Full wwPDB EM Validation Report (i)

Feb 27, 2024 – 06:57 AM EST

PDB ID	:	6VAF
EMDB ID	:	EMD-21139
Title	:	Structure of mono-ubiquitinated FANCD2 bound to non-ubiquitinated FANCI and to DNA
Authors	:	Pavletich, N.P.
Deposited on	:	2019-12-17
Resolution	:	3.90 Å(reported)
Based on initial model	:	3S4W

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:

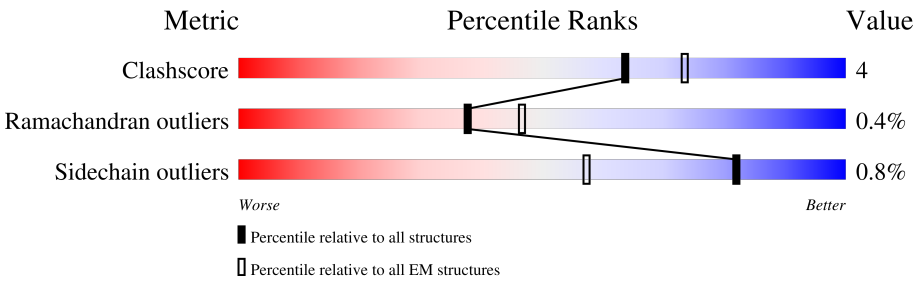
EMDB validation analysis	:	0.0.1.dev70
MolProbity	:	4.02b-467
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ	:	1.9.13
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 3.90 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	EM structures (#Entries)
Clashscore	158937	4297
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for >=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	1328	14% 79% 10% 10%
2	B	1451	12% 70% 9% 20%
3	D	76	12% 78% 22%
4	S	29	59% 34% 66%
5	T	29	45% 45% 55%

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 41243 atoms, of which 20677 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 Fanconi anemia, complementation group I.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
1	A	1198	19364	6077	9876	1592	1763	56	0	0

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	136	VAL	ALA	conflict	UNP B7ZMF2
A	476	ASN	SER	conflict	UNP B7ZMF2
A	638	GLU	LYS	conflict	UNP B7ZMF2
A	657	GLN	LYS	conflict	UNP B7ZMF2
A	877	LEU	ILE	conflict	UNP B7ZMF2
A	1235	VAL	ALA	conflict	UNP B7ZMF2
A	1274	SER	ASN	conflict	UNP B7ZMF2

- Molecule 2 is a protein called Fanconi anemia group D2 protein.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
2	B	1155	18808	5992	9520	1537	1706	53	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	654	GLN	HIS	conflict	UNP Q9BXW9
B	693	ASN	ASP	conflict	UNP Q9BXW9

- Molecule 3 is a protein called Ubiquitin.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
3	D	76	1231	378	630	105	117	1	0	0

- Molecule 4 is a DNA chain called DNA (29-MER).

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	P		
4	S	29	923	282	326	111	175	29	0	0

- Molecule 5 is a DNA chain called DNA (29-MER).

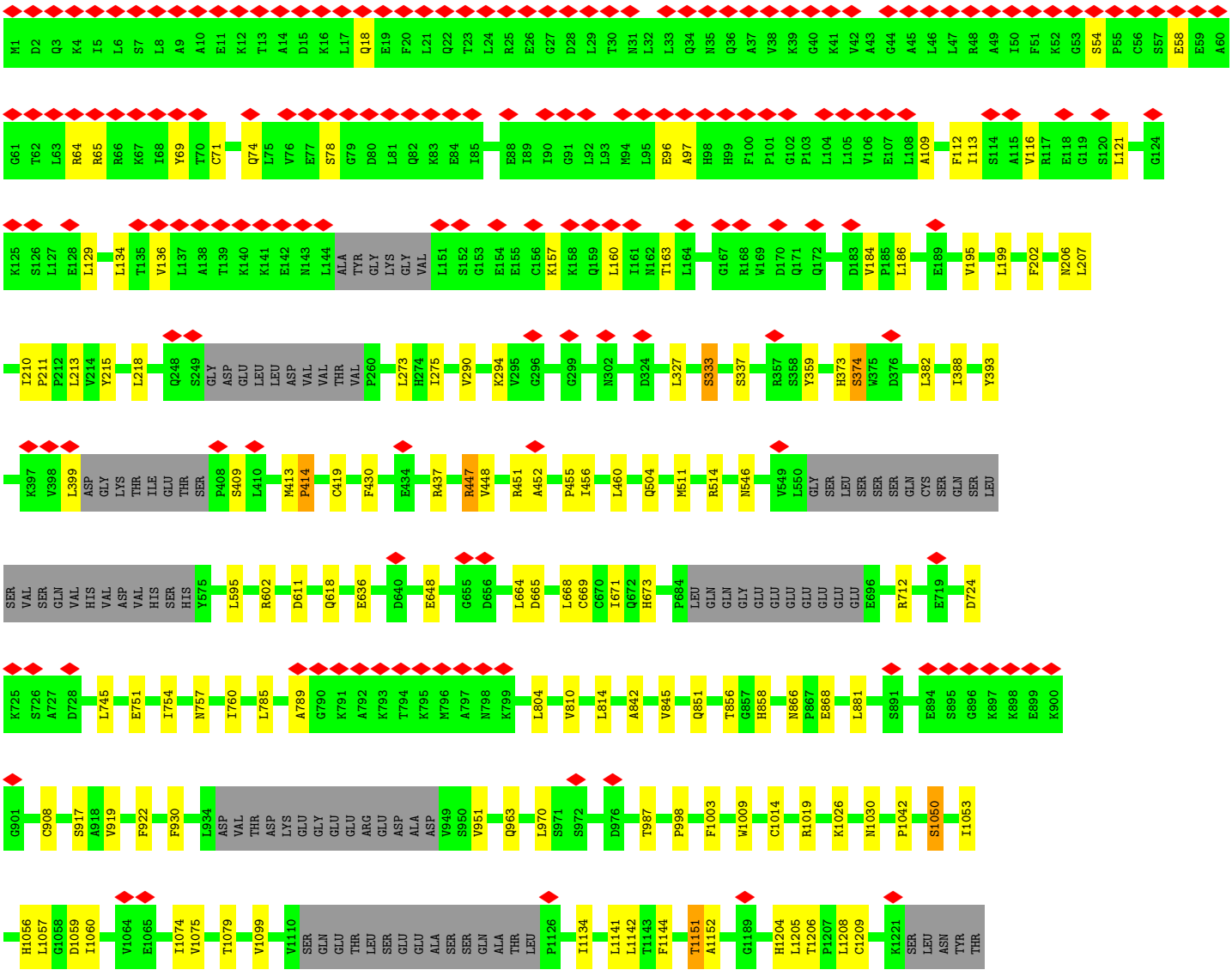
Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	P		
5	T	29	917	280	325	110	173	29	0	0

