Full wwPDB EM Validation Report

Mar 7, 2026 – 01:52 AM UTC

PDB ID	:	9FUX / pdb_00009fux
EMDB ID	:	EMD-50781
Title	:	Cryo-EM structure of the Nipah virus polymerase (L) bound to the tetrameric phosphoprotein (P)
Authors	:	Balikci, E.; Gunl, F.; Carrique, L.; Keown, J.R.; Fodor, E.; Grimes, J.M.
Deposited on	:	2024-06-26
Resolution	:	2.49 Å(reported)
Based on initial model	:	.

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 (i) 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 (i)) were used in the production of this report:

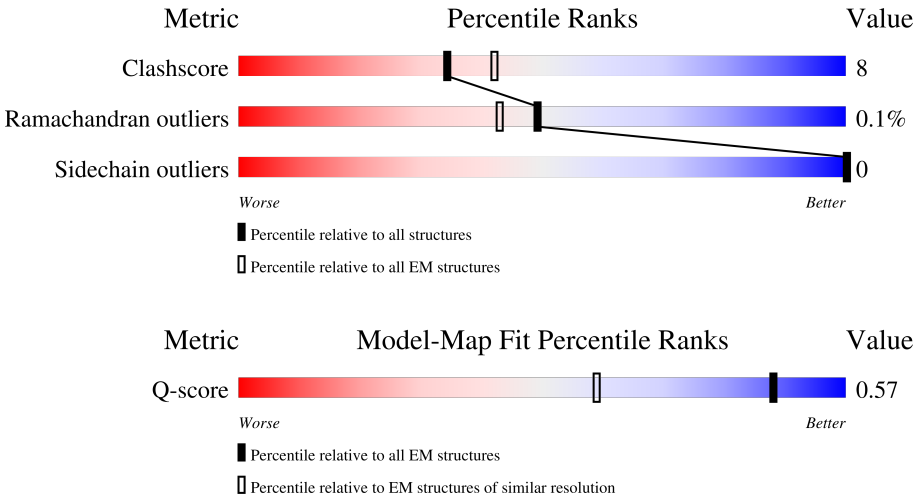
EMDB validation analysis	:	0.0.1.dev132
MolProbity	:	4-5-2 with Phenix2.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics	:	202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ	:	1.9.13
Ideal geometry (proteins)	:	Engh & 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 2.49 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	6237 (2.00 - 2.99)

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 <=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	2243	46% 12% 43%
2	C	708	12% 84%
2	D	708	13% 85%
2	E	708	21% 7% 72%

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Mol	Chain	Length	Quality of chain
2	F	708	11% · 85%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 14556 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 RNA-directed RNA polymerase L.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	1286	Total	C	N	O	S	0	0
			10357	6599	1769	1922	67		

- Molecule 2 is a protein called Phosphoprotein.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	C	116	Total	C	N	O	S	1	0
			927	580	160	181	6		
2	D	108	Total	C	N	O	S	0	0
			861	545	143	166	7		
2	E	198	Total	C	N	O	S	0	0
			1591	999	265	319	8		
2	F	103	Total	C	N	O	S	0	0
			818	516	135	160	7		

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		AltConf
3	A	2	Total	Zn	0
			2	2	

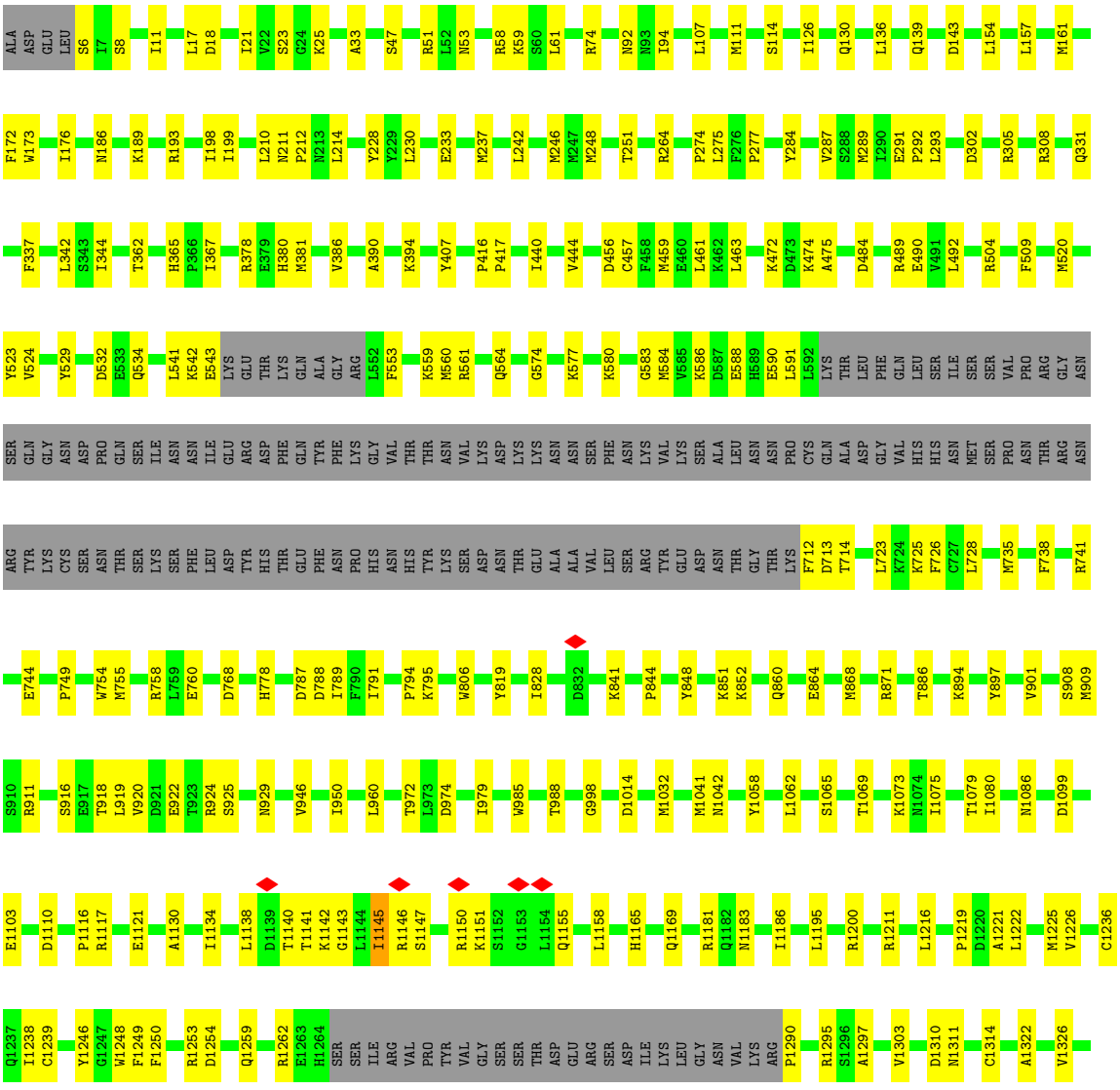
3 Residue-property plots (i)

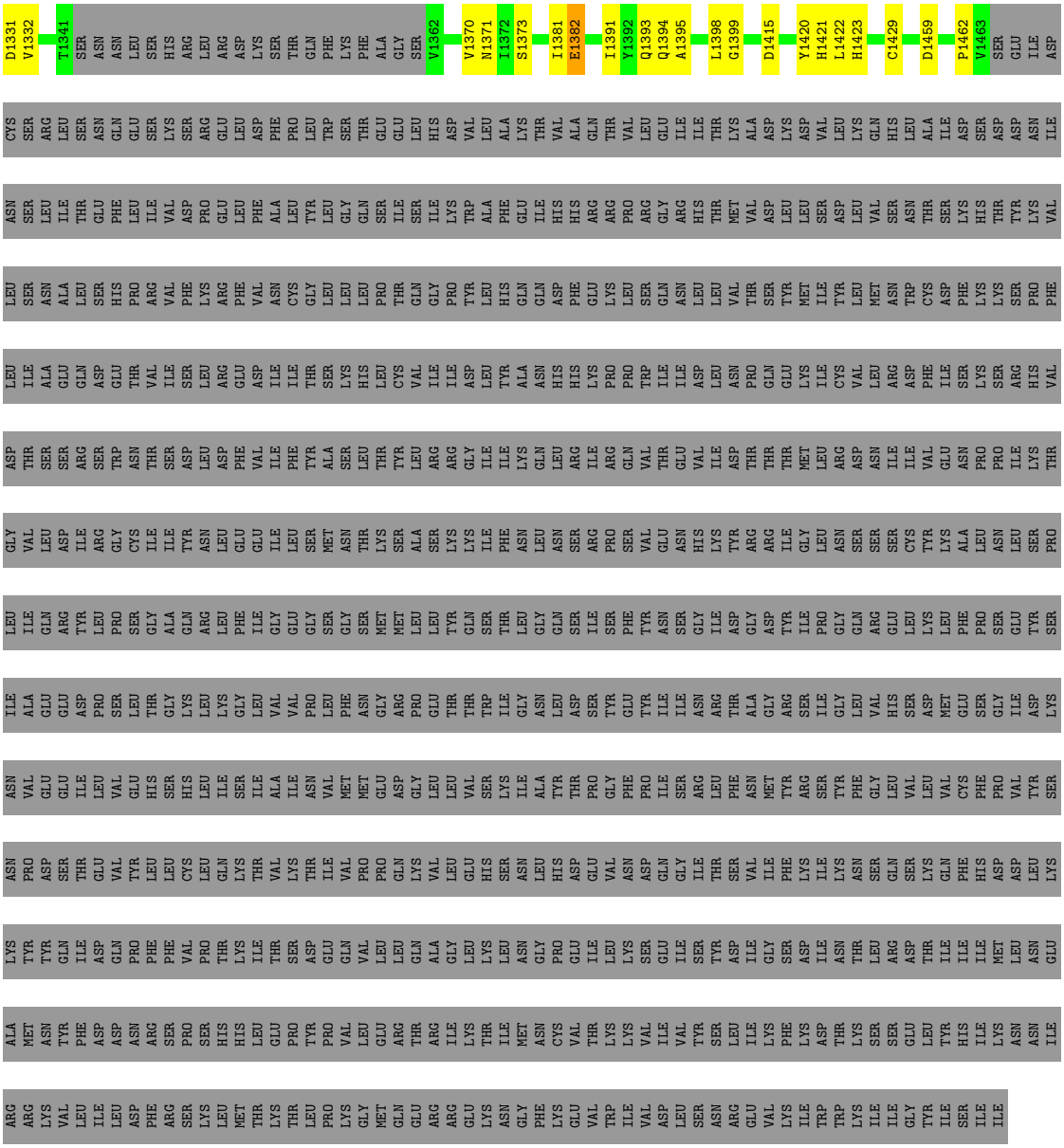
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion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

