ILE:HG13	2:A:53:VAL:HG21	1.90	0.52
3:E:396:THR:N	3:E:432:SER:O	2.28	0.52
3:E:701:VAL:HG12	3:E:702:LEU:HD23	1.92	0.52
3:E:337:LYS:NZ	3:E:494:PHE:O	2.30	0.52
2:A:368:PHE:HD1	2:A:403:ILE:HG21	1.74	0.52
3:E:329:PHE:HD1	3:E:330:ASP:H	1.57	0.52
3:E:555:PRO:HA	3:E:627:LEU:HD13	1.91	0.52
3:E:707:LEU:HB3	3:E:743:PHE:H	1.75	0.52
2:A:190:ASP:OD1	2:A:191:LYS:N	2.43	0.52
3:E:63:MET:HG2	3:E:514:LYS:HB3	1.92	0.52
3:E:160:ARG:HB2	3:E:162:TYR:CZ	2.45	0.52
1:D:60:LYS:HA	1:D:115:GLY:HA2	1.92	0.52
3:E:18:LEU:O	3:E:34:PHE:N	2.34	0.51
3:E:8:TRP:CD2	3:E:20:LEU:HD13	2.45	0.51
3:E:448:ASN:HD21	3:E:451:ILE:HD12	1.75	0.51
3:E:641:LEU:HD23	3:E:642:LYS:HG3	1.92	0.51
2:A:302:TYR:OH	2:A:304:ASN:ND2	2.44	0.51
3:E:223:ARG:NE	3:E:225:GLN:HE21	2.08	0.51
3:E:8:TRP:CZ3	3:E:18:LEU:HD21	2.46	0.51
1:D:78:PRO:HB2	1:D:118:PRO:HB2	1.93	0.51
1:D:135:ILE:HG22	3:E:179:PHE:HZ	1.75	0.51
2:A:377:PHE:CG	3:E:578:LEU:HD13	2.46	0.51
2:A:294:ILE:HA	2:A:297:ILE:HG12	1.93	0.51
3:E:693:ARG:HG2	3:E:722:PHE:HE1	1.76	0.51
3:E:738:PRO:HB2	3:E:740:ASP:OD1	2.11	0.50
3:E:254:LEU:O	3:E:256:PHE:N	2.44	0.50
3:E:265:HIS:CD2	3:E:330:ASP:HB2	2.44	0.50
3:E:474:GLY:O	3:E:477:THR:OG1	2.29	0.50
3:E:709:ASN:HB3	3:E:739:ILE:HG23	1.93	0.50
3:E:744:ARG:H	3:E:760:ASP:CG	2.20	0.50
2:A:86:VAL:HG12	2:A:130:VAL:HG22	1.93	0.50
2:A:406:ASN:HB2	2:A:409:GLN:HG3	1.93	0.50
3:E:41:VAL:HA	3:E:89:TRP:O	2.12	0.50
3:E:40:TYR:HB2	3:E:91:ILE:HD11	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:183:ILE:HD12	3:E:245:LEU:HD21	1.94	0.50
2:A:376:ARG:CZ	3:E:577:ARG:HD3	2.42	0.50
3:E:330:ASP:OD1	3:E:331:LEU:N	2.45	0.50
3:E:25:ARG:HB3	3:E:156:PHE:CE2	2.47	0.49
2:A:251:ASP:N	2:A:251:ASP:OD1	2.44	0.49
3:E:791:PHE:HE2	3:E:813:ALA:HB2	1.77	0.49
2:A:221:GLU:OE1	2:A:221:GLU:N	2.45	0.49
2:A:367:ASP:OD1	2:A:367:ASP:N	2.43	0.49
3:E:139:CYS:SG	3:E:289:PRO:HB2	2.53	0.49
3:E:446:ASP:O	3:E:448:ASN:ND2	2.46	0.49
2:A:88:ILE:HG13	2:A:96:VAL:HG23	1.95	0.49
2:A:259:ARG:O	2:A:262:THR:OG1	2.31	0.49
3:E:373:ASP:N	3:E:376:THR:OG1	2.42	0.49
2:A:83:PHE:CZ	2:A:248:LYS:HG3	2.47	0.49
3:E:405:VAL:HA	3:E:419:LEU:HD23	1.93	0.49
1:D:61:ARG:O	1:D:153:SER:N	2.45	0.49
2:A:318:GLN:HA	2:A:321:ASN:ND2	2.28	0.49
2:A:320:THR:OG1	2:A:327:GLU:OE1	2.31	0.49
2:A:399:SER:O	2:A:403:ILE:HG12	2.12	0.49
3:E:106:GLU:H	3:E:106:GLU:CD	2.21	0.49
3:E:164:PHE:CD2	3:E:260:VAL:HB	2.48	0.49
1:D:157:CYS:HB3	1:D:163:PHE:CG	2.47	0.49
2:A:14:LEU:O	2:A:18:LYS:HG2	2.13	0.49
3:E:549:ASP:HB2	3:E:792:GLU:HG2	1.94	0.49
3:E:554:TYR:HB2	3:E:555:PRO:HD3	1.95	0.49
3:E:586:LEU:HA	3:E:589:GLN:NE2	2.28	0.49
3:E:121:PRO:HG2	3:E:483:ALA:O	2.13	0.48
1:D:62:VAL:O	1:D:116:VAL:HA	2.12	0.48
3:E:8:TRP:HE3	3:E:18:LEU:HD11	1.77	0.48
3:E:397:VAL:HB	3:E:401:ILE:HB	1.95	0.48
1:D:40:ASP:HB3	1:D:126:LYS:HG3	1.95	0.48
1:D:158:LEU:HD23	1:D:182:PRO:HD3	1.95	0.48
2:A:401:GLU:O	2:A:405:ARG:HG2	2.13	0.48
3:E:574:SER:HB3	3:E:579:GLU:HB3	1.94	0.48
3:E:617:ARG:HH11	3:E:617:ARG:HB2	1.78	0.48
3:E:44:ASP:OD1	3:E:89:TRP:NE1	2.47	0.48
2:A:420:VAL:HG23	2:A:423:ARG:HH22	1.77	0.48
3:E:457:ARG:O	3:E:460:ILE:HG22	2.14	0.48
2:A:226:ASP:N	2:A:234:ILE:O	2.42	0.48
3:E:166:ASP:HB3	3:E:463:ALA:HB2	1.94	0.48
2:A:204:GLU:OE1	2:A:204:GLU:N	2.46	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:130:THR:HG22	1:D:131:LYS:HG3	1.96	0.47
3:E:133:THR:HB	3:E:141:HIS:HB3	1.94	0.47
3:E:524:GLN:HB3	3:E:526:PHE:CZ	2.49	0.47
1:D:13:ILE:HG23	1:D:50:PHE:HB3	1.95	0.47
3:E:214:GLN:NE2	3:E:218:ASP:OD2	2.46	0.47
3:E:509:LEU:HD11	3:E:625:ARG:HG2	1.97	0.47
3:E:509:LEU:HD23	3:E:622:THR:HG23	1.96	0.47
3:E:717:PRO:HA	3:E:731:PRO:HB3	1.95	0.47
2:A:408:ASN:HA	2:A:411:VAL:HG22	1.95	0.47
3:E:136:ASN:ND2	3:E:290:ASP:O	2.47	0.47
3:E:760:ASP:OD1	3:E:760:ASP:N	2.47	0.47
2:A:228:ILE:HG13	2:A:232:ILE:HG22	1.97	0.47
3:E:286:PHE:CZ	3:E:297:LEU:HD12	2.49	0.47
1:D:145:LEU:HD12	1:D:145:LEU:HA	1.75	0.47
2:A:406:ASN:HB2	2:A:409:GLN:CG	2.45	0.47
3:E:283:LYS:O	3:E:285:ILE:HG13	2.14	0.47
3:E:509:LEU:O	3:E:513:THR:HG23	2.15	0.47
3:E:591:TYR:HB3	3:E:596:TYR:CG	2.49	0.47
1:D:149:THR:HA	1:D:152:VAL:HG22	1.96	0.47
1:D:174:VAL:HA	2:A:44:LYS:HD2	1.96	0.47
2:A:90:SER:HB3	2:A:93:LYS:HD2	1.96	0.47
2:A:377:PHE:CZ	3:E:578:LEU:HB2	2.49	0.47
3:E:46:ILE:O	3:E:50:LEU:HG	2.14	0.47
3:E:114:TYR:CD2	3:E:487:VAL:HG12	2.47	0.47
3:E:409:ASP:O	3:E:416:LYS:N	2.30	0.47
3:E:543:ASN:HA	3:E:762:GLN:HG2	1.97	0.47
1:D:159:GLY:O	1:D:163:PHE:HD2	1.98	0.47
2:A:347:LYS:HG3	2:A:348:ILE:HD13	1.96	0.47
3:E:0:ASP:HB3	3:E:24:CYS:SG	2.55	0.47