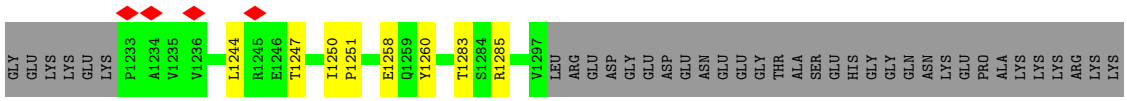
3 Residue-property plots (i)

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: Fanconi anemia, complementation group I

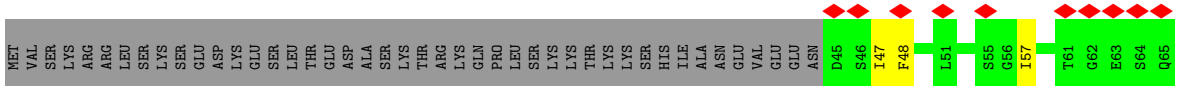
Chain A:

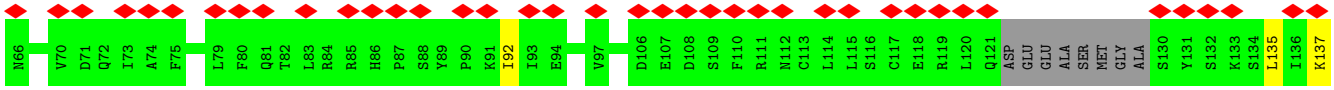


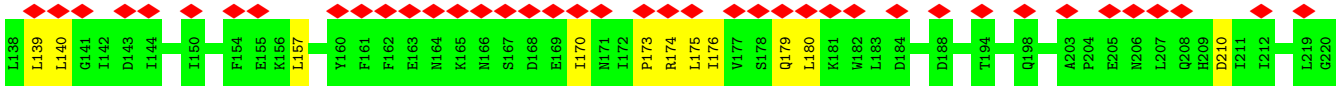


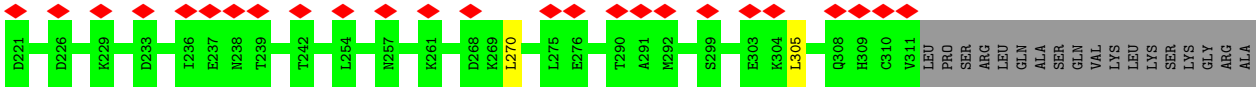
- Molecule 2: Fanconi anemia group D2 protein

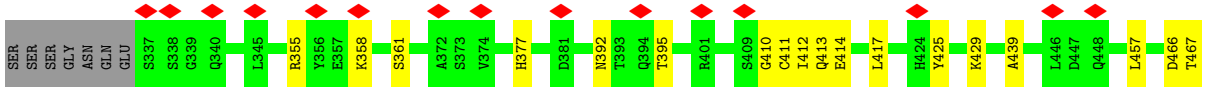
Chain B:

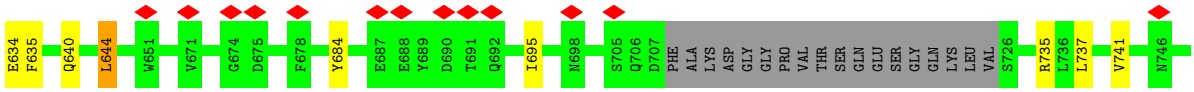


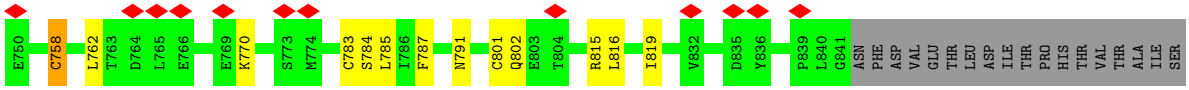


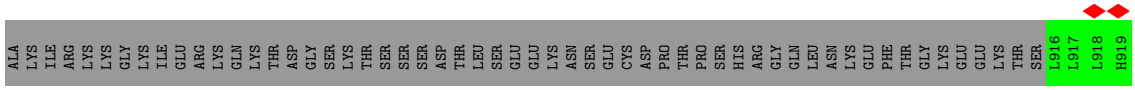


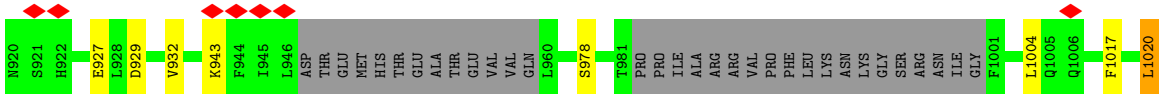


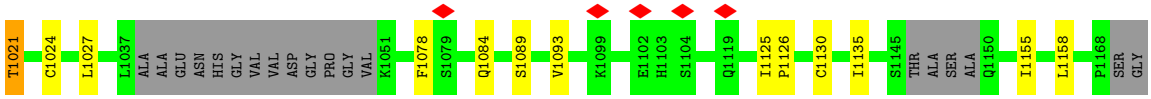


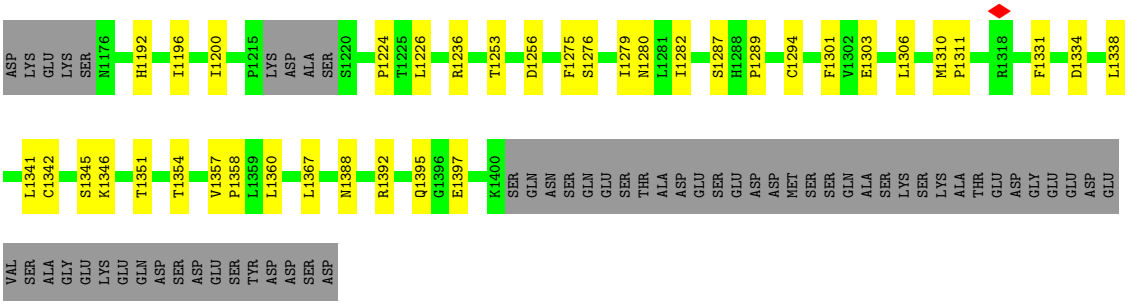






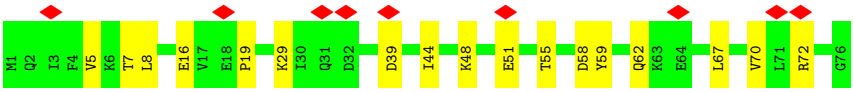






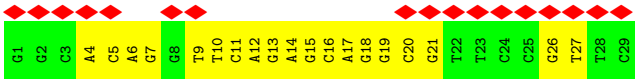
- Molecule 3: Ubiquitin

Chain D:



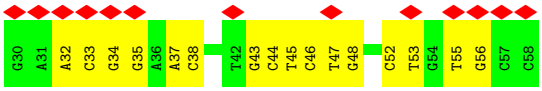
- Molecule 4: DNA (29-MER)

Chain S:



- Molecule 5: DNA (29-MER)

Chain T:



4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	57993	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/Å^2$)	51	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.098	Depositor
Minimum map value	-0.052	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.003	Depositor
Recommended contour level	0.012	Depositor
Map size (Å)	278.64832, 278.64832, 278.64832	wwPDB
Map dimensions	256, 256, 256	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.08847, 1.08847, 1.08847	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	A	0.78	2/9633 (0.0%)	0.98	4/12990 (0.0%)
2	B	0.70	0/9462	0.89	1/12795 (0.0%)
3	D	0.60	0/607	0.72	0/816
4	S	0.52	0/669	0.90	0/1031
5	T	0.50	0/663	0.91	0/1020
All	All	0.72	2/21034 (0.0%)	0.93	5/28652 (0.0%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	6
2	B	0	8
All	All	0	14

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	337	SER	CA-CB	-6.30	1.43	1.52
1	A	333	SER	CA-CB	-5.33	1.45	1.52

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	447	ARG	NE-CZ-NH2	-7.13	116.73	120.30
1	A	922	PHE	CB-CA-C	-5.45	99.50	110.40
1	A	1260	TYR	CB-CG-CD2	-5.36	117.78	121.00
2	B	1236	ARG	NE-CZ-NH1	5.13	122.86	120.30
1	A	1151	THR	CA-CB-OG1	-5.12	98.26	109.00

There are no chirality outliers.

All (14) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1079	THR	Peptide
1	A	359	TYR	Peptide
1	A	373	HIS	Peptide
1	A	414	PRO	Peptide
1	A	452	ALA	Peptide
1	A	951	VAL	Peptide
2	B	1020	LEU	Peptide
2	B	1224	PRO	Peptide
2	B	377	HIS	Peptide
2	B	411	CYS	Peptide
2	B	483	SER	Peptide
2	B	517	ILE	Peptide
2	B	644	LEU	Peptide
2	B	758	CYS	Peptide

5.2 Too-close contacts (i)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	9488	9876	9845	76	0
2	B	9288	9520	9468	68	0
3	D	601	630	629	12	0
4	S	597	326	326	14	0
5	T	592	325	325	12	0
All	All	20566	20677	20593	177	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 4.