Chain A:



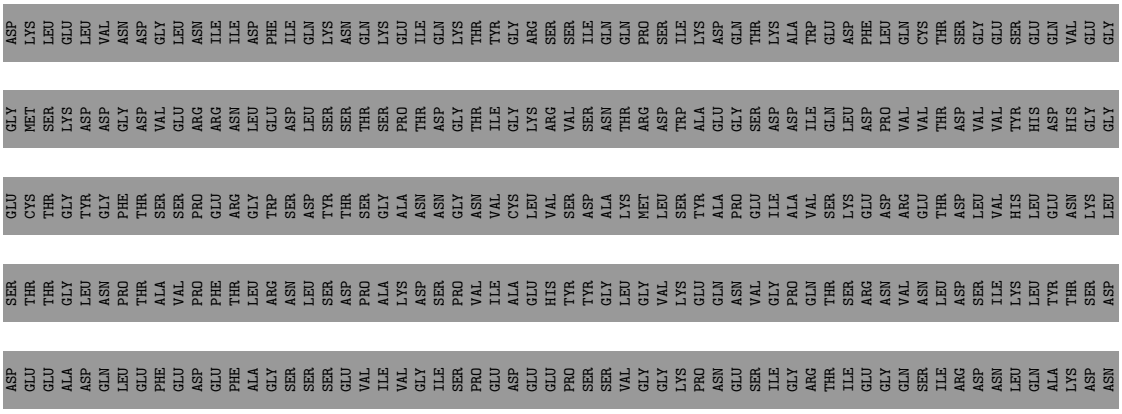


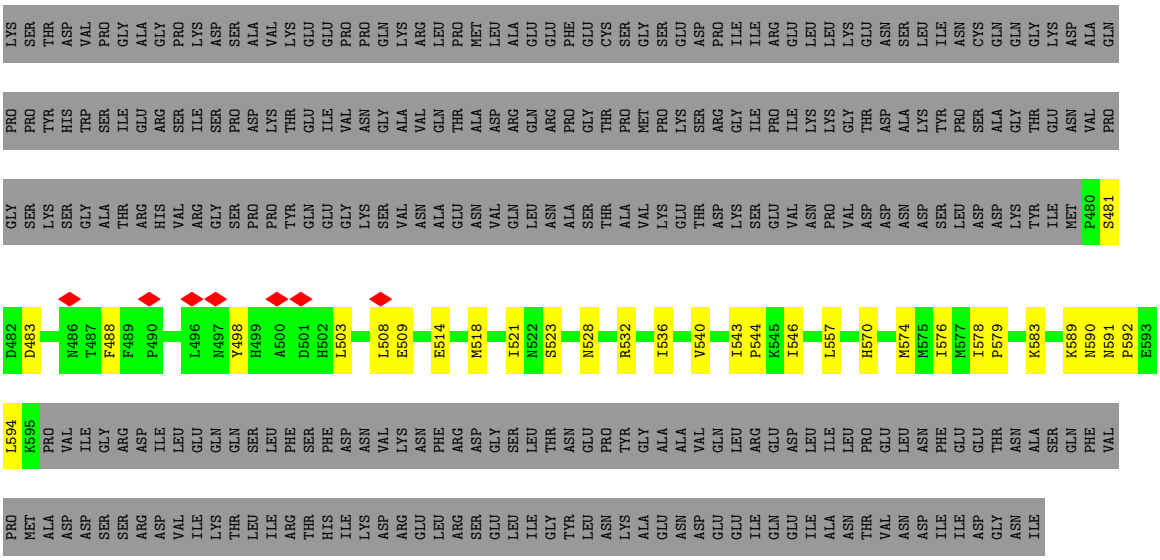


• Molecule 2: Phosphoprotein

Chain C:



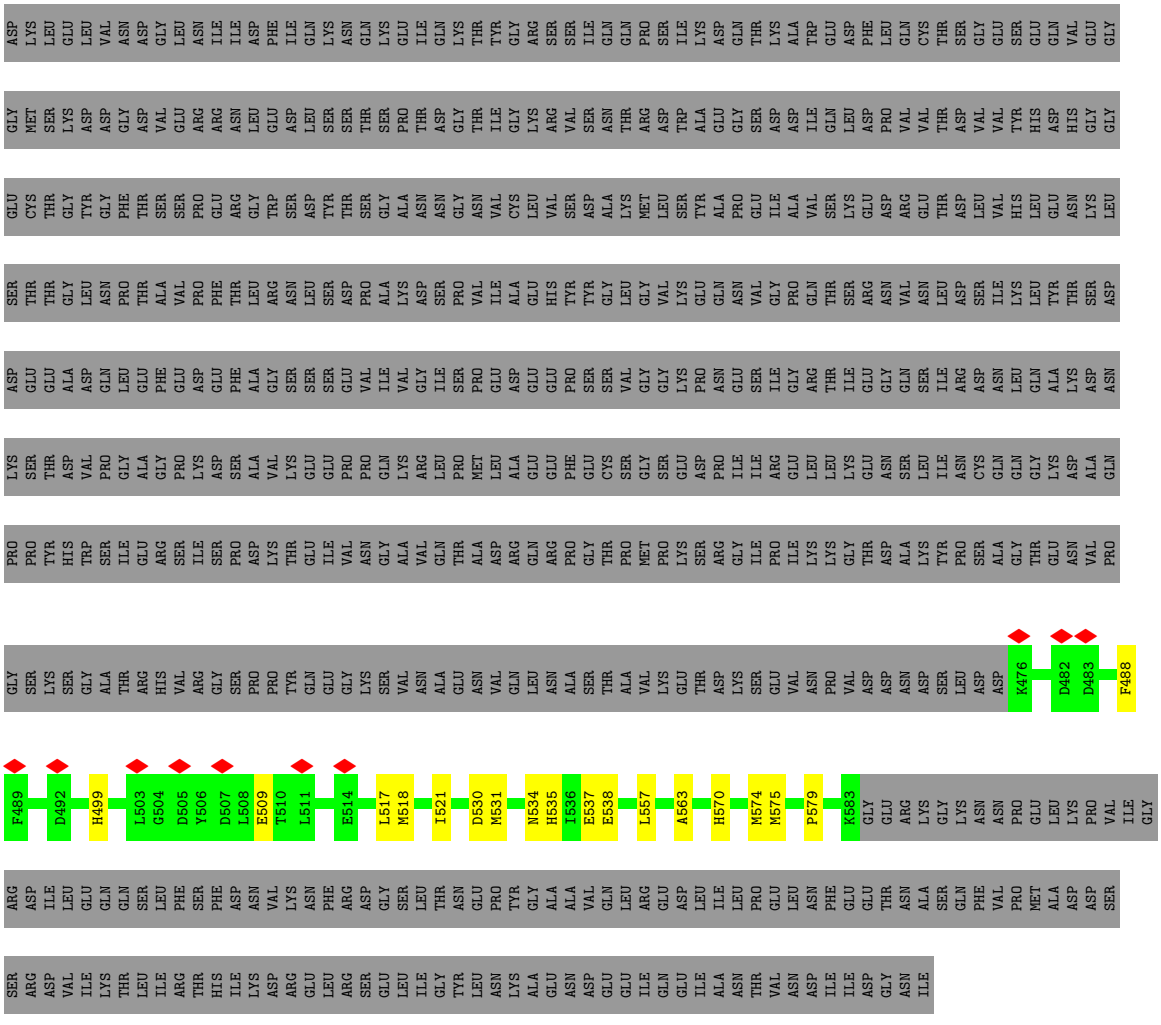




• Molecule 2: Phosphoprotein

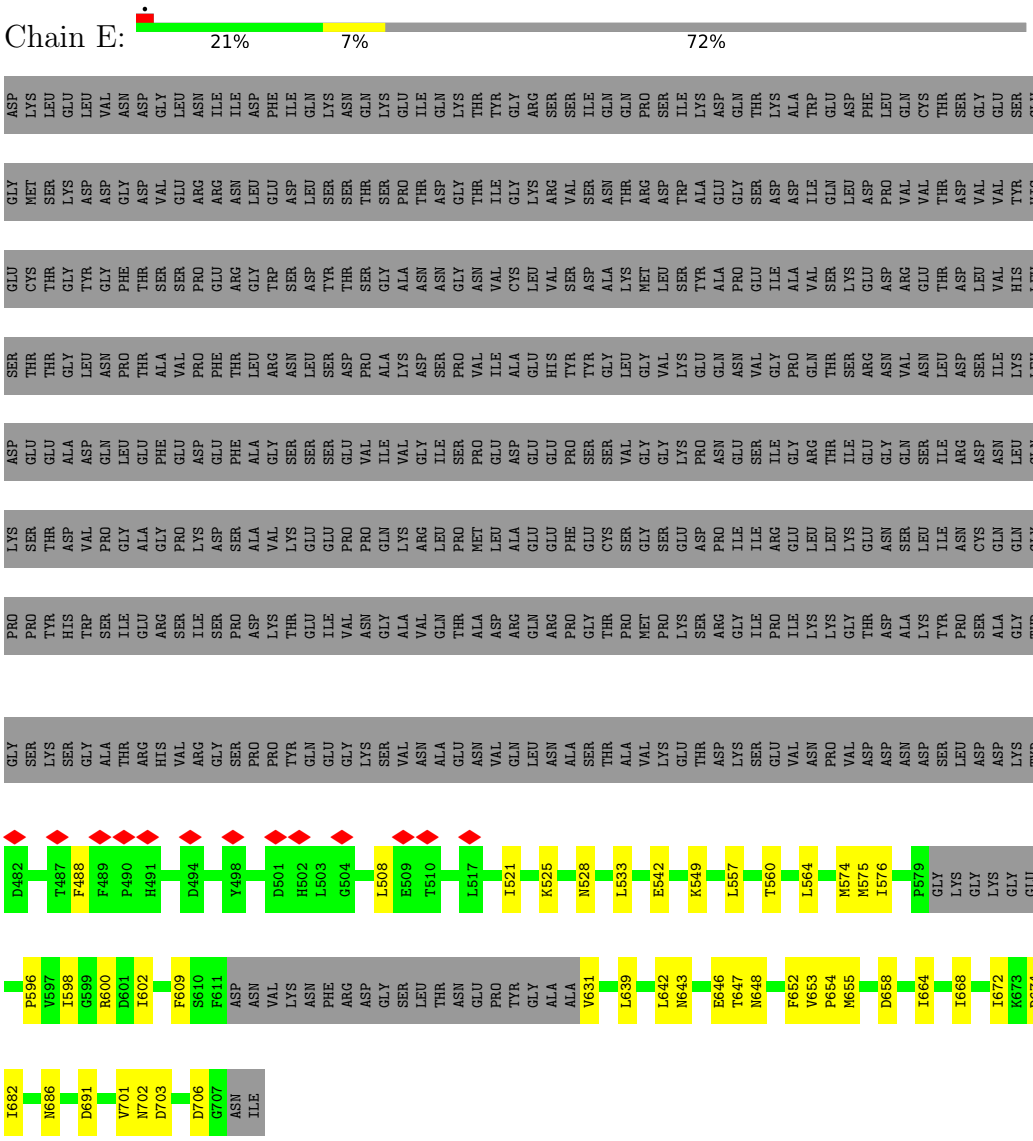
Chain D:





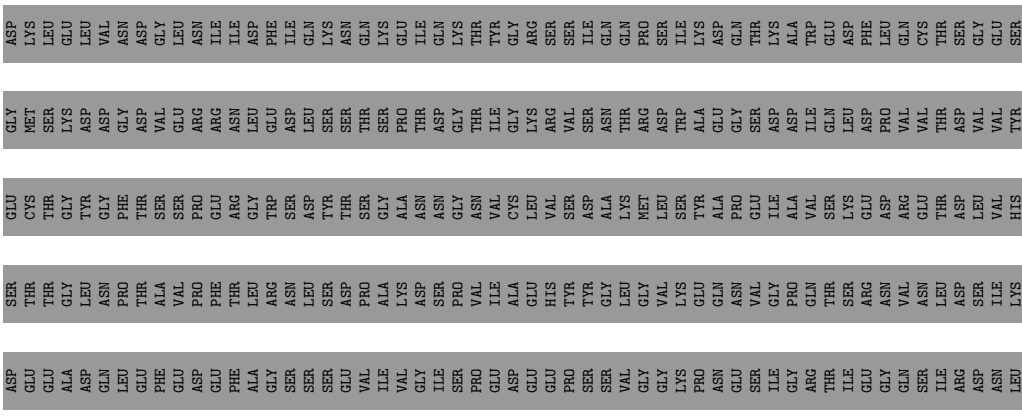
• Molecule 2: Phosphoprotein

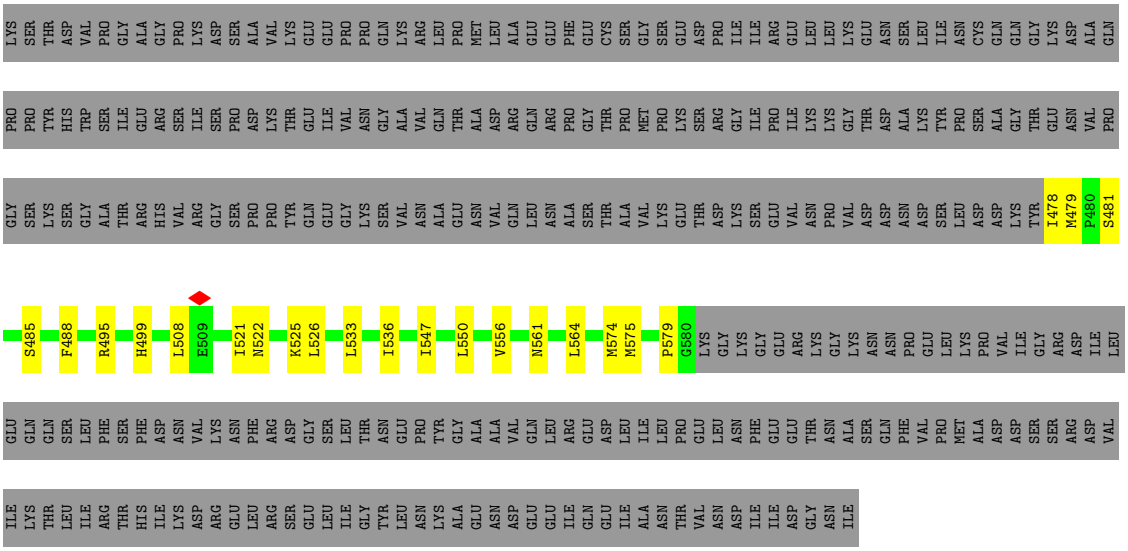
Chain E:



• Molecule 2: Phosphoprotein

Chain F:




S485
F488
R495
H499
L508
E509
I521
N522
K525
L526
L533
I536
I547
L550
V556
N561
L564
M574
M575
P579
G580
I478
M479
P480
S481

4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	490675	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/Å^2$)	50	Depositor
Minimum defocus (nm)	1400	Depositor
Maximum defocus (nm)	2400	Depositor
Magnification	Not provided	
Image detector	TFS FALCON 4i (4k x 4k)	Depositor
Maximum map value	66.063	Depositor
Minimum map value	-54.750	Depositor
Average map value	-0.001	Depositor
Map value standard deviation	0.908	Depositor
Recommended contour level	3	Depositor
Map size (Å)	373.9136, 373.9136, 373.9136	wwPDB
Map dimensions	512, 512, 512	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.7303, 0.7303, 0.7303	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZN

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	0.26	0/10571	0.40	0/14294
2	C	0.19	0/941	0.30	0/1266
2	D	0.16	0/875	0.27	0/1180
2	E	0.17	0/1611	0.34	0/2176
2	F	0.16	0/831	0.33	0/1124
All	All	0.23	0/14829	0.38	0/20040

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

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	10357	0	10379	167	0
2	C	927	0	931	33	0
2	D	861	0	868	22	0
2	E	1591	0	1582	39	0
2	F	818	0	816	20	0
3	A	2	0	0	0	0
All	All	14556	0	14576	237	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 8.