3:E:199:LEU:HD13	3:E:460:ILE:HG12	1.97	0.47
3:E:440:LEU:HD22	3:E:455:MET:HE3	1.97	0.47
2:A:61:TYR:HA	2:A:207:ARG:O	2.15	0.47
3:E:169:CYS:HA	3:E:184:SER:N	2.29	0.47
3:E:397:VAL:N	3:E:401:ILE:O	2.34	0.47
3:E:634:ARG:NH1	3:E:661:LYS:HD3	2.30	0.47
3:E:163:LEU:HD22	3:E:256:PHE:CE2	2.50	0.46
2:A:227:ARG:HA	2:A:233:PHE:HD1	1.80	0.46
2:A:363:TYR:CE1	2:A:403:ILE:HD12	2.50	0.46
2:A:418:THR:HG22	2:A:422:LYS:HE3	1.98	0.46
3:E:12:HIS:CE1	3:E:321:ASN:HD22	2.33	0.46
3:E:730:ASN:HD21	3:E:732:ILE:HB	1.80	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:66:ILE:HD13	3:E:566:GLU:HB2	1.97	0.46
3:E:323:ASN:O	3:E:323:ASN:ND2	2.49	0.46
3:E:711:MET:SD	3:E:734:LYS:HE2	2.56	0.46
2:A:407:PHE:O	2:A:411:VAL:HG13	2.15	0.46
3:E:107:PHE:CD2	3:E:503:LYS:HB3	2.51	0.46
3:E:116:SER:HG	3:E:146:ARG:NH2	2.10	0.46
3:E:399:GLU:OE1	3:E:432:SER:OG	2.32	0.46
3:E:579:GLU:O	3:E:583:ASN:ND2	2.36	0.46
3:E:498:ALA:HA	3:E:501:VAL:HG22	1.96	0.46
3:E:288:SER:OG	3:E:290:ASP:OD1	2.28	0.46
1:D:16:HIS:CD2	1:D:18:ASP:H	2.34	0.46
1:D:69:PRO:HD3	1:D:122:TYR:O	2.15	0.46
3:E:158:ILE:HD11	3:E:160:ARG:HH21	1.81	0.46
3:E:365:VAL:HG22	3:E:366:ARG:H	1.80	0.46
1:D:2:ASN:HB2	1:D:15:TYR:CE1	2.51	0.46
3:E:106:GLU:O	3:E:503:LYS:NZ	2.28	0.46
3:E:107:PHE:CE2	3:E:507:LEU:HB2	2.51	0.46
1:D:43:SER:OG	1:D:70:TYR:O	2.26	0.46
1:D:52:GLN:HB3	1:D:119:TRP:CZ3	2.51	0.46
3:E:426:ASN:OD1	3:E:427:ASP:N	2.49	0.46
3:E:163:LEU:HD13	3:E:256:PHE:CE2	2.51	0.45
3:E:36:HIS:CG	3:E:37:TYR:N	2.84	0.45
3:E:368:MET:SD	3:E:423:THR:HA	2.57	0.45
1:D:40:ASP:OD2	1:D:133:HIS:NE2	2.46	0.45
3:E:175:PHE:CG	3:E:176:PRO:HD2	2.51	0.45
3:E:623:ILE:HD13	3:E:671:MET:SD	2.57	0.45
3:E:163:LEU:HD22	3:E:256:PHE:CD2	2.51	0.45
2:A:198:SER:HA	2:A:209:VAL:HG12	1.99	0.45
1:D:163:PHE:HD1	1:D:166:ILE:HG21	1.81	0.45
3:E:406:ILE:HD11	3:E:420:LEU:HB2	1.99	0.45
2:A:59:ASP:N	2:A:59:ASP:OD1	2.45	0.45
2:A:234:ILE:HD12	2:A:235:PRO:HD2	1.99	0.45
2:A:383:LYS:HE3	2:A:422:LYS:NZ	2.31	0.45
3:E:192:ASP:OD1	3:E:193:LEU:N	2.47	0.45
3:E:281:GLY:N	3:E:282:GLU:OE1	2.50	0.45
3:E:62:LYS:HG2	3:E:85:VAL:HG12	1.98	0.45
1:D:122:TYR:CZ	1:D:134:ALA:HB1	2.52	0.45
1:D:166:ILE:H	1:D:166:ILE:HG13	1.55	0.45
1:D:43:SER:C	1:D:74:GLY:HA3	2.43	0.44
3:E:225:GLN:N	3:E:229:GLU:OE1	2.48	0.44
3:E:455:MET:HA	3:E:458:TYR:CD2	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:8:TRP:HZ3	3:E:18:LEU:HD21	1.81	0.44
3:E:793:ALA:HB1	3:E:810:LYS:HG3	2.00	0.44
2:A:372:VAL:CG1	3:E:578:LEU:H	2.31	0.44
3:E:8:TRP:CH2	3:E:20:LEU:HD22	2.52	0.44
3:E:121:PRO:HG2	3:E:484:SER:HA	1.99	0.44
3:E:185:HIS:HB3	3:E:203:ILE:HA	2.00	0.44
1:D:2:ASN:HB2	1:D:15:TYR:CZ	2.52	0.44
3:E:673:PHE:HD1	3:E:675:ASN:H	1.66	0.44
1:D:44:PRO:HB2	1:D:48:LYS:HB2	2.00	0.44
1:D:61:ARG:NH2	1:D:210:ILE:HD13	2.33	0.44
2:A:155:ALA:HB1	2:A:160:LEU:HD11	1.99	0.44
3:E:300:TYR:OH	3:E:321:ASN:HB3	2.18	0.44
3:E:623:ILE:HD11	3:E:678:LEU:HB3	2.00	0.44
3:E:87:ASP:OD1	3:E:87:ASP:N	2.51	0.44
3:E:403:CYS:SG	3:E:419:LEU:HB3	2.58	0.44
2:A:299:ARG:HG3	2:A:405:ARG:NH1	2.32	0.44
2:A:394:CYS:SG	2:A:400:ILE:HG21	2.58	0.44
3:E:155:ARG:NH2	3:E:652:ILE:HG21	2.33	0.44
3:E:546:LEU:HD23	3:E:794:VAL:HG22	1.98	0.44
1:D:88:SER:HA	1:D:182:PRO:HB2	2.00	0.43
1:D:163:PHE:HB3	1:D:166:ILE:HG12	1.99	0.43
3:E:63:MET:HB3	3:E:514:LYS:O	2.18	0.43
3:E:387:LYS:HG2	3:E:468:TYR:CE1	2.53	0.43
3:E:388:VAL:HG12	3:E:388:VAL:O	2.18	0.43
1:D:156:TYR:HA	1:D:177:ILE:O	2.17	0.43
2:A:117:VAL:HG23	2:A:154:LEU:HD23	2.00	0.43
3:E:194:SER:OG	3:E:196:LYS:HB2	2.18	0.43
1:D:78:PRO:HG2	1:D:118:PRO:C	2.43	0.43
2:A:154:LEU:HG	2:A:158:TYR:CZ	2.54	0.43
2:A:217:PHE:CE2	2:A:265:LYS:HB3	2.53	0.43
3:E:290:ASP:OD1	3:E:290:ASP:N	2.49	0.43
3:E:506:LEU:O	3:E:510:LEU:HG	2.18	0.43
3:E:733:VAL:HG22	3:E:781:ARG:CZ	2.48	0.43
3:E:544:ASN:HD21	3:E:762:GLN:C	2.26	0.43
1:D:161:THR:O	1:D:164:SER:OG	2.28	0.43
2:A:290:ILE:O	2:A:294:ILE:HG13	2.19	0.43
3:E:334:PHE:HE2	3:E:494:PHE:CZ	2.37	0.43
3:E:364:GLY:O	3:E:365:VAL:HG12	2.18	0.43
2:A:295:ASP:OD1	2:A:299:ARG:NH1	2.44	0.43
3:E:327:ILE:HG23	3:E:491:SER:HA	2.00	0.43
3:E:795:TYR:CG	3:E:808:THR:HG21	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:77:ILE:O	2:A:199:TYR:HA	2.19	0.43
3:E:259:VAL:HG23	3:E:326:THR:HG21	2.01	0.43
2:A:348:ILE:HG23	2:A:356:LEU:HD21	1.99	0.43
1:D:205:ASP:HB3	1:D:207:LYS:HG2	2.01	0.42
2:A:125:ASP:HB3	2:A:142:ILE:HG13	2.01	0.42
2:A:375:MET:HA	2:A:390:GLU:N	2.31	0.42
3:E:388:VAL:HG13	3:E:468:TYR:CD2	2.54	0.42
1:D:28:GLU:O	1:D:32:GLU:HG2	2.18	0.42
1:D:60:LYS:HA	1:D:60:LYS:HD3	1.75	0.42
