Full wwPDB EM Validation Report ⓘ

Sep 15, 2026 – 01:40 am BST

PDB ID	:	8OJA / pdb_00008oja
EMDB ID	:	EMD-16909
Title	:	HSV-1 DNA polymerase-processivity factor complex in exonuclease state
Authors	:	Gustavsson, E.; Grunewald, K.; Elias, P.; Hallberg, B.M.
Deposited on	:	2023-03-24
Resolution	:	1.87 Å(reported)
Based on initial model	:	7LUF

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:

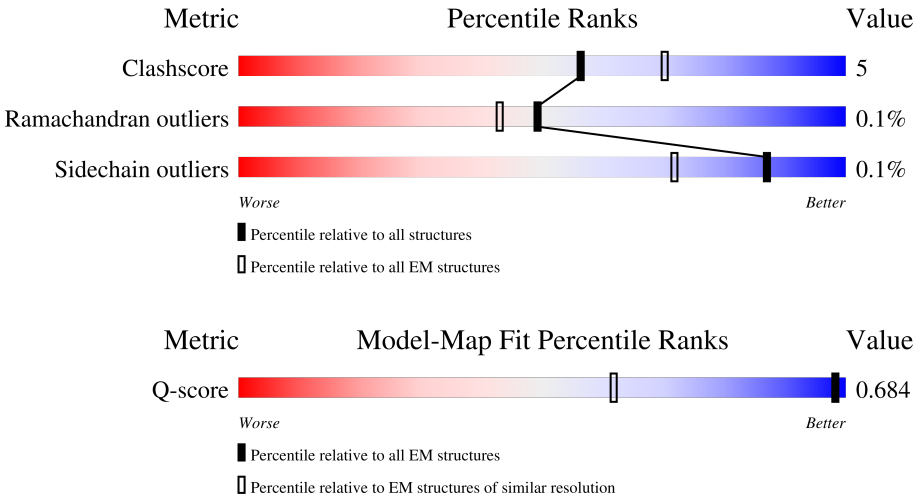
EMDB validation analysis	:	0.0.1.dev133
Mogul	:	1.8.4, CSD as541be (2020)
MolProbity	:	4-5-2 with Phenix2.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics	:	202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ	:	1.9.13
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.50

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 1.87 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	985 (1.37 - 2.37)

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	1235	81% 5% 13%
2	B	488	6% 52% 45%
3	C	47	30% 38% 23% 38%
4	D	68	13% 22% 13% 65%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 11782 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 DNA polymerase catalytic subunit.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	1072	8469	5425	1483	1521	40	4	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	330	ARG	ALA	variant	UNP P04293

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

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
2	B	266	2049	1303	360	379	7	0	0

- Molecule 3 is a DNA chain called DNA (47-MER).

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	N	O	P	S		
3	C	29	583	277	107	169	29	1	0	0

- Molecule 4 is a DNA chain called DNA (68-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	P		
4	D	24	503	236	97	146	24	0	0

- Molecule 5 is CALCIUM ION (CCD ID: CA) (formula: Ca) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		AltConf
			Total	Ca	
5	A	2	2	2	0

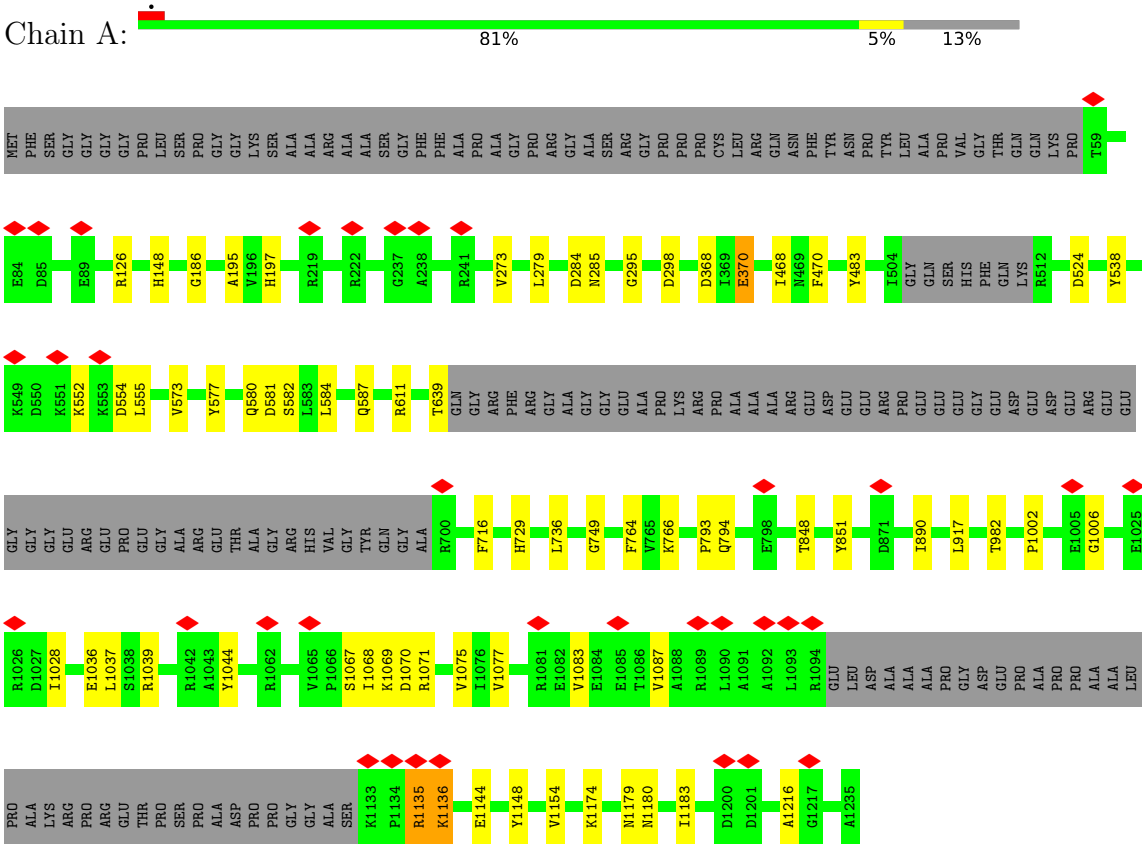
- Molecule 6 is water.