All (237) 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:C:536:ILE:HD11	2:D:537:GLU:HG2	1.51	0.93
1:A:291:GLU:HG2	1:A:292:PRO:HD3	1.53	0.90
1:A:474:LYS:HD3	1:A:561:ARG:HH21	1.43	0.84
1:A:916:SER:OG	1:A:929:ASN:OD1	1.99	0.79
1:A:1147:SER:HA	1:A:1151:LYS:HD2	1.64	0.79
1:A:381:MET:HE3	1:A:794:PRO:HB2	1.66	0.78
1:A:1145:ILE:HG22	1:A:1146:ARG:H	1.49	0.77
1:A:139:GLN:NE2	1:A:143:ASP:OD2	2.19	0.75
1:A:58:ARG:HH12	1:A:490:GLU:HG3	1.53	0.74
2:C:591:ASN:HA	2:C:594:LEU:HB3	1.70	0.72
1:A:474:LYS:HD3	1:A:561:ARG:NH2	2.05	0.71
1:A:1381:ILE:HG22	1:A:1382:GLU:HG2	1.74	0.69
2:C:523:SER:OG	2:F:525:LYS:NZ	2.26	0.69
2:E:631:VAL:HG11	2:F:579:PRO:HG2	1.75	0.68
2:E:682:ILE:O	2:E:686:ASN:ND2	2.24	0.67
1:A:308:ARG:NH1	2:E:702:ASN:OD1	2.27	0.67
1:A:864:GLU:HG2	2:E:642:LEU:HD12	1.77	0.67
2:C:540:VAL:HG21	2:F:536:ILE:HG13	1.77	0.67
1:A:1326:VAL:HG12	1:A:1459:ASP:HB3	1.77	0.66
2:F:533:LEU:HA	2:F:536:ILE:HG22	1.77	0.65
1:A:1080:ILE:HD11	1:A:1138:LEU:HD11	1.79	0.64
1:A:524:VAL:HA	1:A:529:TYR:CD1	2.32	0.64
1:A:1311:ASN:H	1:A:1314:CYS:HB2	1.63	0.64
1:A:211:ASN:HB2	1:A:212:PRO:HD2	1.80	0.64
2:E:542:GLU:HB3	2:F:547:ILE:HD11	1.81	0.63
1:A:868:MET:SD	1:A:871:ARG:NH2	2.72	0.62
1:A:284:TYR:HA	1:A:287:VAL:HG22	1.81	0.62
1:A:1219:PRO:HB3	1:A:1429:CYS:HB3	1.82	0.62
1:A:749:PRO:HG3	2:C:589:LYS:HE3	1.80	0.62
2:C:536:ILE:HG22	2:F:536:ILE:HD11	1.82	0.61
1:A:1249:PHE:HB2	1:A:1420:TYR:HB3	1.82	0.61
1:A:186:ASN:OD1	1:A:189:LYS:NZ	2.34	0.60
1:A:725:LYS:HG2	1:A:728:LEU:HD12	1.84	0.60
1:A:509:PHE:O	1:A:1086:ASN:ND2	2.32	0.60
2:C:576:ILE:HD11	2:C:583:LYS:HD3	1.83	0.60
1:A:17:LEU:HD22	1:A:237:MET:HB2	1.84	0.59
2:F:485:SER:HA	2:F:488:PHE:HD2	1.67	0.59
1:A:23:SER:HA	1:A:367:ILE:HD11	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:308:ARG:NH2	2:E:706:ASP:OD1	2.35	0.59
1:A:47:SER:O	1:A:51:ARG:HG3	2.03	0.59
1:A:136:LEU:O	1:A:1420:TYR:OH	2.20	0.58
1:A:1014:ASP:OD1	1:A:1200:ARG:NH2	2.36	0.58
1:A:248:MET:HE1	1:A:362:THR:HG21	1.85	0.58
1:A:744:GLU:OE2	2:C:583:LYS:NZ	2.25	0.58
1:A:755:MET:HE1	1:A:806:TRP:HZ2	1.69	0.57
1:A:1042:ASN:O	1:A:1181:ARG:NH1	2.37	0.57
1:A:788:ASP:OD1	1:A:789:ILE:N	2.31	0.57
1:A:1226:VAL:HG13	1:A:1238:ILE:HD13	1.86	0.56
1:A:972:THR:OG1	1:A:974:ASP:OD1	2.18	0.56
1:A:1140:THR:HG23	1:A:1142:LYS:H	1.71	0.56
2:E:681:LEU:HD13	2:E:701:VAL:HG12	1.88	0.56
1:A:475:ALA:HB2	1:A:504:ARG:NH1	2.21	0.55
1:A:1145:ILE:HG22	1:A:1146:ARG:N	2.19	0.55
1:A:924:ARG:NH1	1:A:998:GLY:O	2.25	0.55
2:C:532:ARG:NH1	2:D:530:ASP:OD1	2.39	0.55
1:A:1099:ASP:O	1:A:1103:GLU:HG3	2.06	0.54
1:A:1236:CYS:HB3	1:A:1239:CYS:HB2	1.88	0.54
1:A:541:LEU:O	1:A:543:GLU:N	2.40	0.54
2:D:488:PHE:HZ	2:E:521:ILE:HG23	1.73	0.54
2:D:579:PRO:HG2	2:E:602:ILE:HD11	1.90	0.54
1:A:860:GLN:O	1:A:864:GLU:HG3	2.08	0.54
1:A:1146:ARG:HB3	1:A:1150:ARG:HB3	1.90	0.54
1:A:1116:PRO:HG2	1:A:1216:LEU:HD11	1.90	0.53
1:A:560:MET:HE3	1:A:560:MET:HA	1.90	0.53
1:A:580:LYS:HG2	1:A:584:MET:HB2	1.91	0.52
2:C:532:ARG:HH21	2:D:537:GLU:CD	2.17	0.52
2:E:525:LYS:HA	2:E:528:ASN:ND2	2.24	0.52
2:E:668:ILE:HG23	2:E:672:ILE:HG13	1.92	0.52
1:A:561:ARG:NH1	1:A:564:GLN:OE1	2.43	0.52
1:A:417:PRO:HB2	1:A:456:ASP:HB3	1.91	0.52
1:A:440:ILE:O	1:A:444:VAL:HG22	2.09	0.52
1:A:529:TYR:O	1:A:758:ARG:NH1	2.42	0.51
1:A:848:TYR:O	1:A:852:LYS:HG2	2.09	0.51
1:A:1041:MET:HB3	1:A:1186:ILE:HD13	1.93	0.51
2:D:570:HIS:CE1	2:E:575:MET:HE1	2.46	0.51
1:A:251:THR:HG21	1:A:909:MET:HG3	1.93	0.51
1:A:979:ILE:HD12	1:A:985:TRP:HB3	1.93	0.51
2:E:674:ASP:HB3	2:E:676:GLU:HG2	1.92	0.51
1:A:25:LYS:NZ	1:A:233:GLU:OE1	2.35	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1248:TRP:HB3	1:A:1373:SER:HB3	1.91	0.51
1:A:472:LYS:HD2	1:A:474:LYS:HE2	1.92	0.51
1:A:574:GLY:O	1:A:577:LYS:HG2	2.11	0.51
1:A:583:GLY:HA2	1:A:828:ILE:HD12	1.92	0.51
1:A:713:ASP:OD1	1:A:714:THR:N	2.44	0.50
1:A:1221:ALA:HB1	1:A:1398:LEU:HD22	1.93	0.50
1:A:248:MET:HE1	1:A:362:THR:CG2	2.41	0.50
1:A:264:ARG:NH1	1:A:344:ILE:O	2.41	0.50
1:A:1138:LEU:HD23	1:A:1138:LEU:O	2.11	0.50
2:C:488:PHE:HE2	2:D:521:ILE:HG12	1.76	0.50
1:A:461:LEU:HB3	1:A:463:LEU:HG	1.93	0.50
1:A:542:LYS:HB2	1:A:553:PHE:HE2	1.75	0.50
1:A:157:LEU:HD11	1:A:950:ILE:HG13	1.94	0.50
1:A:33:ALA:O	1:A:74:ARG:NH2	2.44	0.50
1:A:172:PHE:O	1:A:176:ILE:HG12	2.12	0.50
1:A:1248:TRP:CH2	1:A:1250:PHE:HB2	2.46	0.49
2:E:647:THR:HG1	2:E:653:VAL:H	1.60	0.49
2:C:557:LEU:HG	2:D:557:LEU:HD21	1.94	0.49
1:A:871:ARG:HG3	2:E:639:LEU:HB2	1.95	0.49
1:A:1130:ALA:O	1:A:1134:ILE:HD12	2.13	0.49
1:A:18:ASP:OD1	1:A:18:ASP:N	2.43	0.48
1:A:378:ARG:NH2	1:A:787:ASP:O	2.46	0.48
1:A:53:ASN:ND2	1:A:490:GLU:O	2.46	0.48
1:A:897:TYR:OH	1:A:1259:GLN:OE1	2.11	0.48
1:A:908:SER:HA	1:A:911:ARG:NH1	2.29	0.48
2:E:574:MET:SD	2:F:575:MET:HG2	2.53	0.48
2:C:503:LEU:HD12	2:D:517:LEU:HG	1.96	0.47
2:D:574:MET:SD	2:E:598:ILE:HG23	2.53	0.47
2:D:531:MET:HA	2:D:534:ASN:ND2	2.30	0.47
1:A:293:LEU:HD22	1:A:342:LEU:HD21	1.96	0.47
2:E:643:ASN:HB3	2:E:646:GLU:HB3	1.96	0.47
1:A:291:GLU:OE2	1:A:291:GLU:N	2.45	0.47
1:A:11:ILE:HD11	1:A:919:LEU:HD11	1.97	0.47
1:A:6:SER:O	1:A:8:SER:N	2.45	0.47
1:A:289:MET:HE3	1:A:289:MET:HB3	1.82	0.47
1:A:1058:TYR:OH	1:A:1117:ARG:NH1	2.48	0.47
1:A:1239:CYS:SG	1:A:1423:HIS:CE1	3.08	0.47
2:C:528:ASN:O	2:C:532:ARG:HG3	2.15	0.47
2:C:483:ASP:OD2	2:D:499:HIS:NE2	2.48	0.46
2:C:498:TYR:OH	2:F:478:ILE:HG13	2.15	0.46
1:A:61:LEU:HD13	1:A:193:ARG:HE	1.79	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:481:SER:HA	2:D:518:MET:HG2	1.95	0.46