All (177) 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:B:1017:PHE:O	2:B:1021:THR:OG1	2.07	0.71
1:A:665:ASP:HB2	1:A:804:LEU:HD11	1.76	0.67
4:S:19:DG:H2"	4:S:20:DC:C6	2.30	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:355:ARG:O	2:B:358:LYS:NZ	2.29	0.66
1:A:109:ALA:HB1	1:A:160:LEU:HD21	1.80	0.64
2:B:618:HIS:O	2:B:622:GLU:N	2.32	0.63
1:A:448:VAL:HA	1:A:456:ILE:HD12	1.81	0.63
4:S:19:DG:H2"	4:S:20:DC:C5	2.34	0.62
2:B:47:ILE:HD11	2:B:92:ILE:HD11	1.81	0.62
1:A:1050:SER:HA	1:A:1053:ILE:HD12	1.79	0.62
3:D:39:ASP:O	3:D:72:ARG:NH2	2.33	0.62
1:A:327:LEU:HD13	1:A:388:ILE:CD1	2.31	0.61
1:A:1204:HIS:O	1:A:1208:LEU:HD12	2.01	0.61
2:B:270:LEU:HD21	2:B:305:LEU:HD23	1.83	0.61
3:D:51:GLU:OE1	3:D:59:TYR:OH	2.20	0.59
3:D:16:GLU:O	3:D:29:LYS:NZ	2.37	0.57
1:A:970:LEU:O	1:A:1019:ARG:NH2	2.38	0.57
1:A:1014:CYS:HA	1:A:1074:ILE:HD11	1.87	0.57
1:A:669:CYS:SG	1:A:673:HIS:NE2	2.79	0.56
3:D:8:LEU:HD21	3:D:70:VAL:HG12	1.88	0.55
1:A:511:MET:SD	1:A:514:ARG:NH2	2.79	0.55
4:S:18:DG:H2"	4:S:19:DG:C8	2.42	0.55
2:B:137:LYS:NZ	2:B:179:GLN:O	2.38	0.55
4:S:13:DG:H2"	4:S:14:DA:C8	2.42	0.55
1:A:668:LEU:HA	1:A:671:ILE:HD12	1.89	0.54
2:B:439:ALA:HB2	2:B:457:LEU:HD23	1.90	0.54
2:B:518:LEU:HA	2:B:521:LEU:HD21	1.90	0.54
1:A:636:GLU:O	1:A:712:ARG:NH2	2.41	0.53
2:B:640:GLN:OE1	2:B:735:ARG:NH2	2.41	0.53
2:B:530:ARG:NH1	2:B:634:GLU:OE2	2.40	0.53
1:A:393:TYR:O	1:A:419:CYS:HB2	2.08	0.53
1:A:1141:LEU:O	1:A:1144:PHE:HB3	2.07	0.53
1:A:1244:LEU:O	1:A:1247:THR:HG22	2.09	0.53
2:B:783:CYS:SG	2:B:784:SER:N	2.82	0.53
2:B:816:LEU:HA	2:B:819:ILE:HD12	1.91	0.53
4:S:9:DT:C6	4:S:10:DT:H72	2.44	0.53
2:B:1078:PHE:O	2:B:1084:GLN:NE2	2.42	0.52
5:T:32:DA:H2"	5:T:33:DC:C6	2.44	0.52
1:A:448:VAL:HG13	1:A:460:LEU:HD21	1.91	0.52
2:B:1392:ARG:NH1	2:B:1397:GLU:O	2.43	0.52
2:B:1342:CYS:O	2:B:1345:SER:OG	2.23	0.52
1:A:595:LEU:O	1:A:602:ARG:NE	2.42	0.52
2:B:1020:LEU:HD12	2:B:1024:CYS:HB2	1.92	0.52
2:B:737:LEU:O	2:B:741:VAL:HG23	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:T:55:DT:H2"	5:T:56:DG:C8	2.45	0.51
1:A:664:LEU:HB3	1:A:745:LEU:HD11	1.92	0.51
2:B:586:ALA:HB1	2:B:644:LEU:HD13	1.92	0.51
1:A:113:ILE:O	1:A:116:VAL:HG12	2.11	0.51
3:D:5:VAL:HG22	3:D:67:LEU:HB2	1.93	0.51
4:S:10:DT:H2"	4:S:11:DC:C6	2.46	0.51
1:A:856:THR:OG1	1:A:858:HIS:O	2.28	0.51
1:A:113:ILE:HD11	1:A:160:LEU:HG	1.92	0.50
4:S:17:DA:H2"	4:S:18:DG:C8	2.47	0.50
1:A:184:VAL:HG23	1:A:186:LEU:HD13	1.93	0.50
1:A:1026:LYS:O	1:A:1030:ASN:ND2	2.44	0.50
1:A:785:LEU:O	1:A:789:ALA:N	2.44	0.49
1:A:215:TYR:CD1	3:D:8:LEU:HD22	2.48	0.49
1:A:273:LEU:HD23	3:D:44:ILE:HD13	1.94	0.49
2:B:1125:ILE:CG2	2:B:1130:CYS:HB2	2.41	0.49
2:B:1196:ILE:O	2:B:1200:ILE:HD12	2.12	0.49
2:B:1395:GLN:N	2:B:1395:GLN:OE1	2.46	0.49
5:T:47:DT:H2"	5:T:48:DG:C8	2.48	0.49
1:A:290:VAL:O	1:A:294:LYS:N	2.45	0.49
1:A:327:LEU:HD13	1:A:388:ILE:HD11	1.95	0.49
1:A:751:GLU:HA	1:A:754:ILE:HD12	1.95	0.49
1:A:134:LEU:HD12	1:A:157:LYS:HG3	1.95	0.48
1:A:206:ASN:OD1	1:A:207:LEU:N	2.45	0.48
5:T:45:DT:H2"	5:T:46:DC:C5	2.48	0.48
1:A:409:SER:O	1:A:413:MET:N	2.37	0.48
2:B:787:PHE:O	2:B:791:ASN:ND2	2.45	0.48
2:B:1338:LEU:O	2:B:1341:LEU:HB3	2.14	0.48
2:B:174:ARG:NH1	2:B:210:ASP:O	2.46	0.48
2:B:929:ASP:O	2:B:932:VAL:HG12	2.14	0.48
4:S:11:DC:H2"	4:S:12:DA:C8	2.49	0.48
4:S:6:DA:H2"	4:S:7:DG:C8	2.48	0.48
2:B:1310:MET:N	2:B:1311:PRO:CD	2.77	0.47
2:B:500:VAL:HG12	2:B:507:MET:HG2	1.96	0.47
1:A:919:VAL:HG21	1:A:930:PHE:CE2	2.49	0.47
2:B:140:LEU:HD21	2:B:180:LEU:HD23	1.96	0.47
2:B:1125:ILE:HG23	2:B:1130:CYS:HB2	1.97	0.47
1:A:1142:LEU:CD1	1:A:1205:LEU:HD22	2.44	0.47
2:B:173:PRO:HA	2:B:176:ILE:HG22	1.96	0.46
1:A:814:LEU:HD21	1:A:842:ALA:O	2.16	0.46
1:A:868:GLU:N	1:A:868:GLU:OE1	2.47	0.46
2:B:57:ILE:HD13	2:B:135:LEU:HD13	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:502:LEU:HD23	2:B:503:ASN:HB2	1.98	0.46
2:B:1279:ILE:O	2:B:1282:ILE:HG13	2.16	0.46
5:T:43:DG:H2"	5:T:44:DC:C6	2.50	0.46
2:B:1357:VAL:O	2:B:1360:LEU:N	2.49	0.46
1:A:74:GLN:OE1	1:A:78:SER:OG	2.34	0.46