Mol	Chain	Residues	Atoms	AltConf
6	A	173	Total O 173 173	0
6	B	2	Total O 2 2	0
6	C	1	Total O 1 1	0

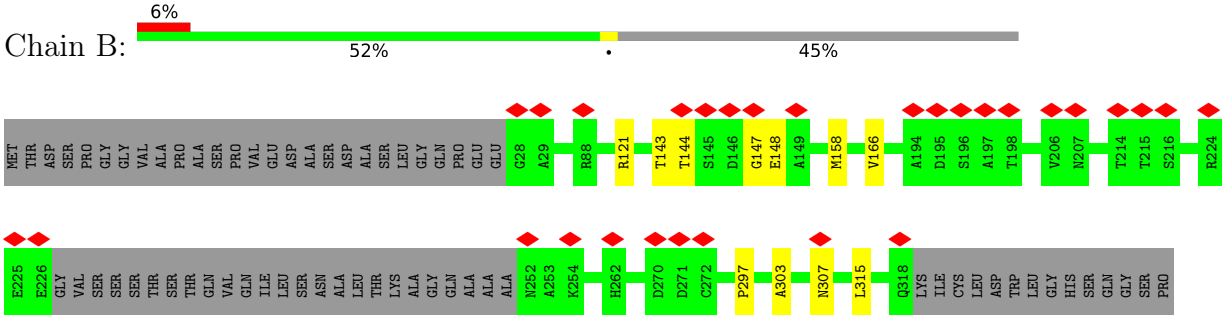
3 Residue-property plots (i)

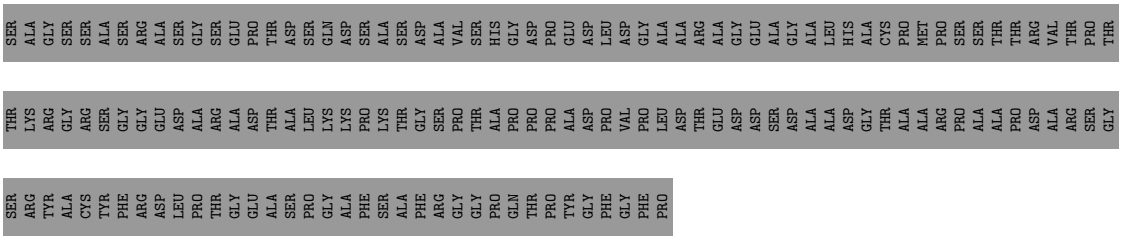
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion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



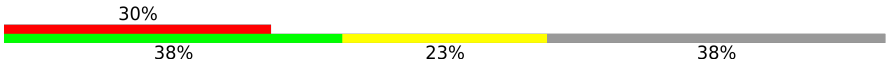
• Molecule 2: DNA polymerase processivity factor

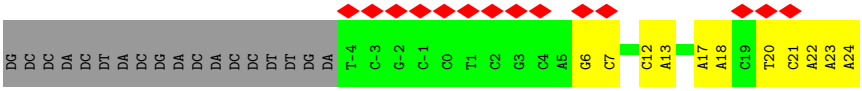




• Molecule 3: DNA (47-MER)

Chain C:

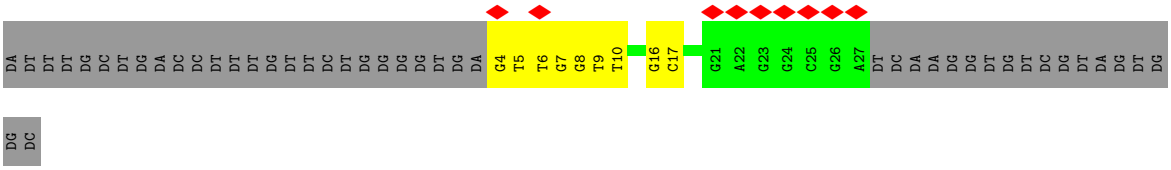




• Molecule 4: DNA (68-MER)

Chain D:





4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	539028	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 ($e^-/Å^2$)	58	Depositor
Minimum defocus (nm)	500	Depositor
Maximum defocus (nm)	1500	Depositor
Magnification	165000	Depositor
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	5.214	Depositor
Minimum map value	-1.336	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.036	Depositor
Recommended contour level	0.3	Depositor
Map size (Å)	353.5, 353.5, 353.5	wwPDB
Map dimensions	700, 700, 700	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.505, 0.505, 0.505	Depositor

5 Model quality

5.1 Standard geometry

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

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	0.62	0/8679	0.93	5/11771 (0.0%)
2	B	0.52	0/2087	0.84	0/2840
3	C	0.38	0/628	0.79	0/963
4	D	0.38	0/565	1.09	5/873 (0.6%)
All	All	0.58	0/11959	0.92	10/16447 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	3

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	D	16	DG	OP1-P-O3'	-15.26	62.23	108.00
4	D	16	DG	O3'-P-O5'	8.70	117.05	104.00
1	A	370	GLU	CB-CG-CD	8.68	127.36	112.60
1	A	126	ARG	NE-CZ-NH1	-7.32	114.18	121.50
1	A	126	ARG	NE-CZ-NH2	6.20	124.78	119.20
4	D	16	DG	OP2-P-O3'	6.16	126.47	108.00
4	D	17	DC	O5'-P-OP1	-5.90	91.30	109.00
4	D	16	DG	P-O3'-C3'	-5.51	111.93	120.20
1	A	298	ASP	CA-CB-CG	5.49	118.09	112.60
1	A	483	TYR	CB-CA-C	5.19	118.91	109.83

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1039	ARG	Sidechain
1	A	1135	ARG	Sidechain
1	A	611	ARG	Sidechain

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	8469	0	8475	78	0
2	B	2049	0	2091	9	0
3	C	583	0	324	18	0
4	D	503	0	270	7	0
5	A	2	0	0	0	0
6	A	173	0	0	13	0
6	B	2	0	0	0	0
6	C	1	0	0	0	0
All	All	11782	0	11160	111	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 5.