1:A:922:GLU:HG3	1:A:925:SER:HB2	1.97	0.46
2:C:578:ILE:HD12	2:F:574:MET:HG2	1.98	0.46
2:C:540:VAL:O	2:C:543:ILE:HG22	2.16	0.46
1:A:489:ARG:NH1	1:A:492:LEU:O	2.49	0.46
1:A:735:MET:HE3	1:A:738:PHE:HB2	1.97	0.46
2:C:543:ILE:HG13	2:C:546:ILE:HD12	1.97	0.46
1:A:161:MET:HE3	1:A:246:MET:HE3	1.98	0.46
1:A:1290:PRO:O	1:A:1295:ARG:NH1	2.49	0.46
1:A:1310:ASP:OD1	1:A:1310:ASP:N	2.48	0.46
2:C:570:HIS:O	2:C:574:MET:HG3	2.16	0.45
1:A:1141:THR:C	1:A:1143:GLY:H	2.24	0.45
2:C:508:LEU:HD21	2:D:509:GLU:HA	1.98	0.45
1:A:331:GLN:HE21	2:E:658:ASP:HB3	1.81	0.45
1:A:456:ASP:OD1	1:A:457:CYS:N	2.46	0.45
1:A:1225:MET:HE3	1:A:1395:ALA:HA	1.99	0.45
2:E:549:LYS:HB3	2:F:550:LEU:HD23	1.97	0.45
1:A:305:ARG:NH2	2:E:703:ASP:OD1	2.46	0.45
1:A:92:ASN:C	1:A:94:ILE:H	2.25	0.45
1:A:723:LEU:HG	1:A:726:PHE:CE2	2.52	0.45
1:A:1121:GLU:HG2	1:A:1303:VAL:HG21	1.98	0.45
2:C:543:ILE:HG22	2:C:544:PRO:HD3	1.99	0.45
1:A:712:PHE:HA	1:A:841:LYS:HE2	1.99	0.45
2:F:479:MET:HE1	2:F:481:SER:OG	2.16	0.45
1:A:198:ILE:O	1:A:199:ILE:HD13	2.17	0.45
2:E:652:PHE:CE2	2:E:654:PRO:HG3	2.52	0.45
2:E:557:LEU:HD23	2:E:557:LEU:HA	1.83	0.45
1:A:1326:VAL:HB	1:A:1462:PRO:HG3	2.00	0.44
1:A:583:GLY:HA2	1:A:828:ILE:HB	1.98	0.44
2:D:563:ALA:HB3	2:E:564:LEU:HD13	2.00	0.44
1:A:21:ILE:HD12	1:A:365:HIS:HB2	1.99	0.44
1:A:894:LYS:HD3	1:A:901:VAL:HG11	2.00	0.44
2:E:560:THR:OG1	2:F:561:ASN:OD1	2.22	0.44
1:A:275:LEU:HD11	1:A:337:PHE:CZ	2.53	0.44
1:A:787:ASP:OD1	1:A:788:ASP:N	2.49	0.44
2:E:479:MET:HE3	2:E:480:PRO:HD2	1.99	0.44
1:A:1250:PHE:HB3	1:A:1371:ASN:HB2	1.99	0.44
1:A:1297:ALA:HB1	1:A:1322:ALA:HB1	2.00	0.44
1:A:1062:LEU:HB2	1:A:1065:SER:HB3	1.99	0.44
1:A:1421:HIS:HB3	1:A:1423:HIS:CD2	2.53	0.44
2:D:488:PHE:HE2	2:E:521:ILE:HG12	1.83	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:484:ASP:OD2	1:A:778:HIS:ND1	2.50	0.44
1:A:523:TYR:OH	1:A:532:ASP:OD2	2.32	0.44
1:A:94:ILE:HG21	1:A:228:TYR:OH	2.18	0.44
1:A:1069:THR:O	1:A:1073:LYS:HG2	2.18	0.44
1:A:173:TRP:CD1	1:A:242:LEU:HD12	2.52	0.43
1:A:459:MET:O	2:C:590:ASN:ND2	2.48	0.43
2:C:509:GLU:HG3	2:F:508:LEU:HD21	1.99	0.43
1:A:1399:GLY:HA3	1:A:1422:LEU:HD22	2.00	0.43
1:A:107:LEU:O	1:A:111:MET:HG2	2.17	0.43
1:A:713:ASP:HB2	1:A:844:PRO:HB3	2.00	0.43
2:E:488:PHE:CZ	2:F:521:ILE:HG23	2.53	0.43
2:E:674:ASP:C	2:E:676:GLU:H	2.26	0.43
1:A:1399:GLY:HA2	1:A:1420:TYR:HE2	1.83	0.43
2:E:576:ILE:HD13	2:E:609:PHE:HE1	1.83	0.43
1:A:11:ILE:HD11	1:A:919:LEU:CD1	2.48	0.43
2:D:535:HIS:O	2:D:538:GLU:HG2	2.19	0.43
1:A:198:ILE:HA	1:A:210:LEU:O	2.19	0.43
1:A:289:MET:C	1:A:292:PRO:HD2	2.44	0.42
1:A:1155:GLN:HB3	1:A:1158:LEU:HB3	2.01	0.42
2:E:525:LYS:HA	2:E:528:ASN:HD21	1.84	0.42
1:A:542:LYS:NZ	1:A:768:ASP:OD2	2.52	0.42
1:A:586:LYS:HB3	1:A:588:GLU:OE1	2.19	0.42
2:D:574:MET:SD	2:E:596:PRO:HB2	2.58	0.42
2:F:522:ASN:O	2:F:526:LEU:HD23	2.19	0.42
1:A:274:PRO:O	1:A:277:PRO:HD2	2.20	0.42
1:A:760:GLU:HG2	1:A:795:LYS:HD3	2.02	0.42
1:A:114:SER:OG	1:A:960:LEU:HD12	2.20	0.42
1:A:1331:ASP:OD1	1:A:1332:VAL:N	2.51	0.42
2:C:514:GLU:O	2:C:518:MET:HG3	2.20	0.42
1:A:741:ARG:HA	1:A:741:ARG:HD3	1.76	0.42
1:A:380:HIS:CE1	1:A:725:LYS:HD2	2.55	0.42
1:A:1041:MET:HE2	1:A:1195:LEU:HD22	2.02	0.42
2:E:533:LEU:HD23	2:E:533:LEU:HA	1.85	0.42
2:C:579:PRO:HG2	2:E:600:ARG:CZ	2.50	0.42
1:A:590:GLU:OE1	1:A:591:LEU:HG	2.19	0.42
1:A:988:THR:HA	1:A:1032:MET:HE1	2.01	0.42
2:E:642:LEU:HD22	2:E:648:ASN:HD22	1.85	0.42
1:A:302:ASP:OD1	1:A:886:THR:OG1	2.35	0.42
1:A:1391:ILE:HG22	1:A:1393:GLN:HG2	2.02	0.42
2:C:532:ARG:HH22	2:D:534:ASN:HB3	1.85	0.42
1:A:126:ILE:O	1:A:130:GLN:HG3	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1250:PHE:O	1:A:1370:VAL:HA	2.19	0.41
1:A:1391:ILE:O	1:A:1394:GLN:HG2	2.20	0.41
2:E:654:PRO:HB3	2:E:664:ILE:HD11	2.01	0.41
2:E:674:ASP:O	2:E:675:ARG:HB3	2.20	0.41
2:C:532:ARG:HH22	2:D:534:ASN:CB	2.33	0.41
2:C:591:ASN:N	2:C:592:PRO:HD2	2.35	0.41
2:E:655:MET:O	2:E:691:ASP:HA	2.19	0.41
1:A:59:LYS:HE2	1:A:59:LYS:HB2	1.89	0.41
1:A:444:VAL:HG21	1:A:819:TYR:CE1	2.56	0.41
1:A:407:TYR:HB3	1:A:416:PRO:HG3	2.01	0.41
1:A:386:VAL:HG12	2:D:575:MET:HE2	2.02	0.41
2:C:488:PHE:CZ	2:D:521:ILE:HG23	2.56	0.41
2:C:521:ILE:HG23	2:F:488:PHE:CE2	2.55	0.41
1:A:390:ALA:O	1:A:394:LYS:HG3	2.20	0.41
1:A:1142:LYS:NZ	1:A:1147:SER:O	2.40	0.41
1:A:1253:ARG:HG3	1:A:1415:ASP:C	2.46	0.41
1:A:1254:ASP:O	1:A:1262:ARG:NH2	2.53	0.41
2:F:495:ARG:HG2	2:F:499:HIS:CE1	2.55	0.41
1:A:534:GLN:HA	1:A:534:GLN:OE1	2.21	0.41
1:A:754:TRP:O	1:A:755:MET:HB3	2.21	0.41
1:A:1110:ASP:HA	1:A:1211:ARG:HH12	1.86	0.41
1:A:1165:HIS:O	1:A:1169:GLN:HG2	2.21	0.41
2:C:557:LEU:HD13	2:F:556:VAL:HG21	2.03	0.41
2:F:564:LEU:HD23	2:F:564:LEU:HA	1.84	0.41
1:A:136:LEU:HD11	1:A:1222:LEU:HD22	2.03	0.40
1:A:1183:ASN:HB3	1:A:1186:ILE:HD12	2.01	0.40
1:A:378:ARG:HG3	1:A:791:ILE:HG23	2.02	0.40
1:A:918:THR:HB	1:A:920:VAL:O	2.21	0.40
1:A:1075:ILE:O	1:A:1079:THR:OG1	2.27	0.40
1:A:214:LEU:HD12	1:A:230:LEU:O	2.22	0.40
1:A:520:MET:SD	1:A:559:LYS:HG2	2.62	0.40
2:E:508:LEU:HD23	2:E:508:LEU:HA	1.92	0.40
1:A:154:LEU:HD21	1:A:946:VAL:HG22	2.03	0.40
1:A:713:ASP:OD2	1:A:851:LYS:NZ	2.32	0.40
1:A:1246:TYR:HB3	1:A:1423:HIS:CG	2.57	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	1276/2243 (57%)	1219 (96%)	55 (4%)	2 (0%)	43	63
2	C	115/708 (16%)	108 (94%)	7 (6%)	0	100	100
2	D	106/708 (15%)	106 (100%)	0	0	100	100
2	E	192/708 (27%)	188 (98%)	4 (2%)	0	100	100
2	F	101/708 (14%)	101 (100%)	0	0	100	100
All	All	1790/5075 (35%)	1722 (96%)	66 (4%)	2 (0%)	49	68