2:A:226:ASP:HB3	2:A:234:ILE:HG23	1.99	0.42
2:A:252:HIS:NE2	2:A:277:ASP:HB3	2.34	0.42
2:A:358:ARG:NH1	2:A:365:GLU:HG2	2.34	0.42
2:A:418:THR:O	2:A:422:LYS:HG3	2.19	0.42
3:E:15:ASN:HB3	3:E:17:PHE:CE2	2.54	0.42
3:E:536:PRO:HA	3:E:749:TYR:HD2	1.84	0.42
1:D:32:GLU:OE1	3:E:180:ILE:HG13	2.18	0.42
3:E:492:MET:HG2	3:E:495:GLU:OE1	2.19	0.42
3:E:168:GLU:CD	3:E:440:LEU:HD13	2.45	0.42
3:E:549:ASP:OD1	3:E:550:TYR:N	2.48	0.42
1:D:191:LYS:HB3	1:D:191:LYS:HE2	1.84	0.42
2:A:369:ILE:HG23	3:E:576:ASN:HB3	2.01	0.42
3:E:164:PHE:O	3:E:165:LEU:HB3	2.19	0.42
3:E:208:LEU:HD23	3:E:208:LEU:HA	1.67	0.42
3:E:388:VAL:HG22	3:E:468:TYR:CG	2.54	0.42
3:E:810:LYS:O	3:E:822:GLU:N	2.37	0.42
3:E:318:PHE:C	3:E:319:HIS:CG	2.96	0.42
1:D:18:ASP:OD2	1:D:58:ARG:HG3	2.19	0.42
2:A:323:LEU:HG	2:A:348:ILE:HD11	2.02	0.42
2:A:355:ASP:O	2:A:359:GLU:HG2	2.19	0.42
3:E:299:ILE:HD13	3:E:320:VAL:HG22	2.02	0.42
3:E:589:GLN:HE21	3:E:589:GLN:HB2	1.54	0.42
3:E:690:ILE:O	3:E:694:MET:HG2	2.20	0.42
3:E:694:MET:HE1	3:E:784:PHE:CZ	2.55	0.42
3:E:749:TYR:O	3:E:756:PHE:N	2.30	0.42
1:D:14:THR:HG22	1:D:54:LYS:HG2	2.02	0.42
2:A:20:LEU:O	2:A:23:SER:OG	2.31	0.42
2:A:33:SER:HA	2:A:36:GLU:OE1	2.20	0.42
2:A:239:THR:OG1	2:A:243:LYS:N	2.44	0.42
3:E:41:VAL:O	3:E:103:THR:HG22	2.20	0.42
3:E:386:ALA:O	3:E:410:ILE:HD13	2.20	0.42
3:E:409:ASP:N	3:E:416:LYS:O	2.36	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:546:LEU:HD22	3:E:771:ALA:HB2	2.02	0.42
2:A:69:PHE:O	2:A:176:PHE:N	2.52	0.42
3:E:336:GLN:HG2	3:E:344:TYR:CG	2.55	0.42
3:E:396:THR:HG23	3:E:402:ILE:HG12	2.02	0.42
3:E:564:SER:O	3:E:683:SER:OG	2.38	0.42
1:D:45:ILE:HG12	1:D:48:LYS:HG3	2.02	0.41
1:D:193:ARG:NE	1:D:196:GLU:OE1	2.49	0.41
2:A:76:TYR:CD2	2:A:170:ILE:HD13	2.55	0.41
2:A:240:LYS:HD3	2:A:240:LYS:HA	1.95	0.41
2:A:376:ARG:HA	2:A:376:ARG:HD3	1.55	0.41
3:E:230:MET:HE1	3:E:237:VAL:HG21	2.01	0.41
3:E:541:PHE:HE1	3:E:800:MET:HG2	1.84	0.41
3:E:627:LEU:HA	3:E:627:LEU:HD23	1.87	0.41
2:A:152:PHE:HE2	2:A:163:PRO:HG2	1.85	0.41
2:A:154:LEU:HD12	2:A:157:GLU:OE2	2.20	0.41
2:A:176:PHE:CE2	2:A:210:VAL:HG21	2.55	0.41
3:E:7:ASN:OD1	3:E:491:SER:OG	2.37	0.41
3:E:31:PHE:O	3:E:139:CYS:HA	2.21	0.41
3:E:188:TYR:N	3:E:188:TYR:CD1	2.88	0.41
3:E:227:LEU:HD13	3:E:247:ILE:HA	2.01	0.41
3:E:410:ILE:HG12	3:E:415:PHE:CB	2.50	0.41
3:E:738:PRO:HD2	3:E:741:TYR:CE1	2.55	0.41
1:D:79:PHE:CZ	1:D:158:LEU:HD13	2.56	0.41
3:E:38:PHE:N	3:E:93:GLU:O	2.43	0.41
3:E:410:ILE:HG12	3:E:415:PHE:HB3	2.02	0.41
3:E:438:VAL:HG21	3:E:458:TYR:CZ	2.55	0.41
3:E:698:LEU:HD12	3:E:698:LEU:HA	1.76	0.41
2:A:131:LEU:HD12	2:A:136:TRP:CH2	2.55	0.41
3:E:183:ILE:HG21	3:E:244:LEU:HD21	2.01	0.41
3:E:604:ARG:HB3	3:E:685:LYS:NZ	2.35	0.41
1:D:52:GLN:HG3	1:D:120:ASN:HD21	1.86	0.41
2:A:75:TYR:HB2	2:A:202:LEU:HD23	2.03	0.41
2:A:137:ILE:HG12	2:A:161:ASN:ND2	2.35	0.41
3:E:439:ASP:OD1	3:E:442:GLN:N	2.41	0.41
3:E:555:PRO:HA	3:E:627:LEU:CD1	2.50	0.41
3:E:708:SER:O	3:E:710:GLY:N	2.54	0.41
2:A:387:PHE:CE2	2:A:411:VAL:HG11	2.56	0.41
3:E:52:PRO:HA	3:E:53:PRO:HD3	1.95	0.41
1:D:34:ALA:O	1:D:38:LEU:HD13	2.20	0.41
1:D:156:TYR:HE1	1:D:179:GLY:HA3	1.85	0.41
2:A:7:LEU:O	2:A:11:LYS:HG2	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:79:LEU:HD21	2:A:165:TYR:HB3	2.01	0.41
2:A:392:THR:O	2:A:395:LEU:HB2	2.21	0.41
3:E:208:LEU:HD22	3:E:212:GLU:OE1	2.21	0.41
3:E:128:ASP:HB3	3:E:131:TYR:HB2	2.03	0.41
3:E:631:LEU:HD22	3:E:631:LEU:HA	1.83	0.41
1:D:15:TYR:CZ	1:D:23:MET:HG3	2.55	0.40
2:A:401:GLU:H	2:A:401:GLU:HG2	1.56	0.40
3:E:20:LEU:HB3	3:E:34:PHE:HE2	1.86	0.40
3:E:266:ASN:HB2	3:E:318:PHE:CD1	2.56	0.40
3:E:577:ARG:O	3:E:580:GLU:HB3	2.21	0.40
3:E:587:LEU:HD12	3:E:612:ILE:HD12	2.02	0.40
3:E:737:LEU:HD23	3:E:737:LEU:HA	1.85	0.40
2:A:113:ALA:HB1	2:A:160:LEU:HD23	2.04	0.40
3:E:55:PHE:N	3:E:93:GLU:OE1	2.35	0.40
3:E:73:SER:O	3:E:572:VAL:HA	2.22	0.40
2:A:87:PHE:HE2	2:A:131:LEU:HB2	1.86	0.40
2:A:320:THR:O	2:A:324:ASP:N	2.55	0.40
2:A:344:VAL:O	2:A:348:ILE:HG12	2.21	0.40
3:E:63:MET:N	3:E:86:ALA:O	2.51	0.40
1:D:109:ASN:ND2	1:D:112:ILE:HG23	2.36	0.40
3:E:210:GLU:HA	3:E:213:ILE:HD12	2.02	0.40
3:E:486:TYR:N	3:E:486:TYR:CD1	2.88	0.40
3:E:499:SER:O	3:E:503:LYS:HE3	2.21	0.40
3:E:559:ILE:HG21	3:E:628:ARG:HB2	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	D	216/242 (89%)	205 (95%)	11 (5%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	A	416/426 (98%)	388 (93%)	26 (6%)	2 (0%)	24	57
3	E	823/1033 (80%)	765 (93%)	56 (7%)	2 (0%)	43	73
All	All	1455/1701 (86%)	1358 (93%)	93 (6%)	4 (0%)	37	67