All (111) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:552:LYS:CE	1:A:584:LEU:HD11	1.19	1.64
1:A:552:LYS:CE	1:A:584:LEU:CD1	2.10	1.28
1:A:552:LYS:HE2	1:A:584:LEU:CD1	1.64	1.25
1:A:555:LEU:HD11	1:A:577[A]:TYR:CE2	1.76	1.19
1:A:555:LEU:HD11	1:A:577[A]:TYR:CZ	1.85	1.11
1:A:555:LEU:HD13	1:A:577[A]:TYR:CG	1.88	1.08
1:A:555:LEU:CD1	1:A:577[A]:TYR:CG	2.38	1.05
1:A:552:LYS:HE3	1:A:584:LEU:HD11	1.06	1.03
1:A:555:LEU:CD1	1:A:577[A]:TYR:CD1	2.43	1.00
1:A:555:LEU:HD13	1:A:577[A]:TYR:CD1	1.96	1.00
1:A:555:LEU:HD11	1:A:577[A]:TYR:CD2	1.95	0.99

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:552:LYS:HE3	1:A:584:LEU:CD1	1.85	0.94
3:C:21:DC:H4'	3:C:22:DA:H5'	1.51	0.92
1:A:552:LYS:HE2	1:A:584:LEU:HD11	0.93	0.90
6:A:1408:HOH:O	3:C:23:DA:H4'	1.69	0.90
1:A:555:LEU:HD11	1:A:577[A]:TYR:CE1	2.06	0.90
1:A:555:LEU:CD1	1:A:577[A]:TYR:CD2	2.58	0.87
1:A:555:LEU:CD1	1:A:577[A]:TYR:CE1	2.60	0.83
1:A:1083:VAL:O	1:A:1087:VAL:HG23	1.78	0.83
1:A:555:LEU:HD21	1:A:577[A]:TYR:CD2	2.14	0.82
1:A:555:LEU:HD11	1:A:577[A]:TYR:CG	2.12	0.79
1:A:555:LEU:CD2	1:A:577[A]:TYR:CD2	2.70	0.74
1:A:555:LEU:HD11	1:A:577[A]:TYR:CD1	2.19	0.73
1:A:639:THR:HG23	6:A:1435:HOH:O	1.89	0.73
1:A:552:LYS:HD2	1:A:580[B]:GLN:CD	2.15	0.71
1:A:370:GLU:OE1	3:C:24:AS:O5'	2.09	0.70
1:A:1135:ARG:O	1:A:1136:LYS:C	2.39	0.66
1:A:368:ASP:OD2	6:A:1403:HOH:O	2.14	0.65
1:A:555:LEU:HD12	1:A:555:LEU:N	2.12	0.64
4:D:5:DT:H2"	4:D:6:DT:C6	2.35	0.62
1:A:197:HIS:CD2	6:A:1442:HOH:O	2.53	0.61
2:B:144:THR:OG1	2:B:148:GLU:HB3	2.00	0.61
1:A:555:LEU:HD13	1:A:577[B]:TYR:CD1	2.36	0.60
1:A:1037:LEU:O	1:A:1069:LYS:HA	2.02	0.59
1:A:1067:SER:HB2	1:A:1070:ASP:OD1	2.02	0.59
1:A:555:LEU:HD13	1:A:577[B]:TYR:CG	2.38	0.58
4:D:6:DT:H2"	4:D:7:DG:C8	2.37	0.58
3:C:24:AS:OP1	3:C:24:AS:H8	2.05	0.57
4:D:9:DT:H2"	4:D:10:DT:H71	1.87	0.57
2:B:297:PRO:HG2	2:B:315:LEU:HB2	1.88	0.56
1:A:284:ASP:OD1	1:A:285:ASN:ND2	2.41	0.54
1:A:555:LEU:CD1	1:A:555:LEU:N	2.71	0.54
1:A:1144:GLU:OE2	1:A:1148:TYR:CD1	2.62	0.54
3:C:20:DT:H2'	3:C:21:DC:H2'	1.89	0.53
3:C:21:DC:C2	3:C:22:DA:C5	2.96	0.53
1:A:793:PRO:HB2	1:A:794:GLN:OE1	2.09	0.52
2:B:121:ARG:HG3	2:B:143:THR:HB	1.92	0.52
1:A:554:ASP:OD1	1:A:577[B]:TYR:CZ	2.63	0.52
3:C:22:DA:N7	3:C:23:DA:N6	2.58	0.51
1:A:195:ALA:C	6:A:1442:HOH:O	2.53	0.51
4:D:8:DG:H2'	4:D:9:DT:C6	2.45	0.51
1:A:851:TYR:HB2	1:A:917:LEU:HD11	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1077:VAL:HG13	1:A:1154:VAL:HG12	1.92	0.51
1:A:1216:ALA:HB2	2:B:166:VAL:HG22	1.93	0.51
1:A:197:HIS:CG	6:A:1442:HOH:O	2.65	0.50
2:B:303:ALA:CB	2:B:307:ASN:OD1	2.59	0.50
4:D:8:DG:H2"	4:D:9:DT:O5'	2.12	0.50
1:A:148:HIS:O	1:A:186:GLY:HA3	2.11	0.50
1:A:716:PHE:HB2	1:A:890:ILE:CG1	2.42	0.49
1:A:468:ILE:HG12	1:A:524:ASP:HB2	1.95	0.48
1:A:716:PHE:HB2	1:A:890:ILE:HG13	1.95	0.48
3:C:21:DC:H4'	3:C:22:DA:C5'	2.34	0.48
1:A:552:LYS:HD2	1:A:580[B]:GLN:NE2	2.29	0.47
1:A:639:THR:CG2	6:A:1435:HOH:O	2.54	0.47
1:A:538:TYR:N	3:C:22:DA:OP2	2.47	0.47
1:A:554:ASP:HA	1:A:577[B]:TYR:CE1	2.49	0.47
1:A:1174:LYS:HG2	1:A:1179:ASN:HA	1.95	0.47
3:C:22:DA:C5	3:C:23:DA:C6	3.03	0.47
1:A:729:HIS:HE1	6:A:1561:HOH:O	1.97	0.46
1:A:736:LEU:HB2	1:A:764:PHE:CE1	2.50	0.46
4:D:7:DG:H2"	4:D:8:DG:C8	2.51	0.46
1:A:1036:GLU:HA	1:A:1071:ARG:HA	1.98	0.45
1:A:1067:SER:O	1:A:1068:ILE:HG13	2.16	0.45
1:A:470:PHE:CD2	3:C:23:DA:H2"	2.52	0.45
1:A:1180:ASN:ND2	1:A:1183:ILE:HD12	2.32	0.44
3:C:17:DA:H2"	3:C:18:DA:C8	2.52	0.44
1:A:1044:TYR:HE2	3:C:18:DA:OP1	2.01	0.44
1:A:1002:PRO:HA	2:B:158:MET:HE1	1.99	0.43
1:A:554:ASP:OD1	1:A:577[B]:TYR:OH	2.36	0.43
1:A:1028:ILE:HG23	1:A:1075:VAL:HG11	2.00	0.43
1:A:749:GLY:O	1:A:766:LYS:HE3	2.19	0.42
1:A:794:GLN:N	1:A:794:GLN:CD	2.77	0.42
4:D:4:DG:H1'	4:D:5:DT:H5'	2.00	0.42
1:A:279:LEU:HD11	1:A:295:GLY:O	2.20	0.42
2:B:297:PRO:HG2	2:B:315:LEU:CB	2.49	0.42
2:B:307:ASN:OD1	2:B:307:ASN:O	2.37	0.42
3:C:6:DG:C4	3:C:7:DC:C5	3.08	0.42
1:A:982:THR:HG22	1:A:1006:GLY:HA2	2.01	0.41
1:A:273:VAL:HG21	1:A:279:LEU:HB2	2.02	0.41
3:C:21:DC:O2	3:C:21:DC:O4'	2.35	0.41
3:C:12:DC:H2"	3:C:13:DA:C8	2.56	0.41
1:A:848:THR:HG21	1:A:890:ILE:HD13	2.02	0.41
1:A:1037:LEU:HD23	1:A:1037:LEU:HA	1.94	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:20:DT:H2'	3:C:21:DC:C2	2.55	0.41
1:A:552:LYS:HE2	1:A:584:LEU:HD12	1.83	0.41
2:B:143:THR:OG1	2:B:147:GLY:HA2	2.21	0.41
1:A:848:THR:CG2	1:A:890:ILE:HD13	2.52	0.40
3:C:6:DG:C5	3:C:7:DC:C5	3.09	0.40
1:A:1067:SER:HB2	1:A:1070:ASP:CG	2.46	0.40
1:A:555:LEU:HD21	1:A:573:VAL:HG12	2.03	0.40
1:A:548:LEU:HD13	1:A:587:GLN:OE1	2.21	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	1068/1235 (86%)	1035 (97%)	32 (3%)	1 (0%)	48 37
2	B	262/488 (54%)	256 (98%)	6 (2%)	0	100 100
All	All	1330/1723 (77%)	1291 (97%)	38 (3%)	1 (0%)	49 37