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1382	GLU
1	A	1145	ILE

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	1156/2045 (56%)	1156 (100%)	0	100	100
2	C	107/624 (17%)	107 (100%)	0	100	100
2	D	100/624 (16%)	100 (100%)	0	100	100
2	E	185/624 (30%)	185 (100%)	0	100	100
2	F	96/624 (15%)	96 (100%)	0	100	100
All	All	1644/4541 (36%)	1644 (100%)	0	100	100

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

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	38	ASN
1	A	139	GLN
1	A	213	ASN
1	A	266	ASN
1	A	331	GLN
1	A	349	ASN
1	A	351	HIS
1	A	536	ASN
1	A	875	HIS
1	A	951	ASN
1	A	968	ASN
1	A	1011	ASN
1	A	1189	ASN
1	A	1324	GLN
1	A	1327	ASN
1	A	1371	ASN
2	C	502	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)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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.

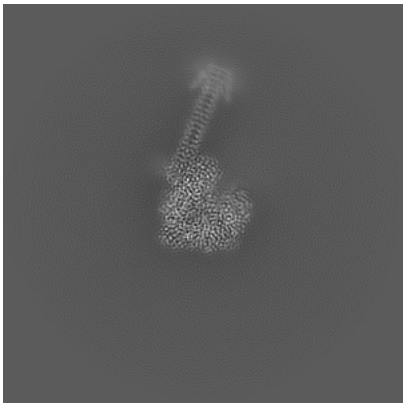
6 Map visualisation

This section contains visualisations of the EMDB entry EMD-50781. These allow visual inspection of the internal detail of the map and identification of artifacts.

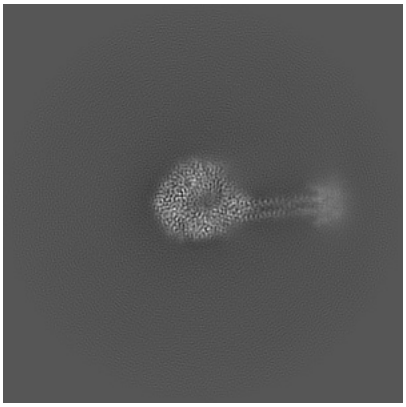
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections

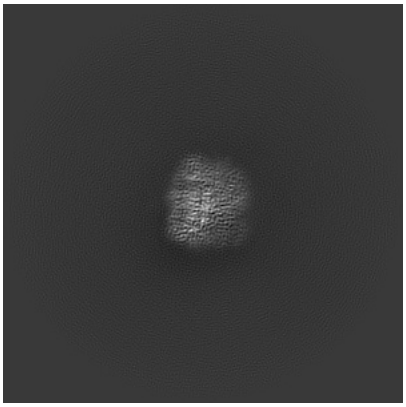
6.1.1 Primary map



X



Y

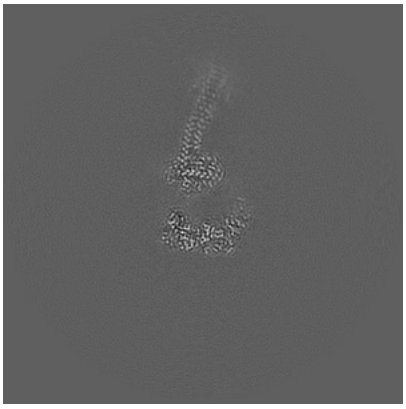


Z

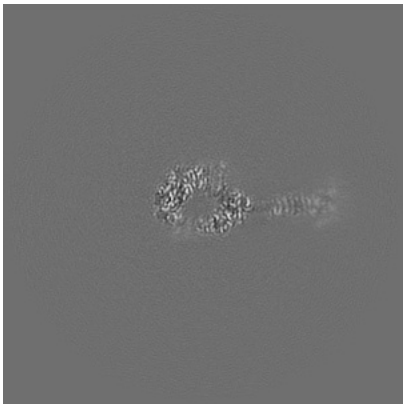
The images above show the map projected in three orthogonal directions.

6.2 Central slices

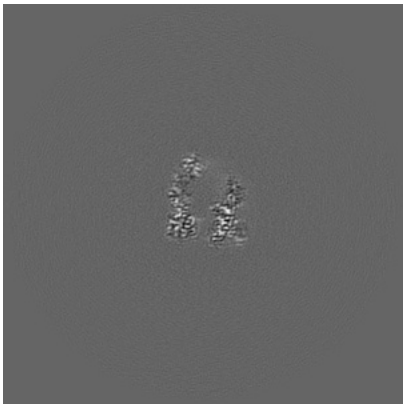
6.2.1 Primary map



X Index: 256



Y Index: 256

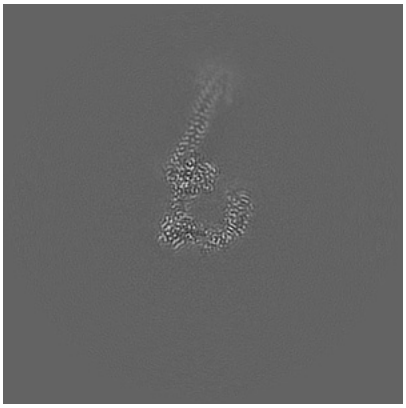


Z Index: 256

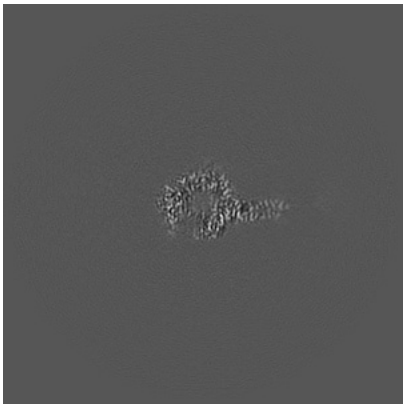
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices (i)

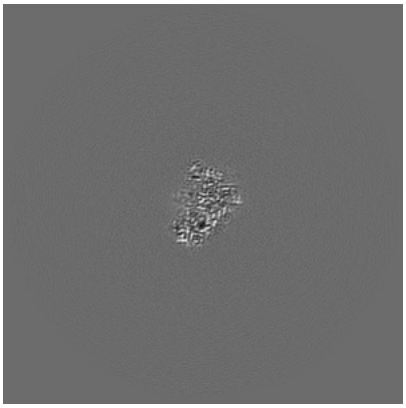
6.3.1 Primary map



X Index: 245



Y Index: 237

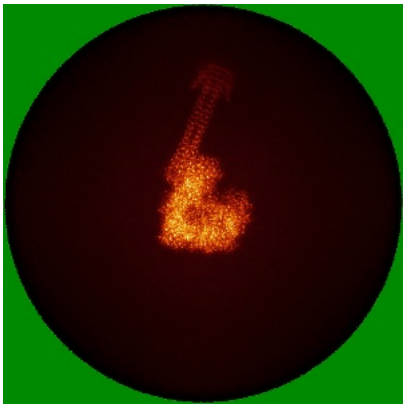


Z Index: 222

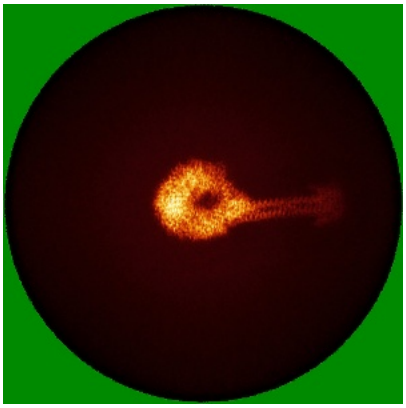
The images above show the largest variance slices of the map in three orthogonal directions.

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

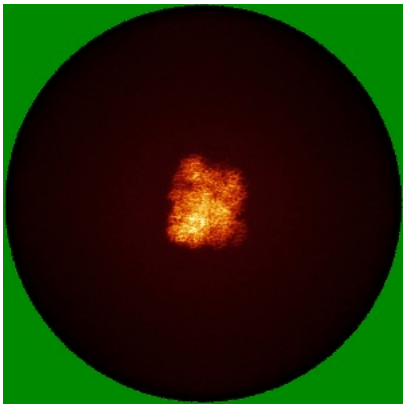
6.4.1 Primary map



X



Y

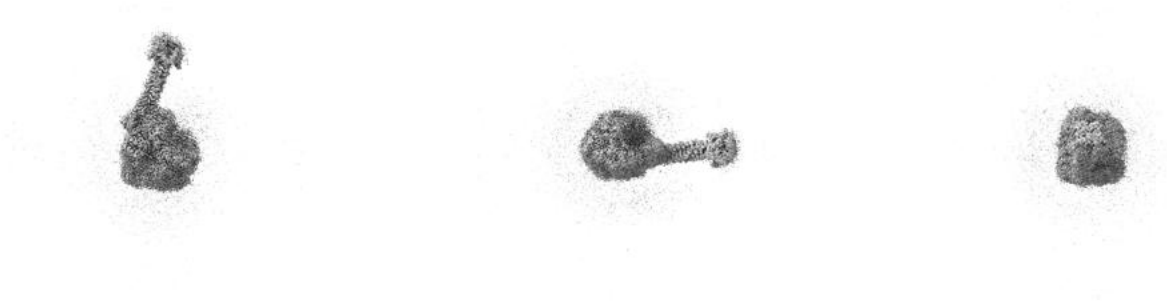


Z

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

6.5 Orthogonal surface views (i)

6.5.1 Primary map



X Y Z

The images above show the 3D surface view of the map at the recommended contour level 3.0. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