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	E	365	VAL
2	A	72	ASP
3	E	709	ASN
2	A	53	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	D	199/215 (93%)	189 (95%)	10 (5%)	22	47
2	A	390/394 (99%)	372 (95%)	18 (5%)	24	48
3	E	755/951 (79%)	690 (91%)	65 (9%)	10	33
All	All	1344/1560 (86%)	1251 (93%)	93 (7%)	16	40

All (93) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	2	ASN
1	D	6	VAL
1	D	22	VAL
1	D	38	LEU
1	D	117	ILE
1	D	120	ASN
1	D	138	ASP
1	D	143	LEU
1	D	166	ILE
1	D	177	ILE

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Mol	Chain	Res	Type
2	A	22	PHE
2	A	57	ILE
2	A	88	ILE
2	A	98	LEU
2	A	251	ASP
2	A	259	ARG
2	A	266	VAL
2	A	300	ASP
2	A	308	PHE
2	A	316	LEU
2	A	351	GLN
2	A	352	SER
2	A	367	ASP
2	A	371	LEU
2	A	372	VAL
2	A	388	ASN
2	A	389	ILE
2	A	420	VAL
3	E	6	ILE
3	E	9	PHE
3	E	16	ARG
3	E	26	ASN
3	E	32	ILE
3	E	41	VAL
3	E	67	ASP
3	E	80	ASP
3	E	100	GLN
3	E	106	GLU
3	E	143	ASP
3	E	153	ILE
3	E	158	ILE
3	E	165	LEU
3	E	168	GLU
3	E	178	VAL
3	E	185	HIS
3	E	188	TYR
3	E	196	LYS
3	E	251	LEU
3	E	253	GLU
3	E	272	ILE
3	E	273	THR
3	E	276	LEU

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Mol	Chain	Res	Type
3	E	279	LEU
3	E	284	ILE
3	E	295	VAL
3	E	329	PHE
3	E	331	LEU
3	E	334	PHE
3	E	335	ILE
3	E	365	VAL
3	E	391	THR
3	E	459	CYS
3	E	462	ASP
3	E	466	CYS
3	E	513	THR
3	E	517	LEU
3	E	526	PHE
3	E	533	VAL
3	E	550	TYR
3	E	563	LEU
3	E	564	SER
3	E	567	THR
3	E	597	ILE
3	E	617	ARG
3	E	622	THR
3	E	631	LEU
3	E	633	GLU
3	E	650	LYS
3	E	654	ASP
3	E	655	SER
3	E	657	GLN
3	E	659	THR
3	E	662	ILE
3	E	665	ASN
3	E	667	VAL
3	E	671	MET
3	E	692	ARG
3	E	734	LYS
3	E	748	VAL
3	E	755	VAL
3	E	784	PHE
3	E	797	ASN
3	E	828	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22)

such sidechains are listed below:

Mol	Chain	Res	Type
1	D	2	ASN
1	D	8	HIS
1	D	16	HIS
1	D	120	ASN
1	D	165	ASN
2	A	32	ASN
2	A	48	GLN
2	A	115	ASN
2	A	304	ASN
2	A	318	GLN
2	A	351	GLN
2	A	373	ASN
3	E	100	GLN
3	E	225	GLN
3	E	538	GLN
3	E	543	ASN
3	E	544	ASN
3	E	583	ASN
3	E	589	GLN
3	E	657	GLN
3	E	703	ASN
3	E	801	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

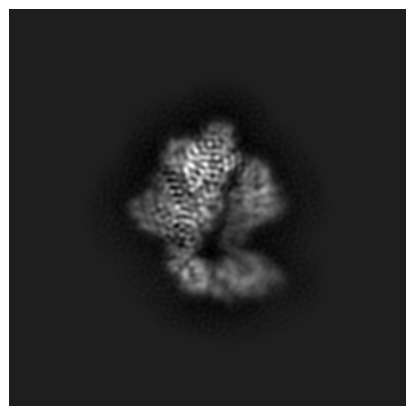
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-18134. These allow visual inspection of the internal detail of the map and identification of artifacts.

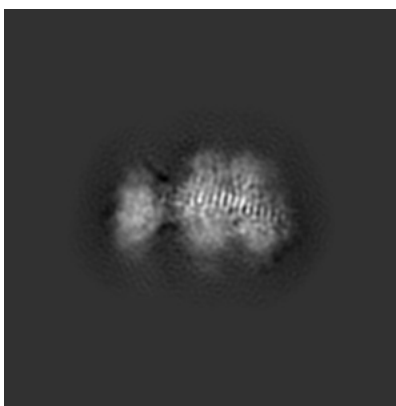
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

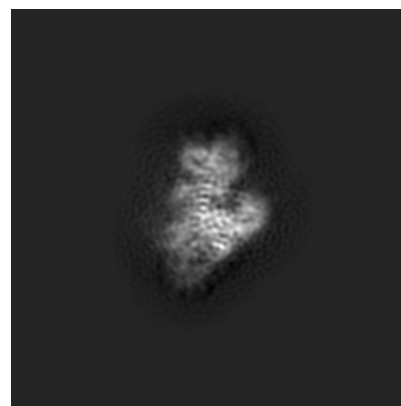
6.1.1 Primary map



X

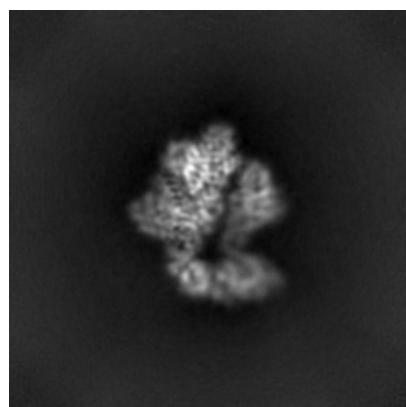


Y

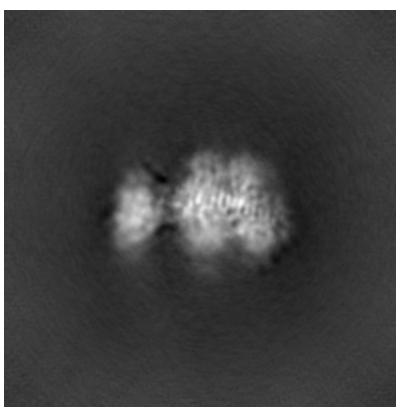


Z

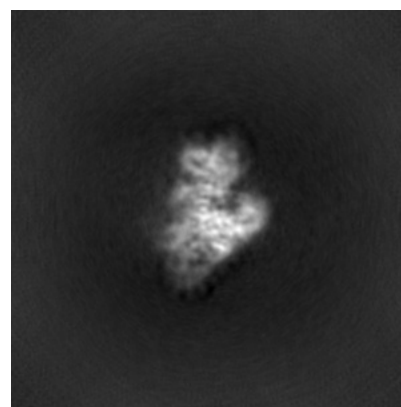
6.1.2 Raw map



X



Y

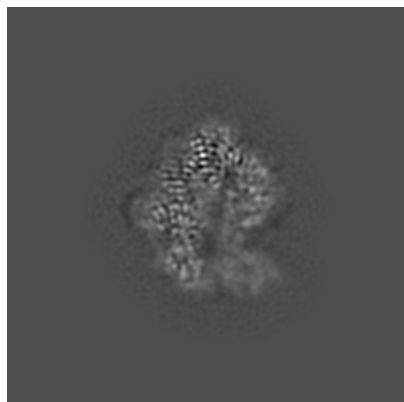


Z

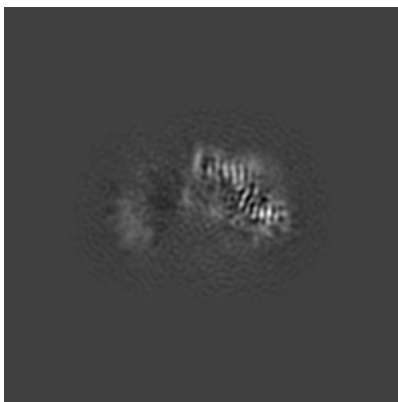
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