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1136	LYS

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	889/996 (89%)	887 (100%)	2 (0%)	87 84
2	B	227/388 (58%)	227 (100%)	0	100 100
All	All	1116/1384 (81%)	1114 (100%)	2 (0%)	87 84

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	581[A]	ASP
1	A	581[B]	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	98	HIS
1	A	135	HIS
1	A	285	ASN
1	A	307	ASN
1	A	408	HIS
1	A	729	HIS
1	A	794	GLN
1	A	806	GLN
1	A	832	HIS
1	A	1180	ASN
2	B	76	GLN
2	B	177	GLN
2	B	282	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)

1 non-standard protein/DNA/RNA residue is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
3	AS	C	24	3,5	20,23,24	0.27	0	29,33,36	0.38	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	AS	C	24	3,5	-	3/7/21/22	0/3/3/3

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	24	AS	C3'-C4'-C5'-O5'
3	C	24	AS	O4'-C4'-C5'-O5'
3	C	24	AS	C2'-C1'-N9-C4

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	24	AS	2	0

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.

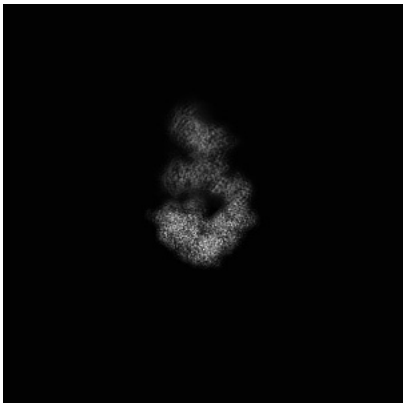
6 Map visualisation

This section contains visualisations of the EMDB entry EMD-16909. These allow visual inspection of the internal detail of the map and identification of artifacts.

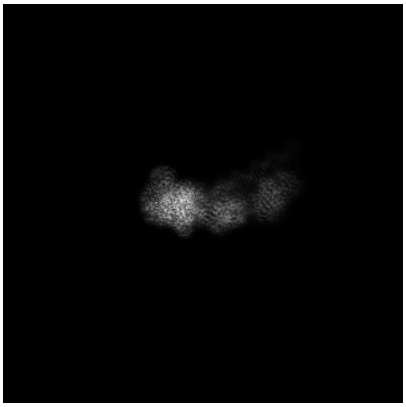
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections

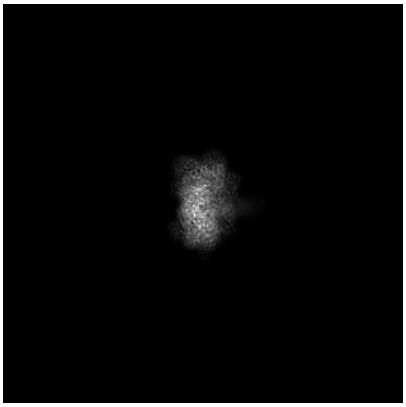
6.1.1 Primary map



X



Y



Z

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

6.2 Central slices

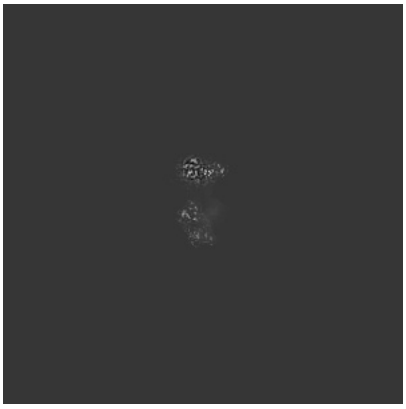
6.2.1 Primary map



X Index: 350



Y Index: 350

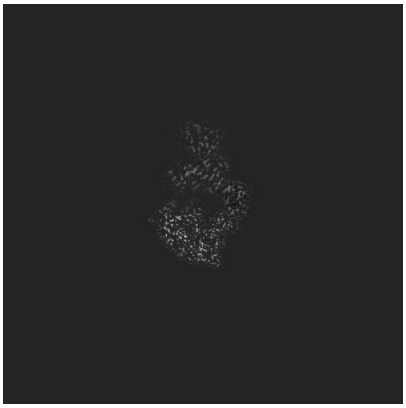


Z Index: 350

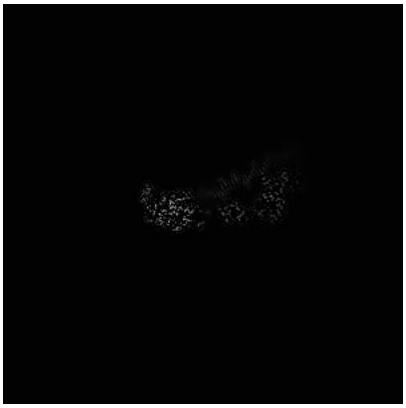
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices (i)