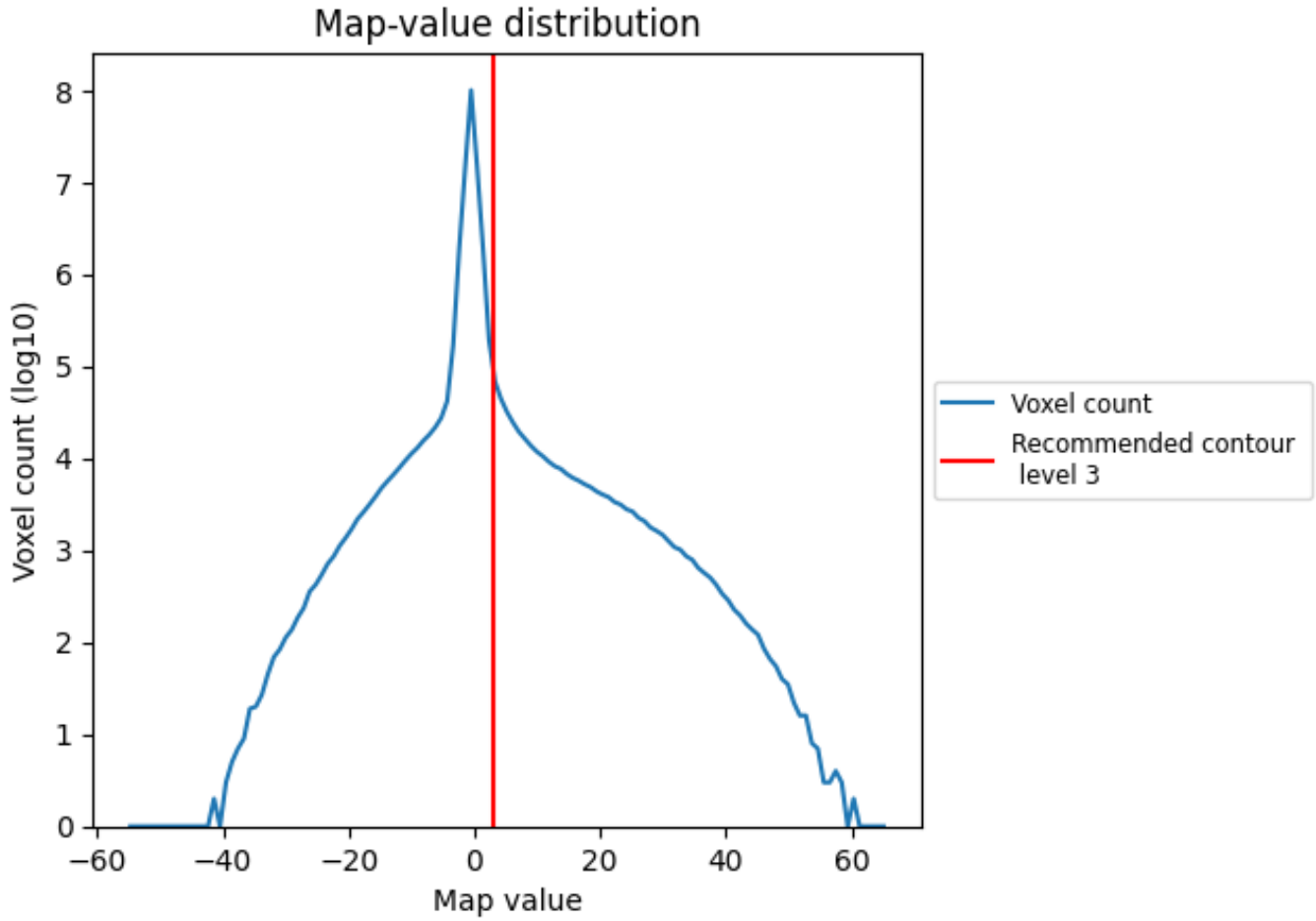
6.6 Mask visualisation (i)

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

7 Map analysis

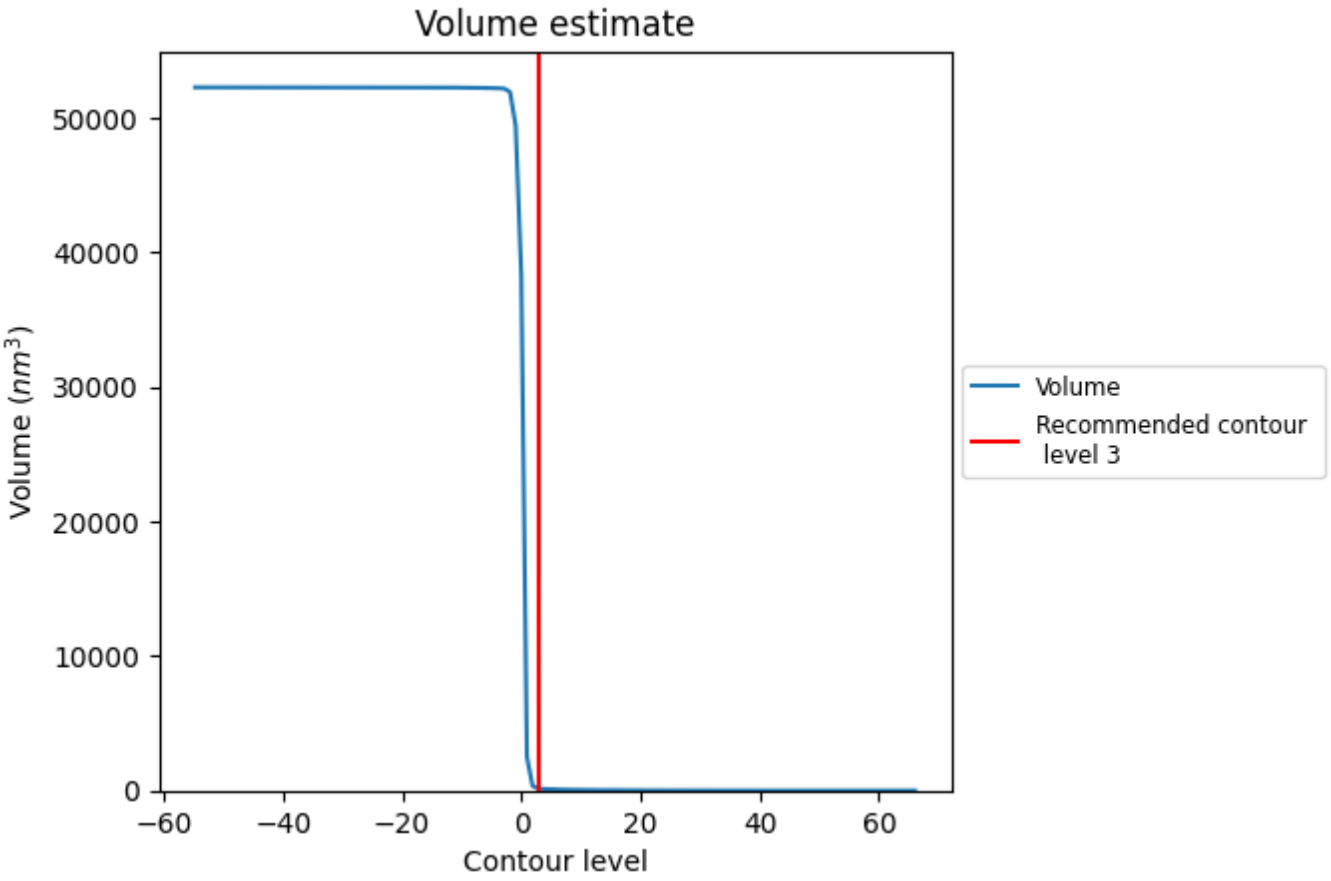
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution



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.