2:B:1024:CYS:HA	2:B:1027:LEU:HD12	1.98	0.46
5:T:44:DC:C6	5:T:45:DT:H72	2.51	0.46
1:A:399:LEU:HD23	1:A:455:PRO:HG2	1.98	0.46
2:B:157:LEU:HD22	2:B:176:ILE:HG21	1.97	0.45
2:B:695:ILE:HG22	2:B:758:CYS:SG	2.57	0.45
2:B:1226:LEU:HD23	2:B:1226:LEU:O	2.15	0.45
5:T:52:DC:H2"	5:T:53:DT:C6	2.51	0.45
3:D:55:THR:N	3:D:58:ASP:OD1	2.47	0.45
1:A:54:SER:O	1:A:65:ARG:NH2	2.50	0.45
2:B:614:LEU:HA	2:B:617:VAL:HG12	1.98	0.45
1:A:810:VAL:HG21	1:A:845:VAL:HG11	1.99	0.45
1:A:134:LEU:HD21	1:A:160:LEU:HD23	1.99	0.45
1:A:210:ILE:O	1:A:213:LEU:HB3	2.17	0.44
1:A:757:ASN:HA	1:A:760:ILE:HD12	1.99	0.44
2:B:927:GLU:OE2	2:B:1004:LEU:HB2	2.17	0.44
1:A:374:SER:O	3:D:48:LYS:NZ	2.40	0.44
2:B:425:TYR:O	2:B:429:LYS:N	2.48	0.44
2:B:1306:LEU:HD21	2:B:1367:LEU:HA	1.99	0.44
1:A:963:GLN:HG3	1:A:1009:TRP:CZ2	2.52	0.44
1:A:195:VAL:O	1:A:199:LEU:HD23	2.17	0.44
1:A:1099:VAL:HG13	1:A:1134:ILE:HG23	2.00	0.44
2:B:1135:ILE:HG12	2:B:1158:LEU:HD23	2.00	0.44
2:B:170:ILE:HB	2:B:175:LEU:HD21	2.00	0.43
2:B:1155:ILE:HG21	2:B:1192:HIS:CD2	2.53	0.43
1:A:1258:GLU:OE1	2:B:1346:LYS:NZ	2.37	0.43
2:B:684:TYR:O	2:B:815:ARG:NH1	2.51	0.43
1:A:430:PHE:CE1	1:A:437:ARG:HB2	2.53	0.43
1:A:116:VAL:HA	1:A:121:LEU:HD11	1.99	0.43
1:A:618:GLN:N	1:A:618:GLN:OE1	2.51	0.43
2:B:1276:SER:O	2:B:1280:ASN:ND2	2.51	0.43
1:A:160:LEU:O	1:A:163:THR:OG1	2.31	0.43
1:A:1283:THR:HG22	2:B:1388:ASN:HD22	1.83	0.43
2:B:1089:SER:O	2:B:1093:VAL:HG23	2.19	0.43
2:B:1303:GLU:HA	2:B:1306:LEU:HD12	1.99	0.43
1:A:1053:ILE:HG22	1:A:1057:LEU:HD12	2.01	0.43
2:B:555:MET:HA	2:B:558:VAL:HG12	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:1342:CYS:SG	2:B:1360:LEU:HD23	2.59	0.43
5:T:52:DC:H2'	5:T:53:DT:H72	2.00	0.43
2:B:762:LEU:HD22	2:B:785:LEU:HD11	2.00	0.42
1:A:998:PRO:HA	1:A:1003:PHE:CD1	2.54	0.42
1:A:987:THR:H	1:A:987:THR:HG1	1.58	0.42
2:B:392:ASN:CB	2:B:395:THR:HG22	2.50	0.42
2:B:466:ASP:OD1	2:B:467:THR:N	2.53	0.42
2:B:414:GLU:O	2:B:417:LEU:N	2.53	0.42
1:A:210:ILE:N	1:A:211:PRO:CD	2.83	0.42
2:B:410:GLY:O	2:B:413:GLN:NE2	2.53	0.42
5:T:45:DT:C2	5:T:46:DC:C4	3.07	0.42
1:A:69:TYR:OH	1:A:96:GLU:O	2.29	0.42
1:A:382:LEU:HD23	1:A:382:LEU:HA	1.91	0.42
5:T:37:DA:H2"	5:T:38:DC:C5	2.55	0.42
1:A:134:LEU:HD21	1:A:160:LEU:CD2	2.50	0.42
1:A:851:GLN:NE2	1:A:851:GLN:HA	2.34	0.42
3:D:19:PRO:O	3:D:55:THR:OG1	2.38	0.42
2:B:1354:THR:HA	2:B:1357:VAL:HG23	2.02	0.41
2:B:1357:VAL:O	2:B:1358:PRO:C	2.59	0.41
4:S:20:DC:H2"	4:S:21:DG:C8	2.55	0.41
5:T:45:DT:H2"	5:T:46:DC:C6	2.55	0.41
2:B:1253:THR:O	2:B:1256:ASP:HB3	2.21	0.41
3:D:7:THR:OG1	3:D:8:LEU:N	2.53	0.41
1:A:648:GLU:N	1:A:648:GLU:OE1	2.53	0.41
1:A:919:VAL:HG21	1:A:930:PHE:HE2	1.83	0.41
1:A:1056:HIS:ND1	1:A:1075:VAL:HG11	2.35	0.41
1:A:1059:ASP:OD1	1:A:1060:ILE:N	2.53	0.41
1:A:97:ALA:O	1:A:136:VAL:HG11	2.21	0.41
2:B:1020:LEU:HD12	2:B:1024:CYS:CB	2.50	0.41
4:S:4:DA:H2"	4:S:5:DC:C6	2.56	0.41
5:T:34:DG:H2"	5:T:35:DG:C8	2.55	0.41
1:A:1151:THR:HG22	1:A:1152:ALA:N	2.36	0.41
1:A:1206:THR:O	1:A:1209:CYS:N	2.53	0.41
2:B:1287:SER:HB2	2:B:1289:PRO:HD2	2.03	0.41
4:S:16:DC:H2"	4:S:17:DA:C8	2.56	0.41
1:A:1250:ILE:HB	1:A:1251:PRO:CD	2.50	0.41
2:B:135:LEU:HD12	2:B:139:LEU:HD13	2.03	0.41
2:B:801:CYS:SG	2:B:802:GLN:N	2.94	0.41
1:A:202:PHE:CE1	1:A:213:LEU:HD21	2.56	0.41
2:B:1275:PHE:CD1	2:B:1301:PHE:CD2	3.09	0.41
3:D:62:GLN:N	3:D:62:GLN:OE1	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:18:GLN:NE2	1:A:58:GLU:OE1	2.54	0.40
2:B:1331:PHE:O	2:B:1334:ASP:OD1	2.39	0.40
1:A:218:LEU:HD11	1:A:275:ILE:HG22	2.03	0.40
1:A:665:ASP:OD1	1:A:665:ASP:N	2.55	0.40
1:A:881:LEU:HD22	1:A:908:CYS:HA	2.02	0.40
4:S:14:DA:H2"	4:S:15:DG:C8	2.56	0.40
1:A:504:GLN:HB3	1:A:546:ASN:HD21	1.86	0.40
2:B:1279:ILE:CD1	2:B:1294:CYS:HB2	2.51	0.40
1:A:112:PHE:HZ	1:A:129:LEU:HD12	1.86	0.40
4:S:26:DG:H2"	4:S:27:DT:C6	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	1180/1328 (89%)	1066 (90%)	109 (9%)	5 (0%)	34	71
2	B	1131/1451 (78%)	1027 (91%)	100 (9%)	4 (0%)	34	71
3	D	74/76 (97%)	71 (96%)	3 (4%)	0	100	100
All	All	2385/2855 (84%)	2164 (91%)	212 (9%)	9 (0%)	38	71