6.2.1 Primary map



X Index: 128

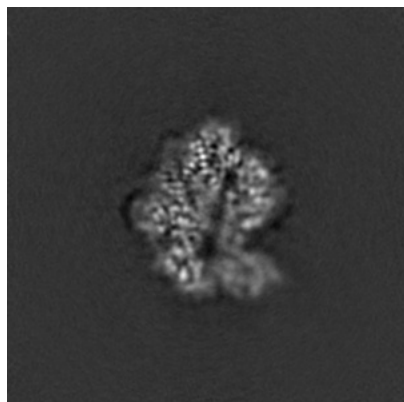


Y Index: 128

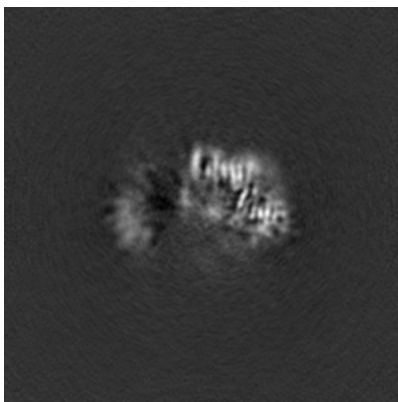


Z Index: 128

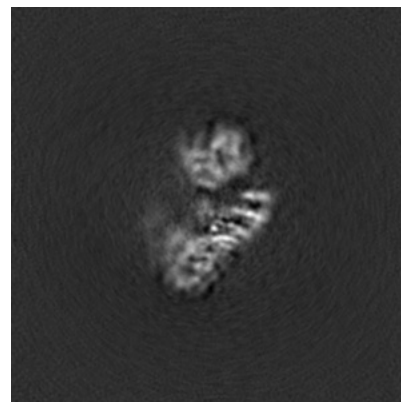
6.2.2 Raw map



X Index: 128



Y Index: 128

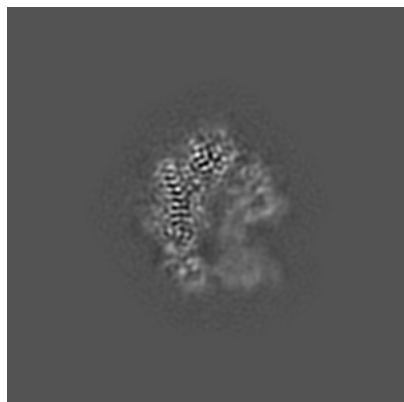


Z Index: 128

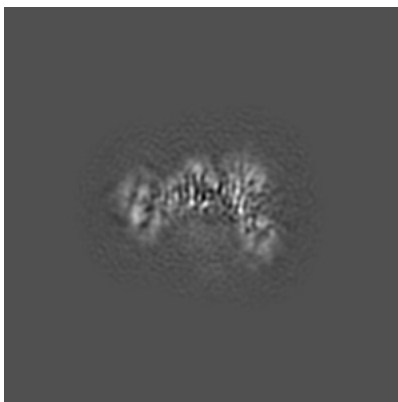
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

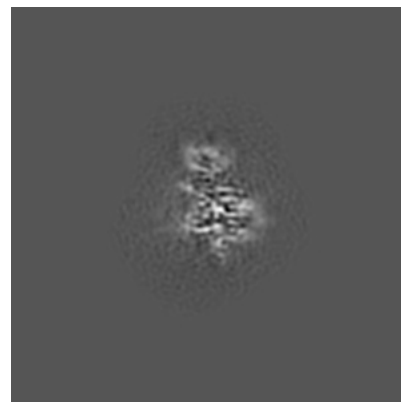
6.3.1 Primary map



X Index: 133

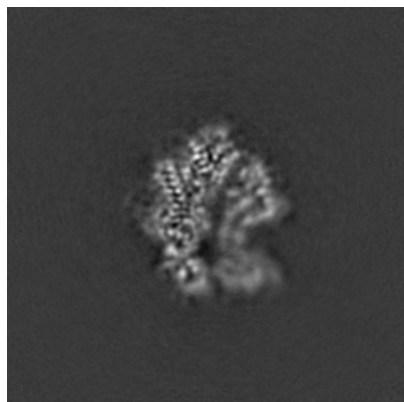


Y Index: 114

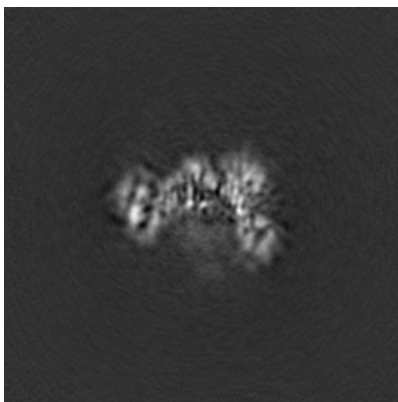


Z Index: 152

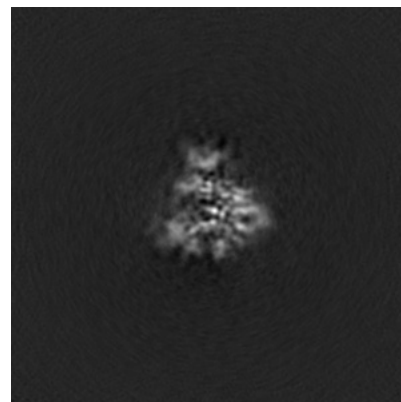
6.3.2 Raw map



X Index: 132



Y Index: 114

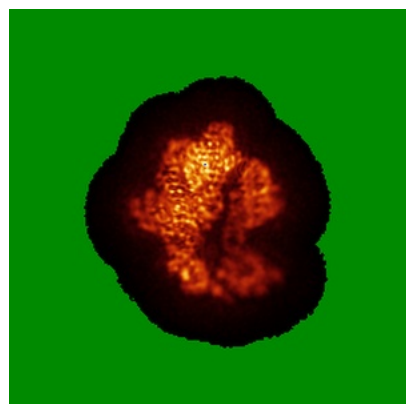


Z Index: 155

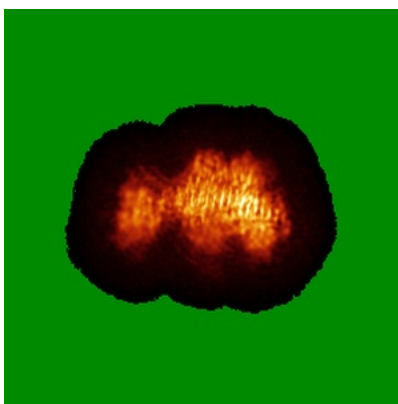
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) ⓘ

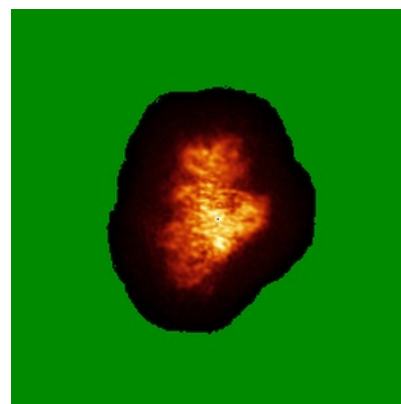
6.4.1 Primary map



X



Y

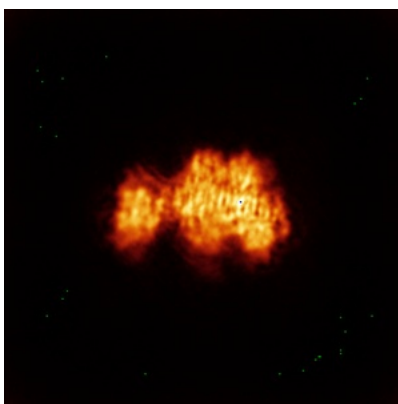


Z

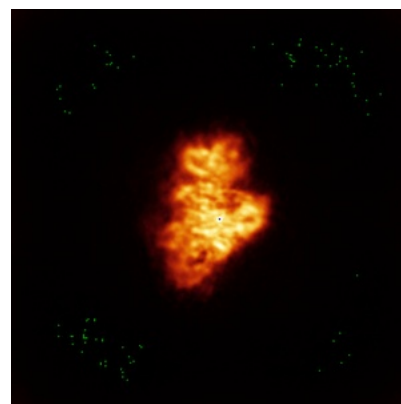
6.4.2 Raw map



X



Y

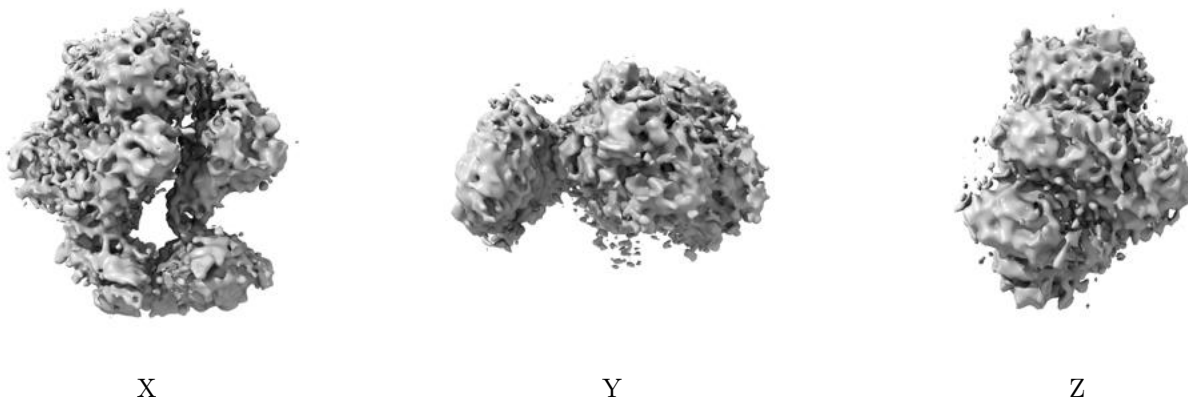


Z

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.