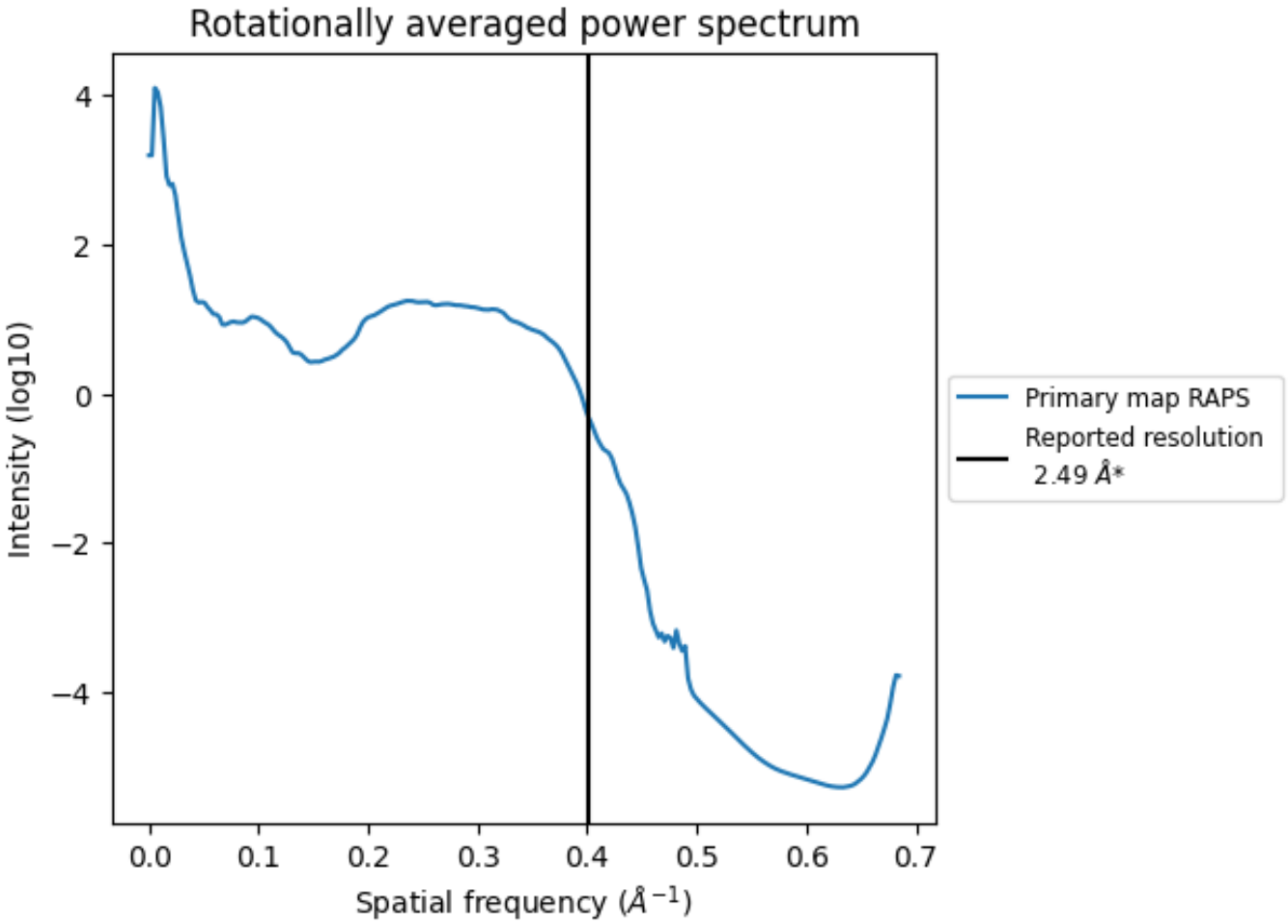
7.2 Volume estimate



The volume at the recommended contour level is 146 nm^3; this corresponds to an approximate mass of 132 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 (i)



*Reported resolution corresponds to spatial frequency of 0.402 Å^{-1}

8 Fourier-Shell correlation (i)

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

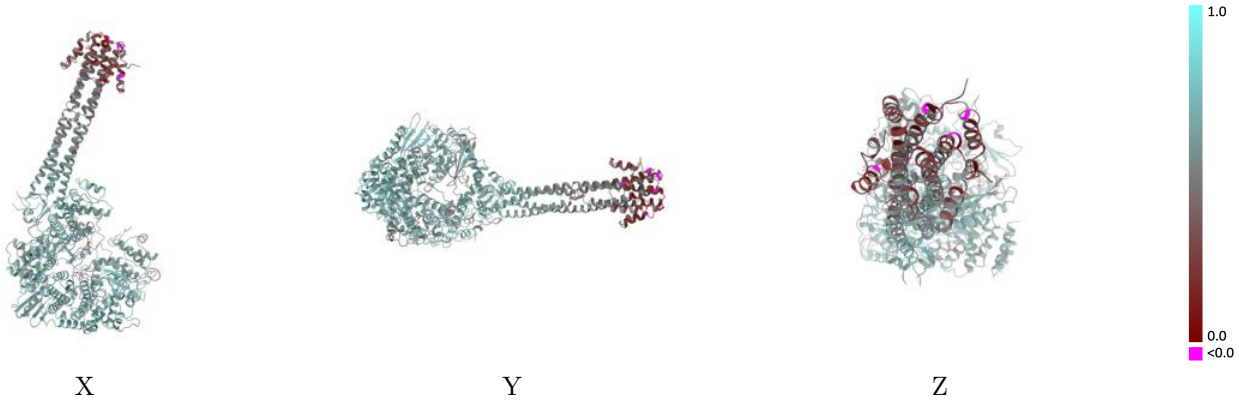
9 Map-model fit

This section contains information regarding the fit between EMDB map EMD-50781 and PDB model 9FUX. Per-residue inclusion information can be found in section 3 on page 5.

9.1 Map-model overlay

This section was not generated.

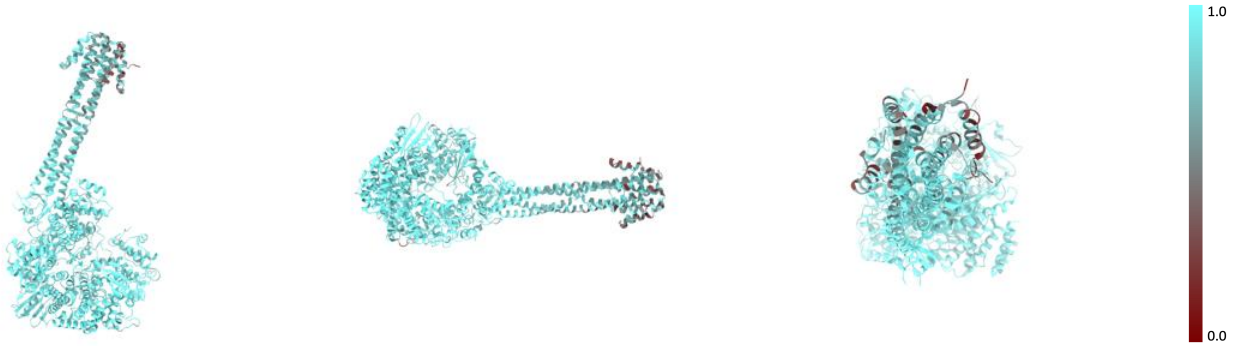
9.2 Q-score mapped to coordinate model



X Y Z

The images above show the model with each residue coloured according 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.

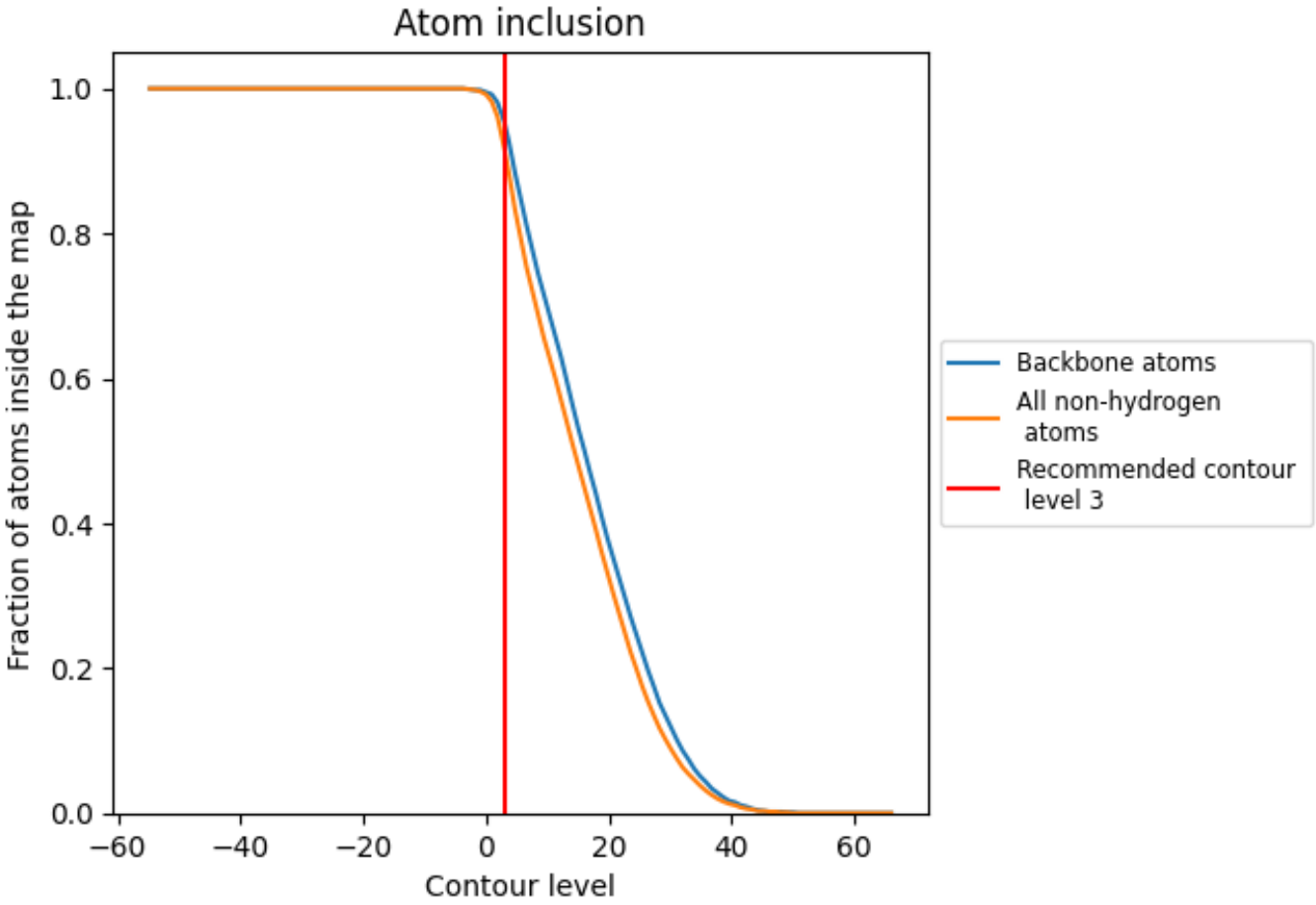
9.3 Atom inclusion mapped to coordinate model



X Y Z

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 (3).

9.4 Atom inclusion (i)



At the recommended contour level, 95% of all backbone atoms, 92% 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 (3) and Q-score for the entire model and for each chain.



Chain	Atom inclusion	Q-score
All	0.9170	0.5700
A	0.9620	0.6150
C	0.7940	0.4400
D	0.7770	0.4350
E	0.8260	0.4960
F	0.8190	0.4330