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	451	ARG
1	A	866	ASN
2	B	412	ILE
1	A	724	ASP
1	A	1042	PRO
1	A	374	SER
2	B	1351	THR

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Mol	Chain	Res	Type
2	B	48	PHE
2	B	1126	PRO

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	1093/1205 (91%)	1084 (99%)	9 (1%)	81	89
2	B	1067/1324 (81%)	1059 (99%)	8 (1%)	84	90
3	D	68/68 (100%)	68 (100%)	0	100	100
All	All	2228/2597 (86%)	2211 (99%)	17 (1%)	82	89

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	64	ARG
1	A	71	CYS
1	A	333	SER
1	A	414	PRO
1	A	447	ARG
1	A	611	ASP
1	A	917	SER
1	A	1050	SER
1	A	1285	ARG
2	B	361	SER
2	B	585	MET
2	B	633	ASP
2	B	635	PHE
2	B	770	LYS
2	B	943	LYS
2	B	978	SER
2	B	1021	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	18	GLN
1	A	82	GLN
1	A	342	GLN
1	A	739	ASN
1	A	757	ASN
1	A	865	GLN
1	A	924	GLN
1	A	963	GLN
1	A	1005	GLN
1	A	1017	ASN
1	A	1137	GLN
2	B	65	GLN
2	B	340	GLN
2	B	394	GLN
2	B	545	ASN
2	B	1129	GLN
2	B	1280	ASN
2	B	1388	ASN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

There are no ligands in this entry.