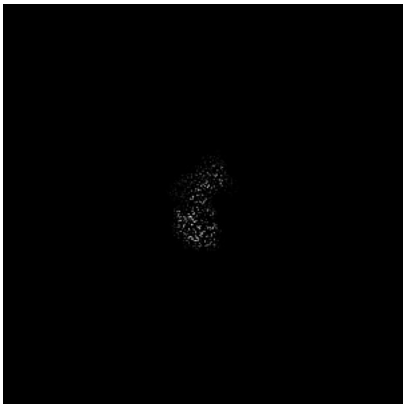
6.3.1 Primary map



X Index: 341



Y Index: 337

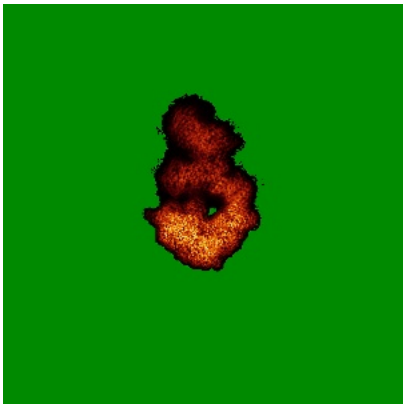


Z Index: 305

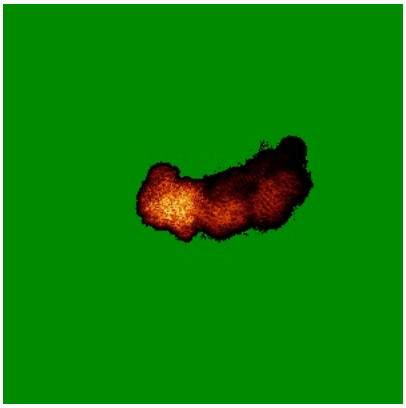
The images above show the largest variance slices of the map in three orthogonal directions.

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

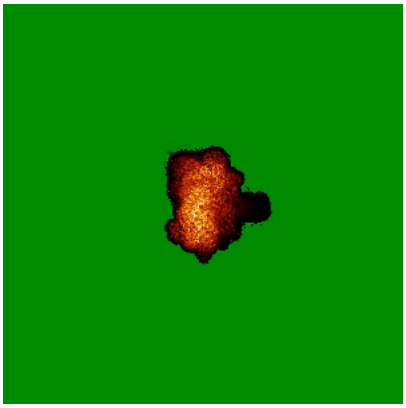
6.4.1 Primary map



X



Y

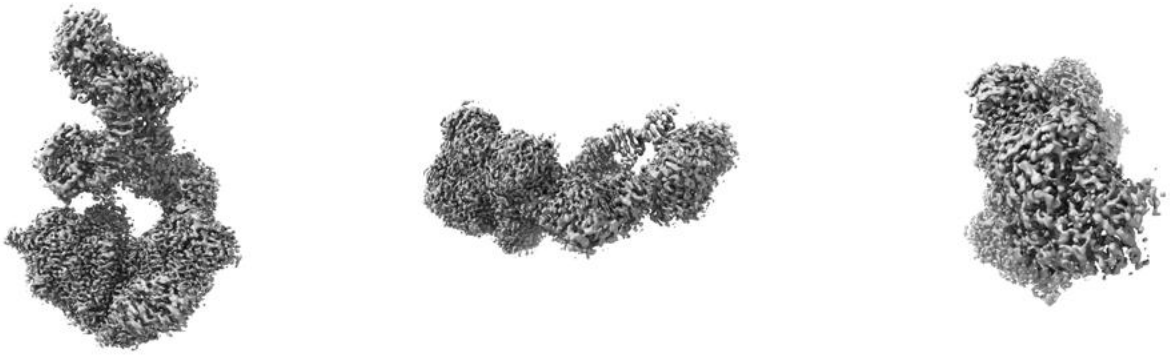


Z

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 0.3. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

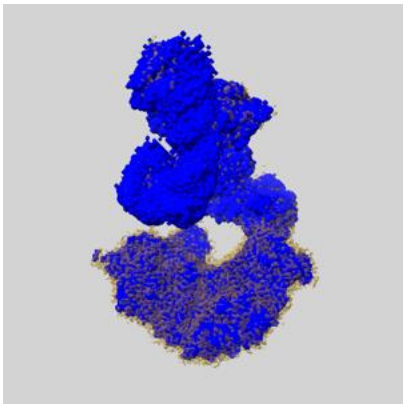
6.6 Mask visualisation (i)

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

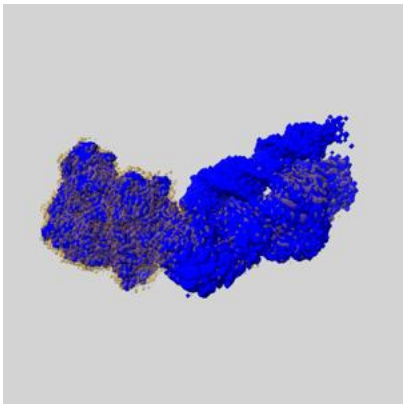
A mask typically either:

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

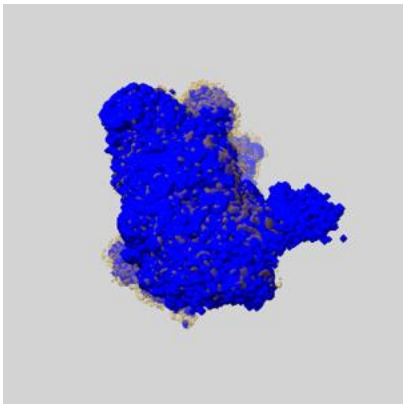
6.6.1 emd_16909_msk_1.map (i)