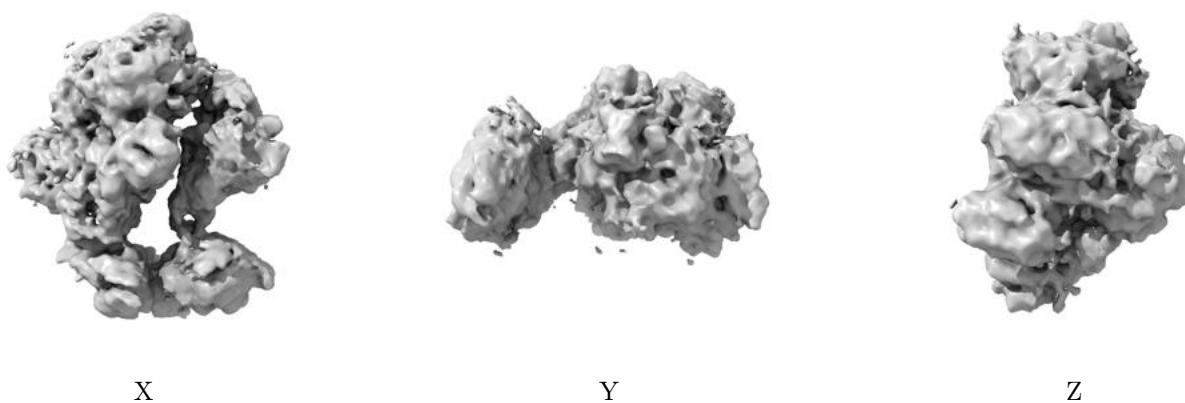
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.224. 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.

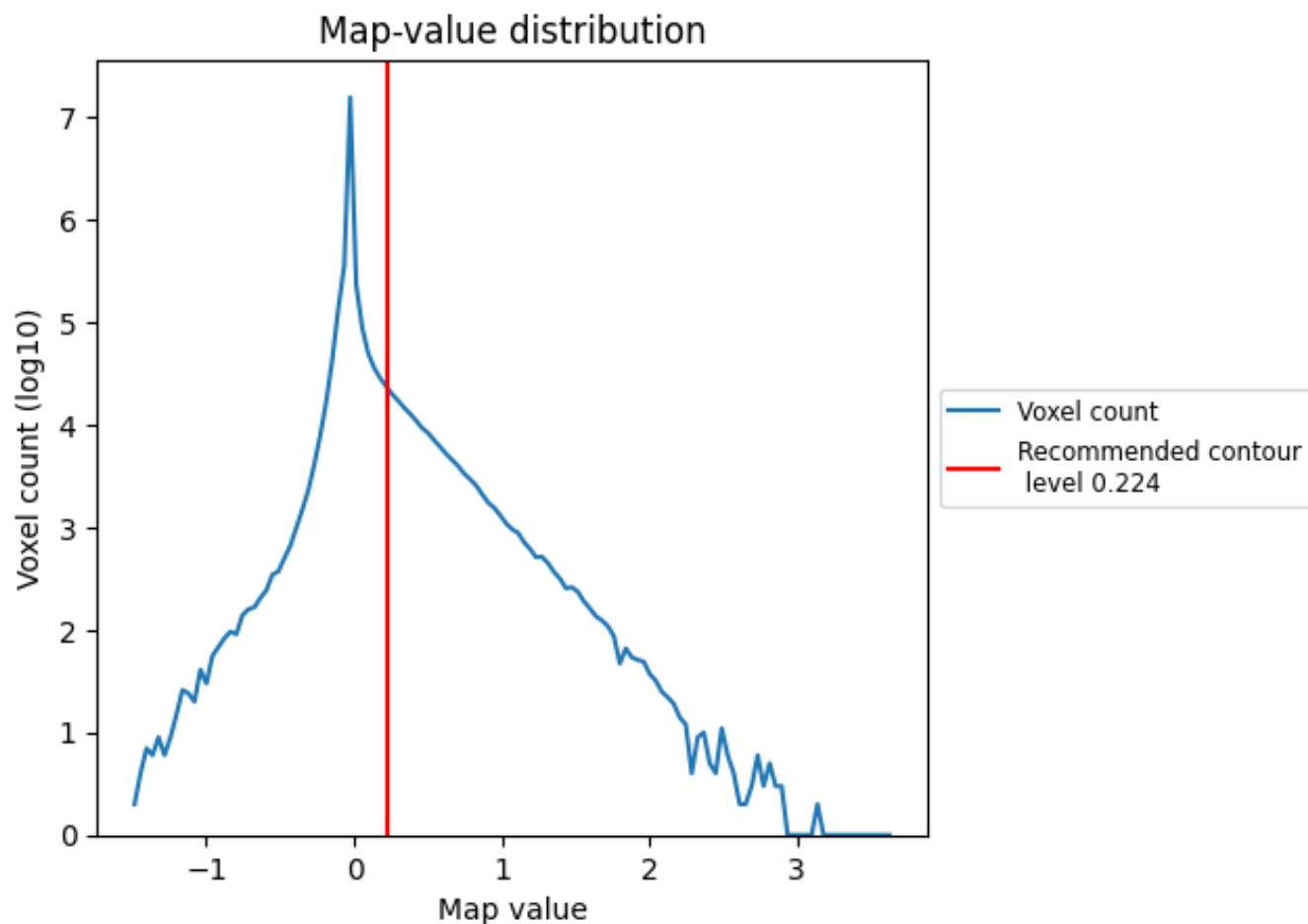
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

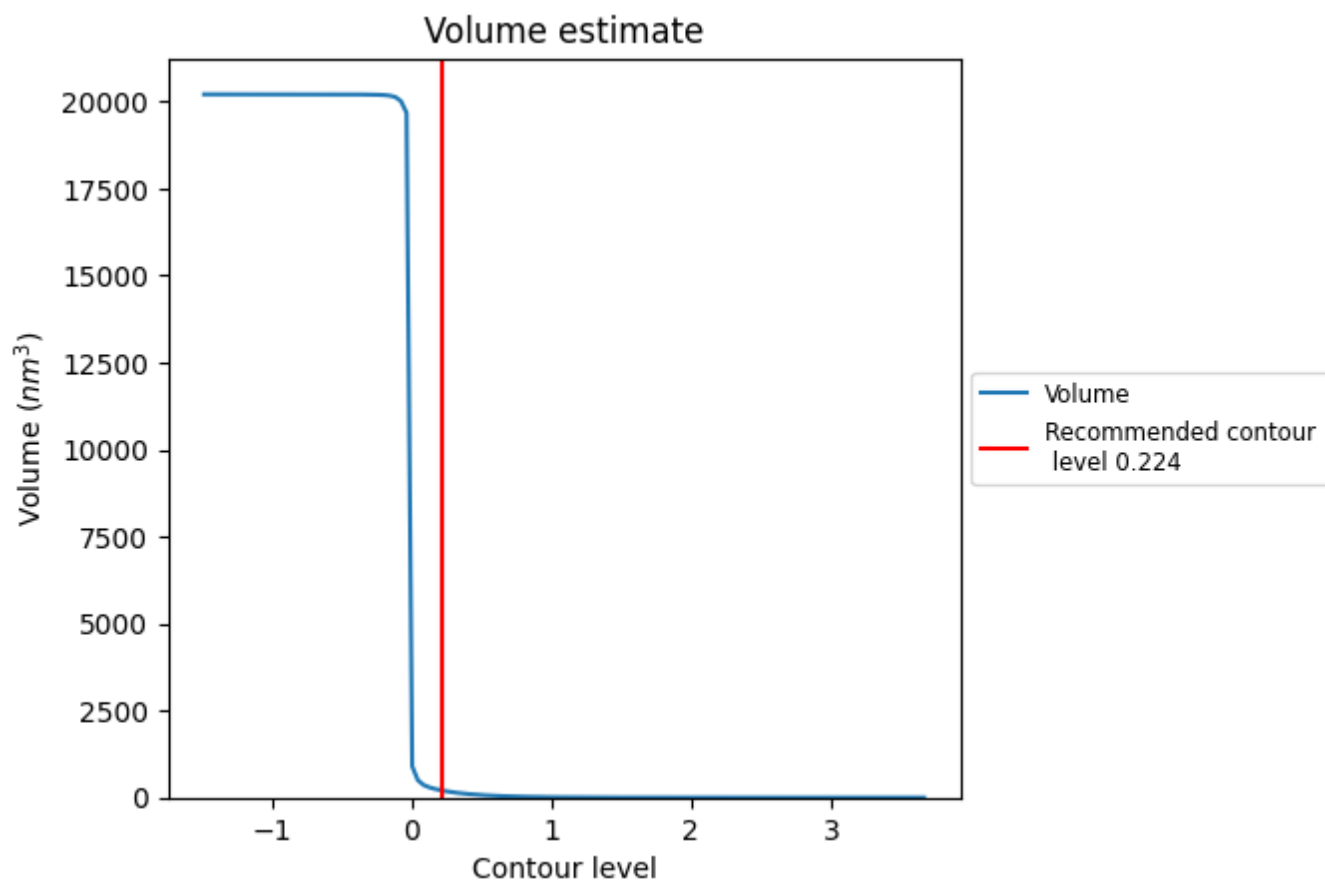
This section contains the results of statistical analysis of the map.

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.