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

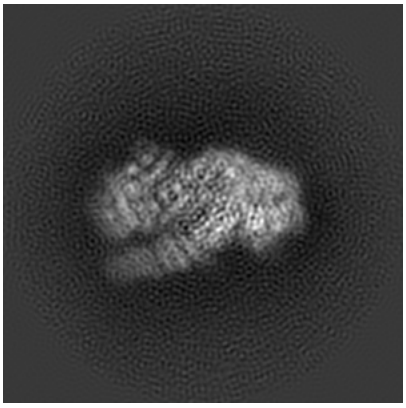
6 Map visualisation (i)

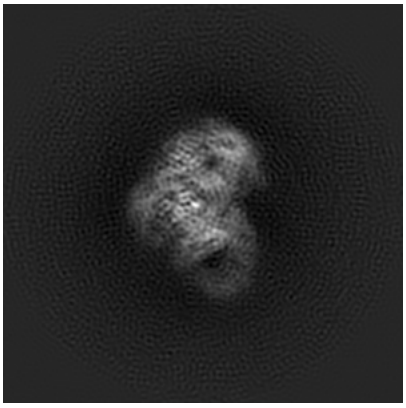
This section contains visualisations of the EMDB entry EMD-21139. These allow visual inspection of the internal detail of the map and identification of artifacts.

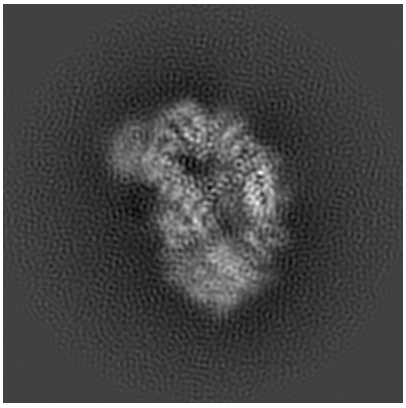
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections (i)

6.1.1 Primary map


X

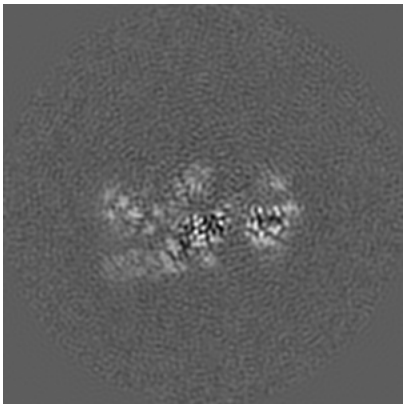

Y

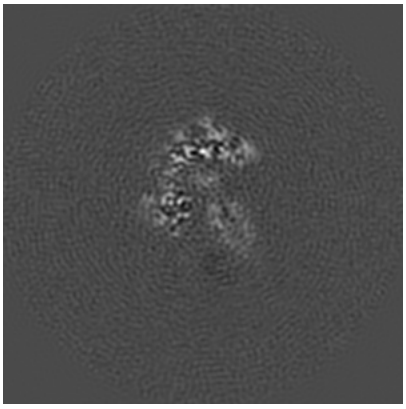

Z

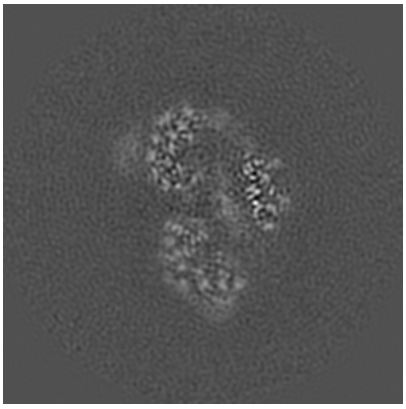
The images above show the map projected in three orthogonal directions.

6.2 Central slices (i)

6.2.1 Primary map


X Index: 128

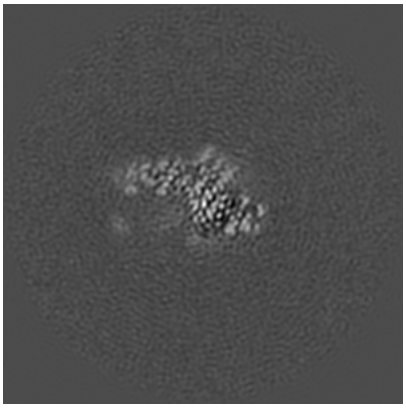

Y Index: 128


Z Index: 128

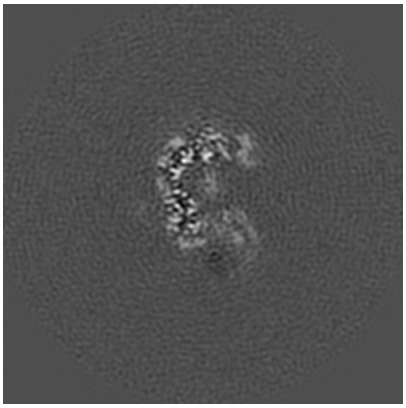
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices (i)