X



Y

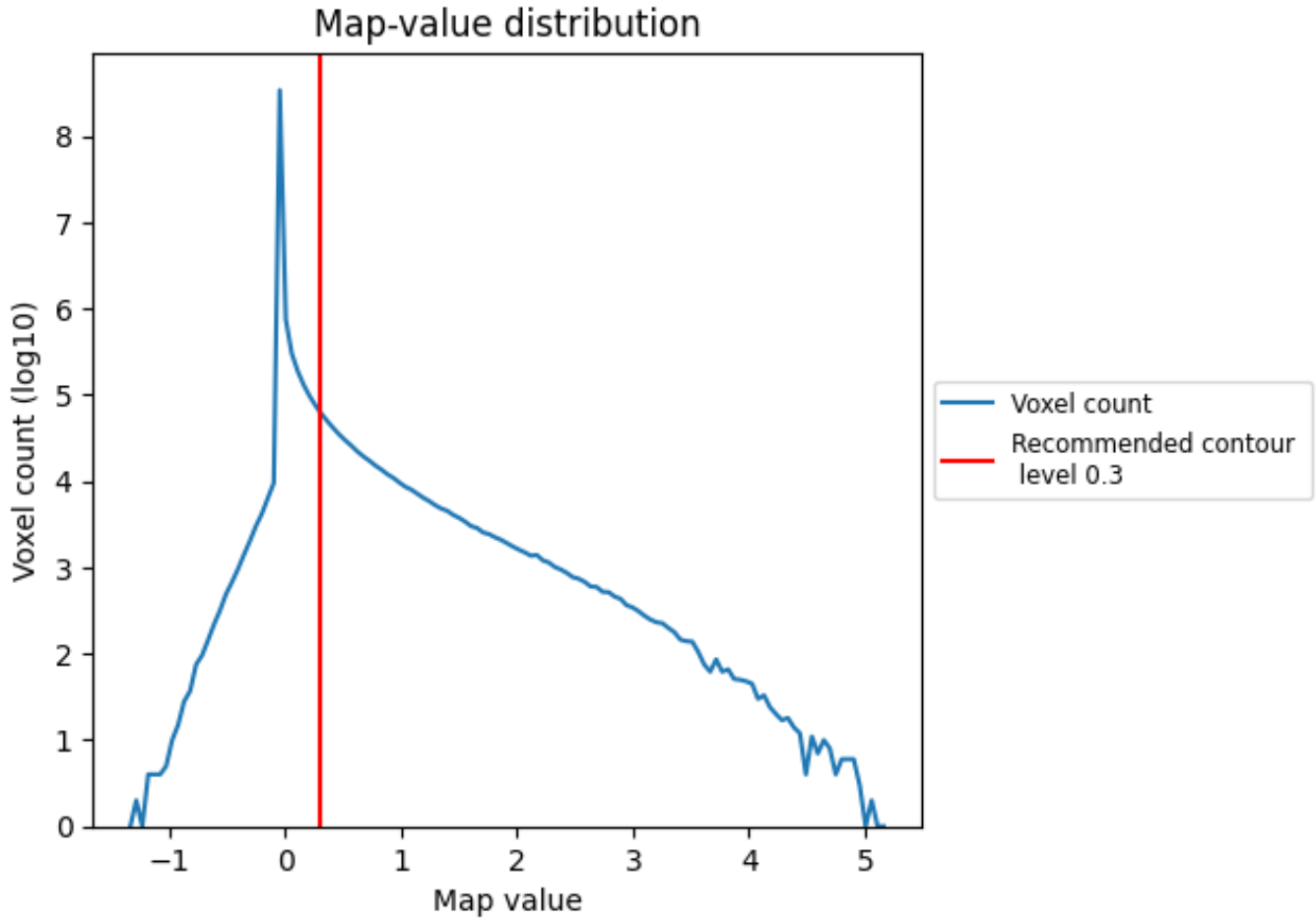


Z

7 Map analysis

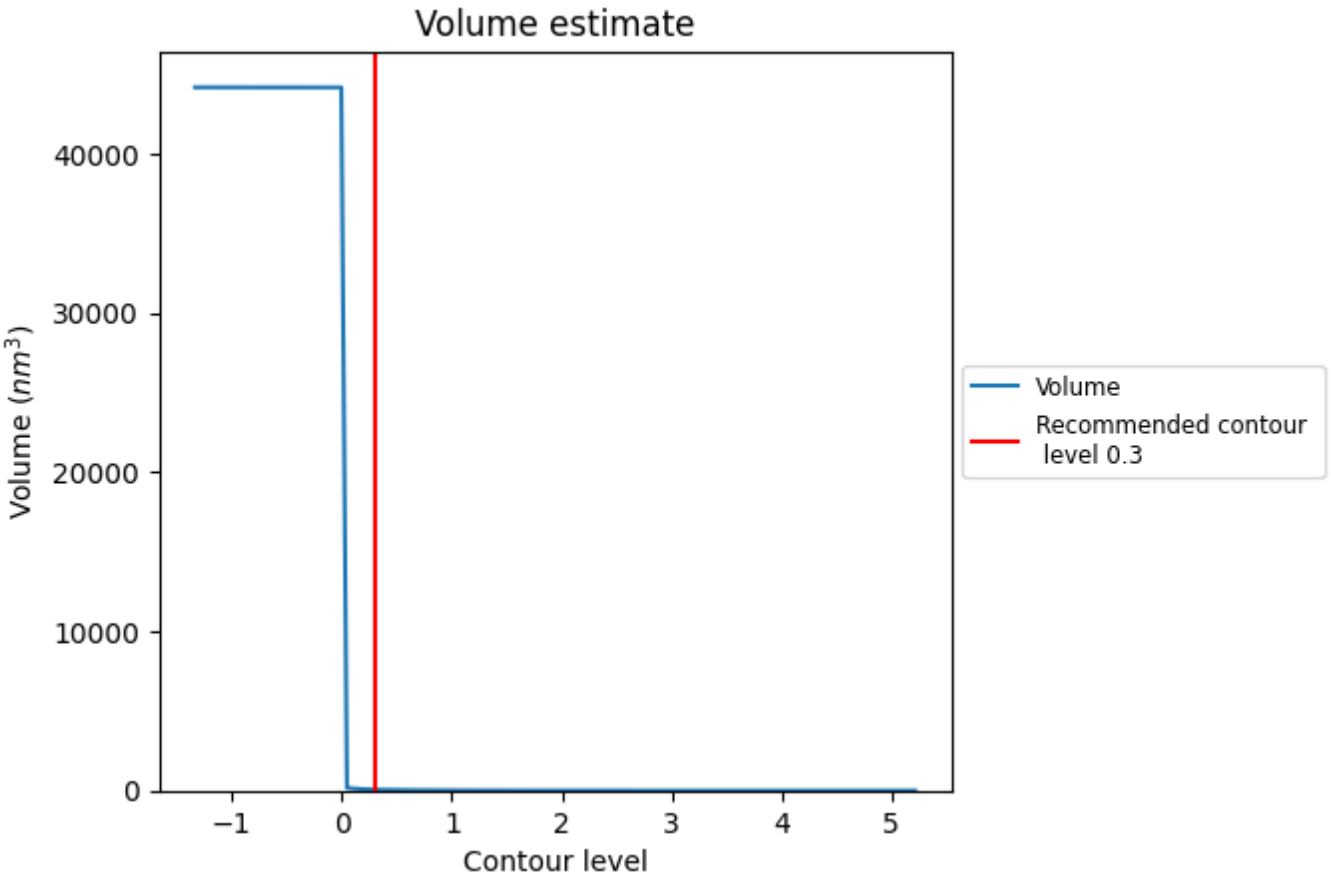
This section contains the results of statistical analysis of the map.