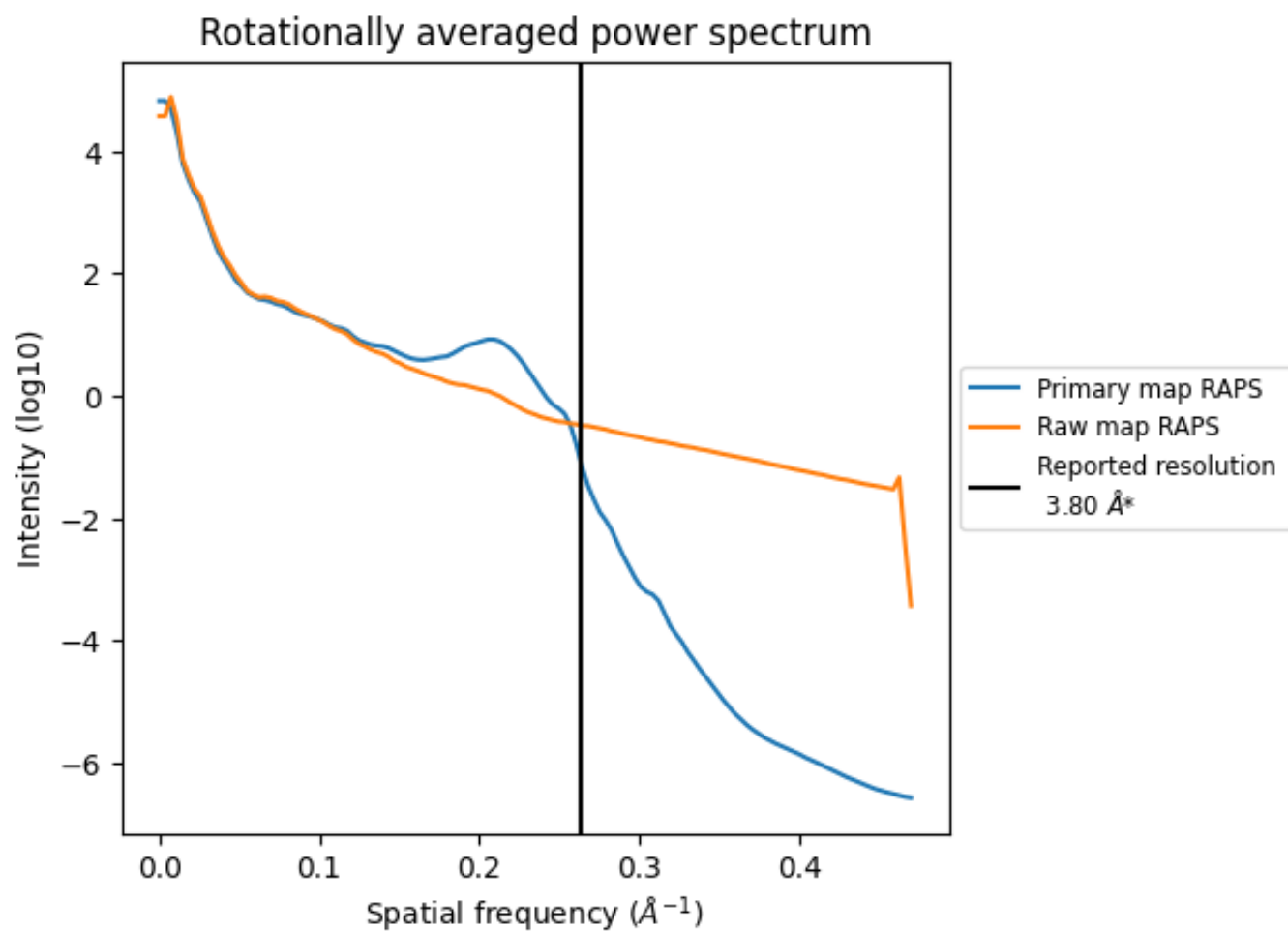
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 199 nm^3 ; this corresponds to an approximate mass of 179 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 ⓘ

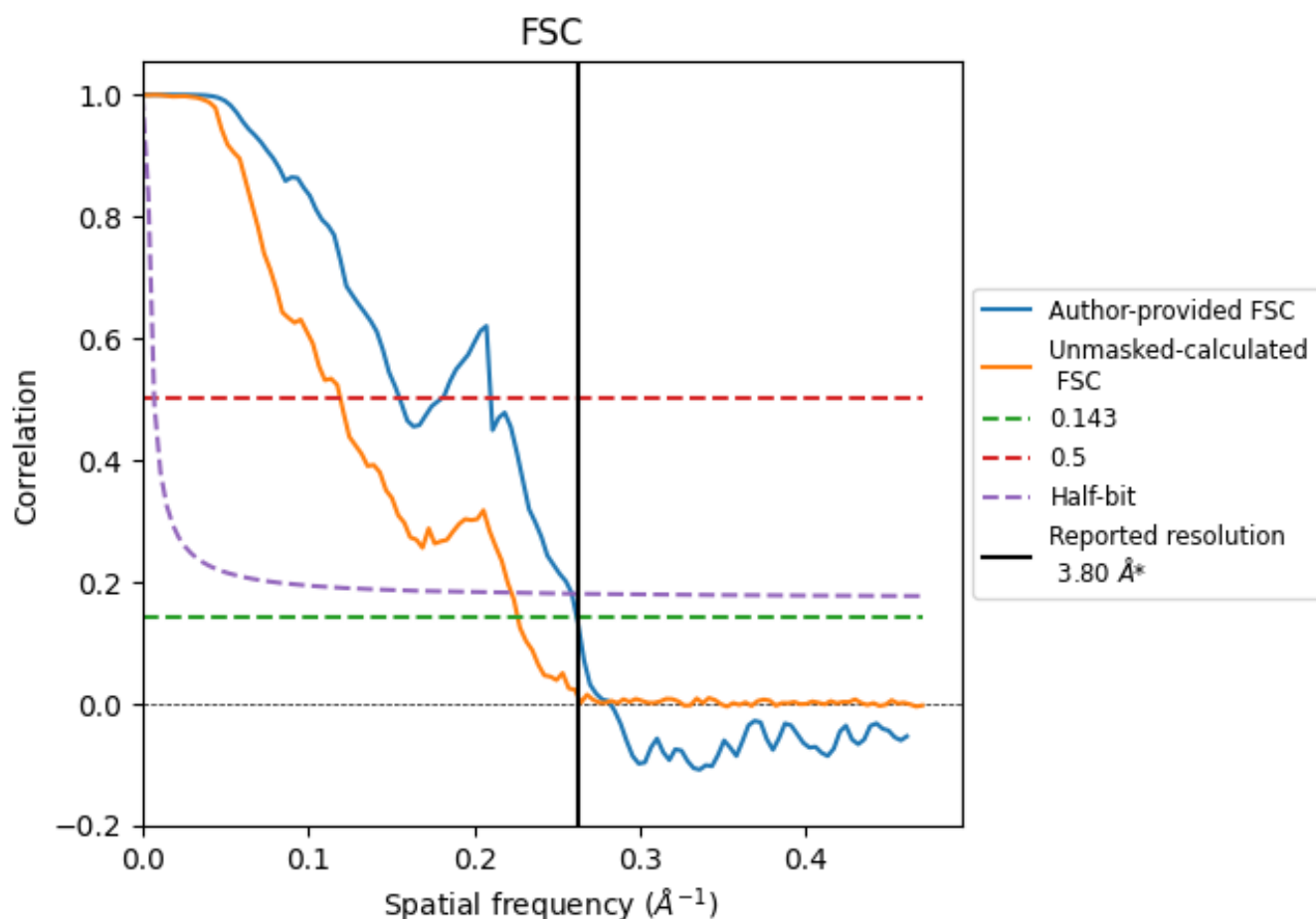


*Reported resolution corresponds to spatial frequency of 0.263 Å⁻¹

8 Fourier-Shell correlation [i](#)

Fourier-Shell Correlation (FSC) is the most commonly used method to estimate the resolution of single-particle and subtomogram-averaged maps. The shape of the curve depends on the imposed symmetry, mask and whether or not the two 3D reconstructions used were processed from a common reference. The reported resolution is shown as a black line. A curve is displayed for the half-bit criterion in addition to lines showing the 0.143 gold standard cut-off and 0.5 cut-off.