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.

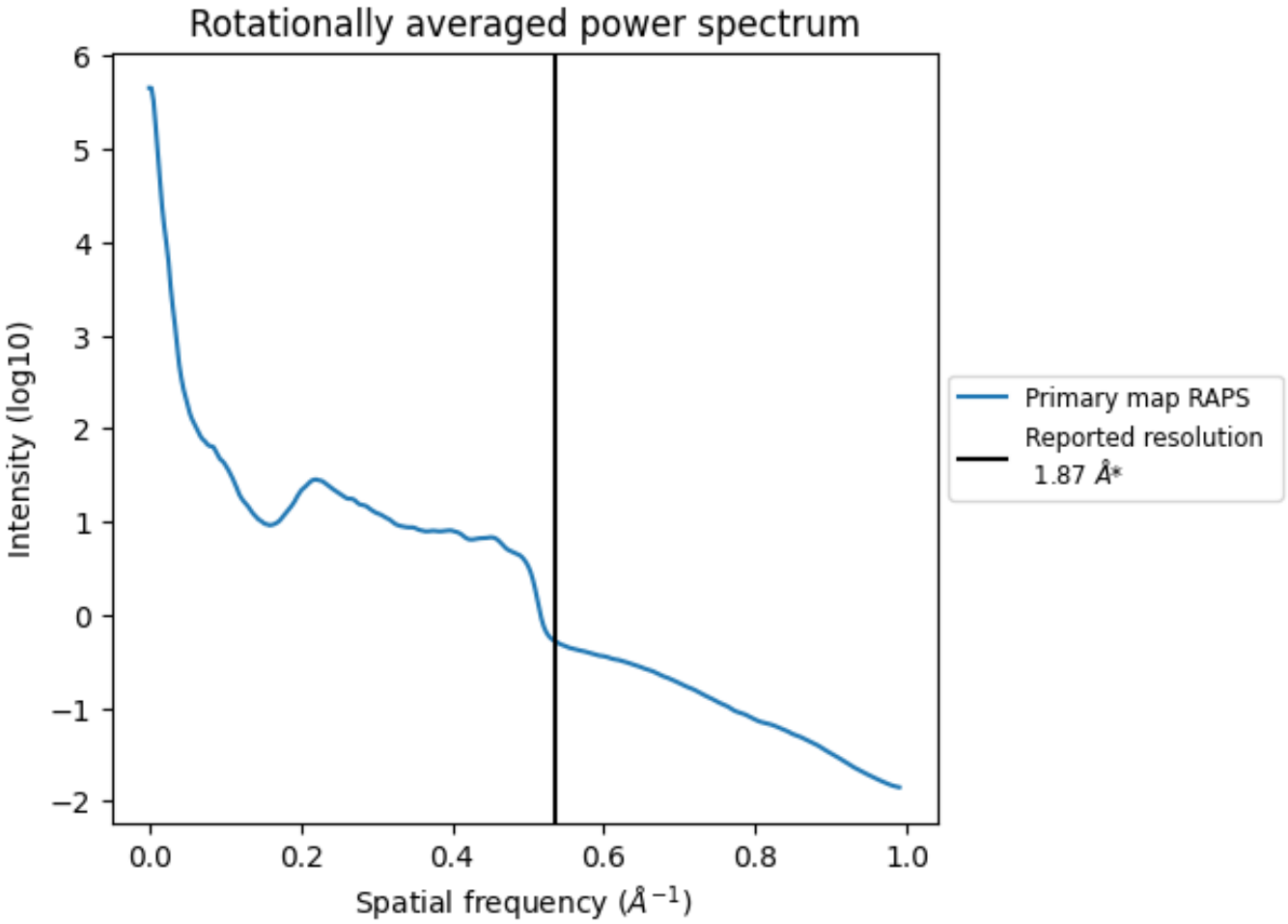
7.2 Volume estimate



The volume at the recommended contour level is 62 nm^3; this corresponds to an approximate mass of 56 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 0.535 Å^{-1}

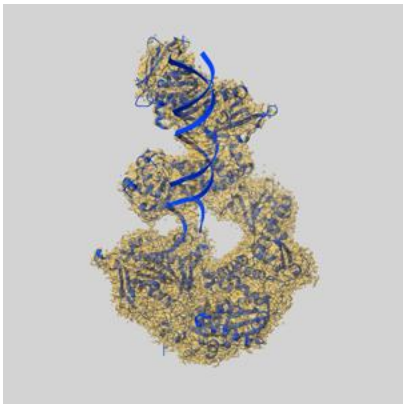
8 Fourier-Shell correlation (i)

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

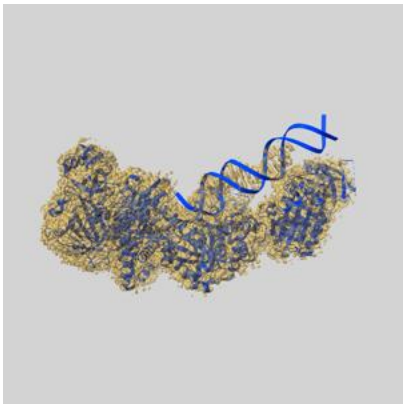
9 Map-model fit

This section contains information regarding the fit between EMDB map EMD-16909 and PDB model 8OJA. Per-residue inclusion information can be found in section 3 on page 5.

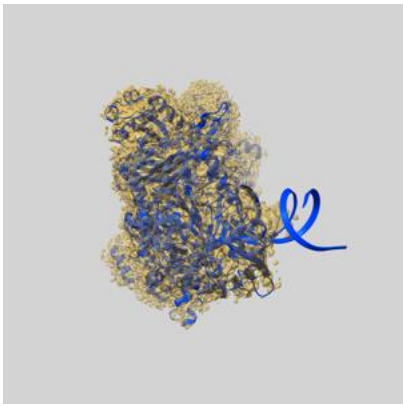
9.1 Map-model overlay



X



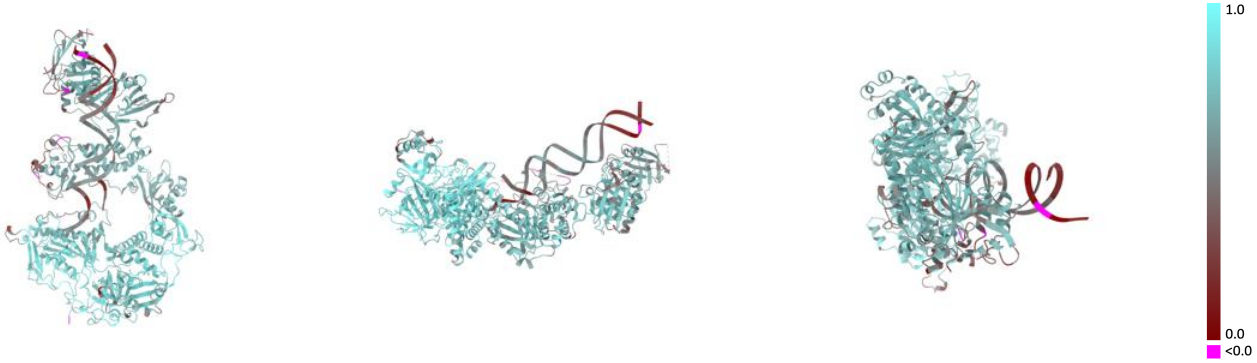
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.3 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.