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.263 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

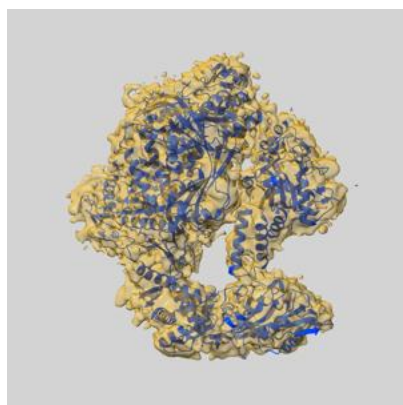
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.80	-	-
Author-provided FSC curve	3.82	6.44	3.86
Unmasked-calculated*	4.42	8.38	4.49

*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 4.42 differs from the reported value 3.8 by more than 10 %

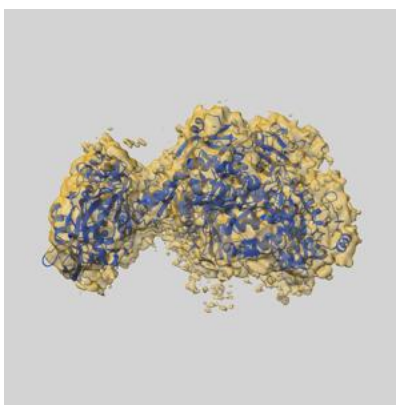
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-18134 and PDB model 8Q3R. Per-residue inclusion information can be found in section [3](#) on page [5](#).

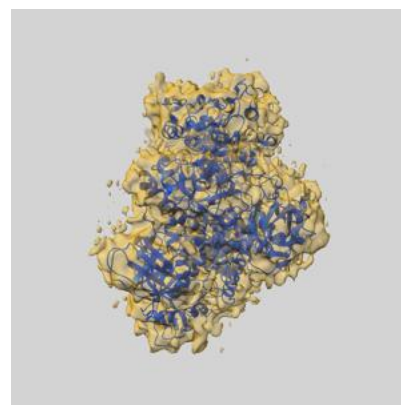
9.1 Map-model overlay [i](#)



X



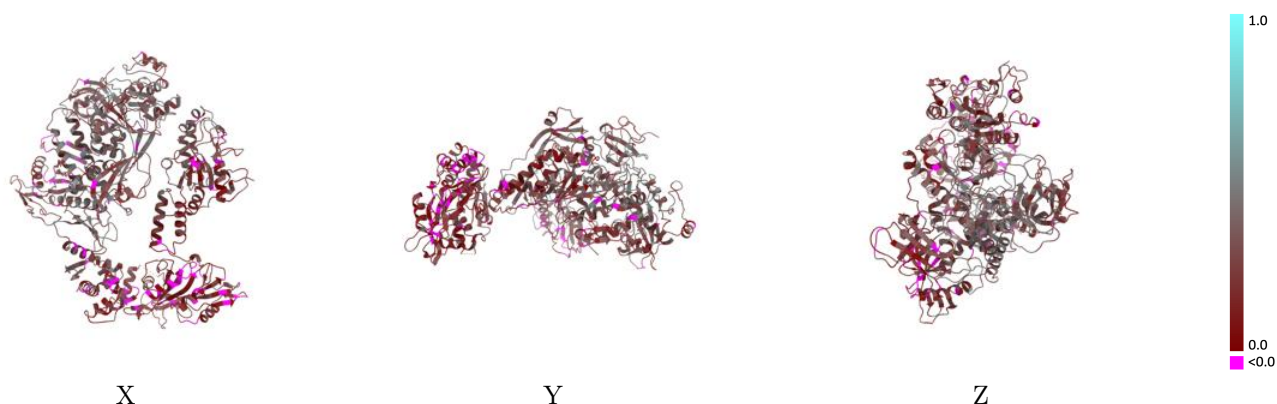
Y



Z

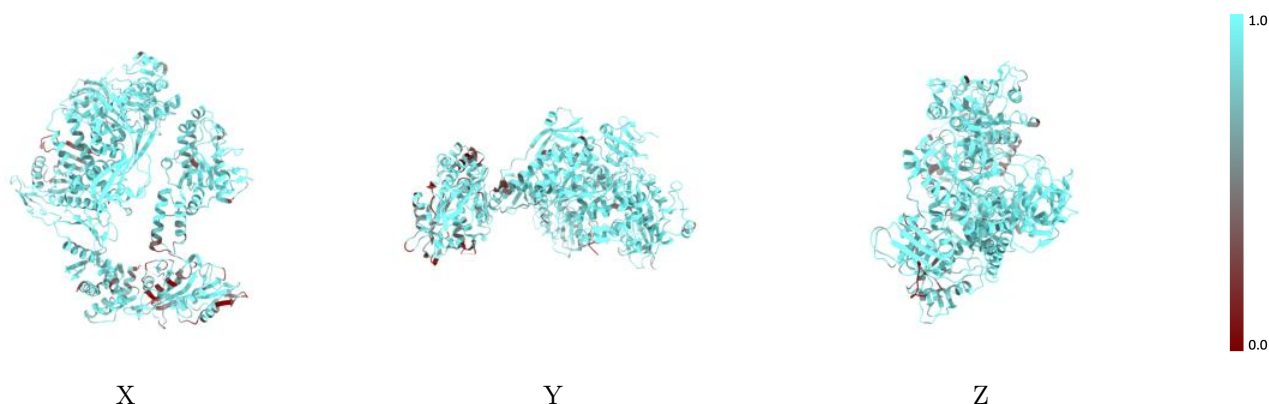
The images above show the 3D surface view of the map at the recommended contour level 0.224 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



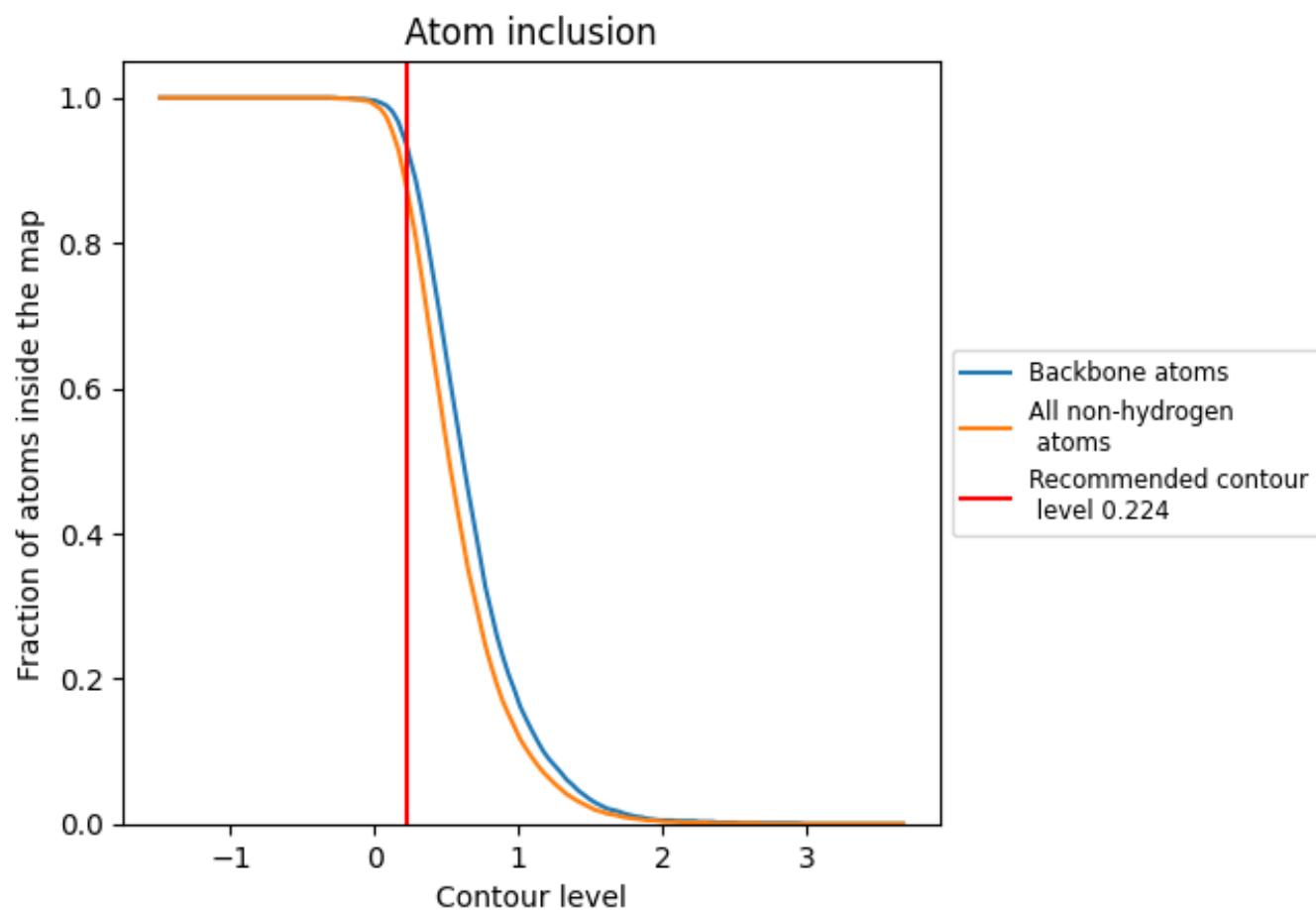
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.224).

9.4 Atom inclusion [i](#)



At the recommended contour level, 93% of all backbone atoms, 88% 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 (0.224) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8750	0.2600
A	0.7800	0.1560
D	0.9250	0.2520
E	0.9110	0.3160