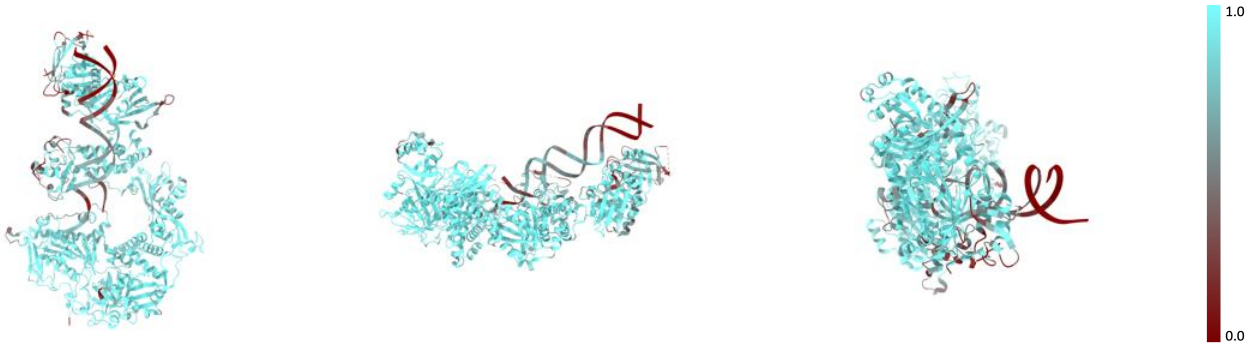
9.2 Q-score mapped to coordinate model (i)



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.

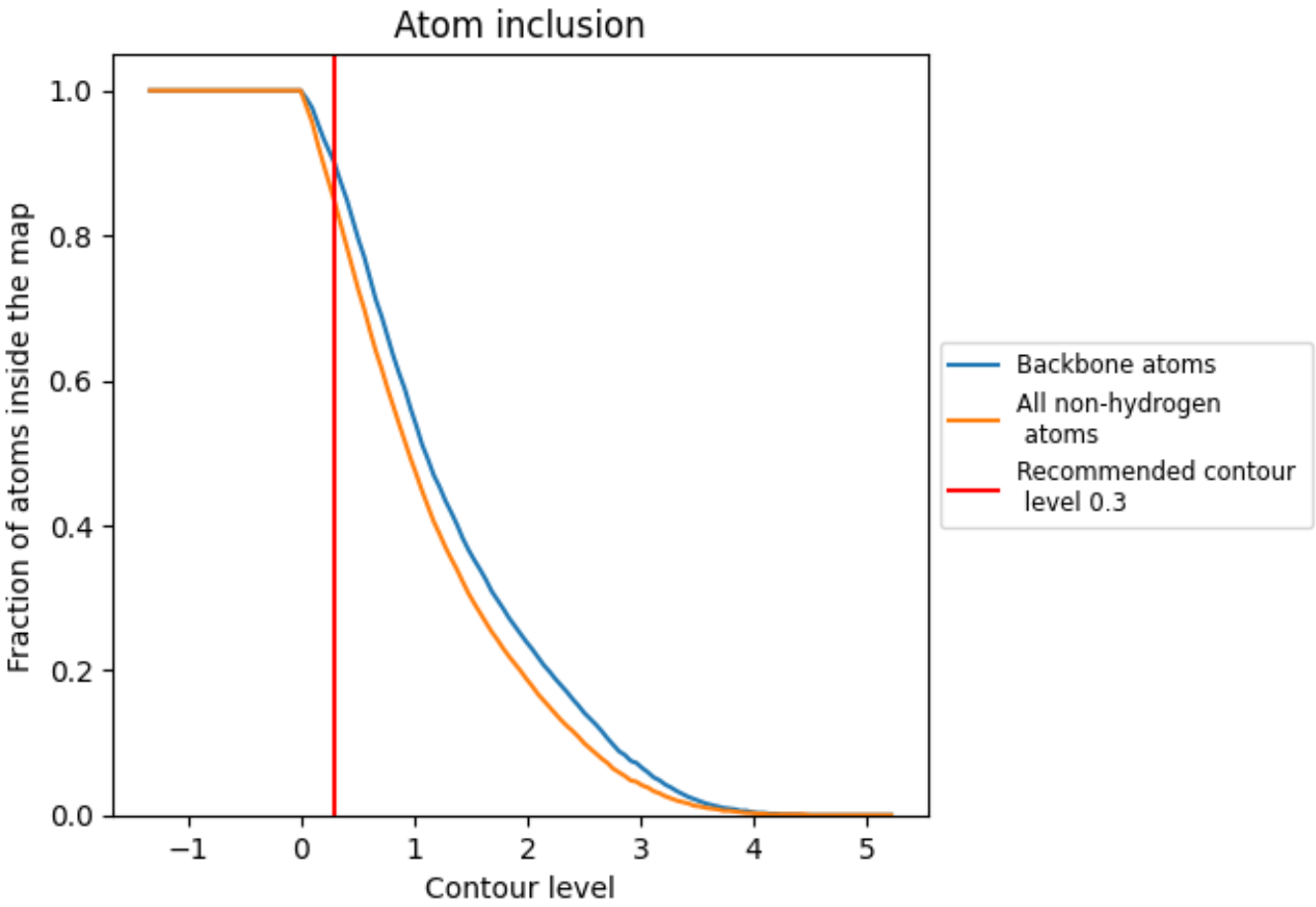
9.3 Atom inclusion mapped to coordinate model (i)



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 (0.3).

9.4 Atom inclusion (i)



At the recommended contour level, 90% of all backbone atoms, 84% 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 (0.3) and Q-score for the entire model and for each chain.



Chain	Atom inclusion	Q-score
All	0.8440	0.6840
A	0.9150	0.7360
B	0.8000	0.6280
C	0.3770	0.3500
D	0.4310	0.4240