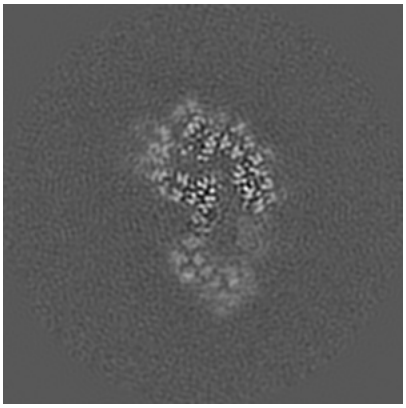
6.3.1 Primary map



X Index: 161



Y Index: 137

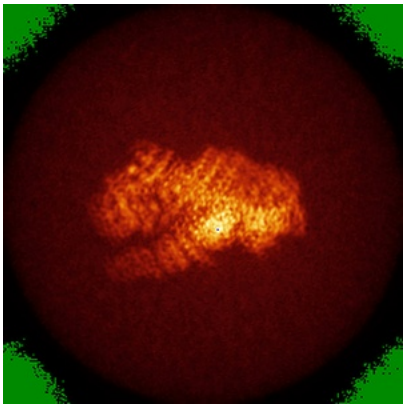


Z Index: 116

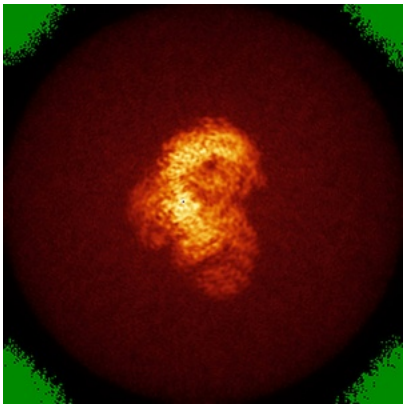
The images above show the largest variance slices of the map in three orthogonal directions.

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

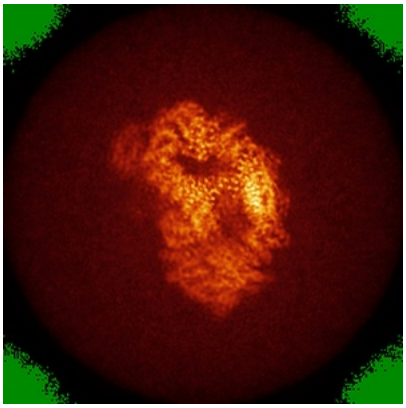
6.4.1 Primary map



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

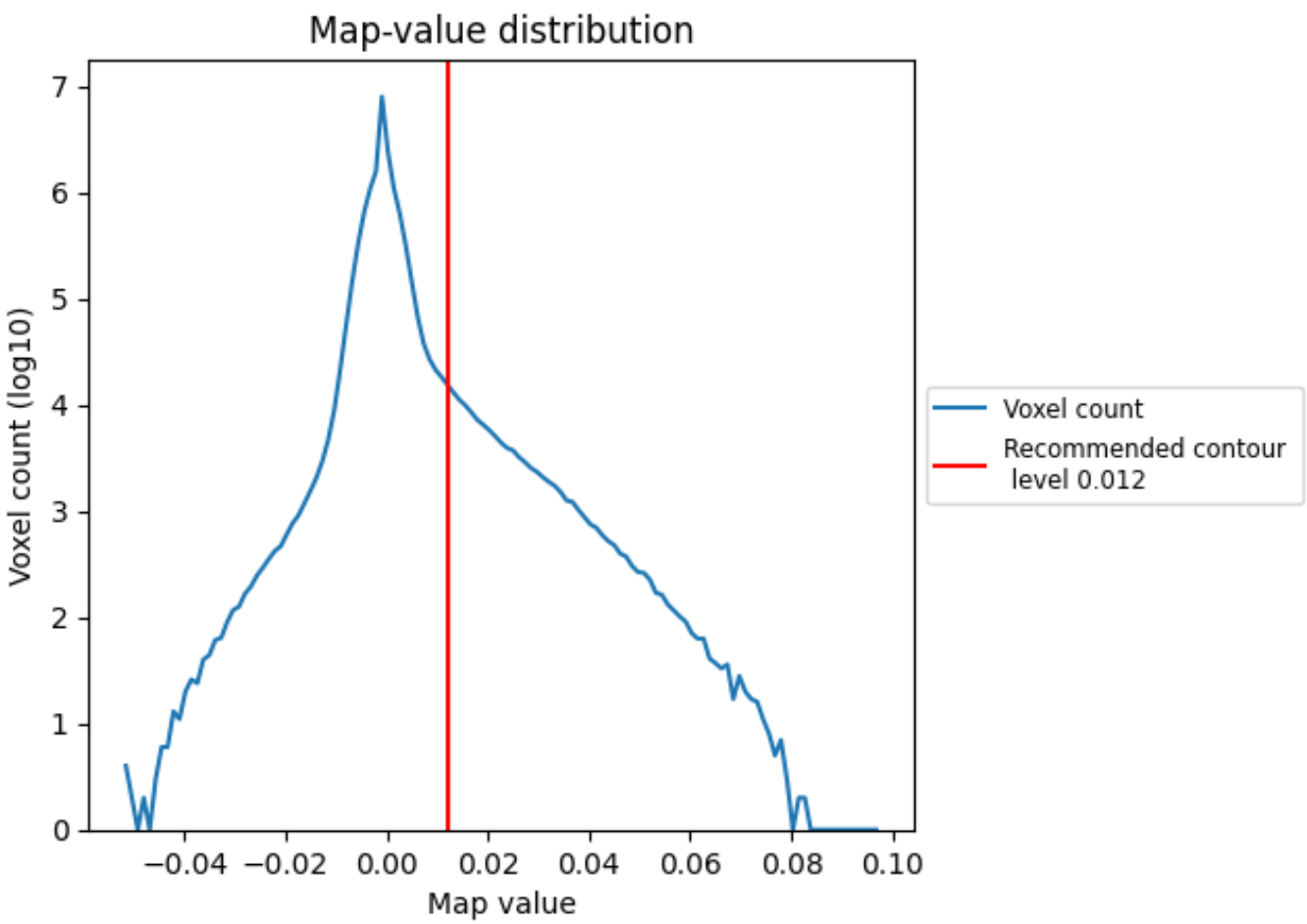
6.6 Mask visualisation (i)

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

7 Map analysis (i)

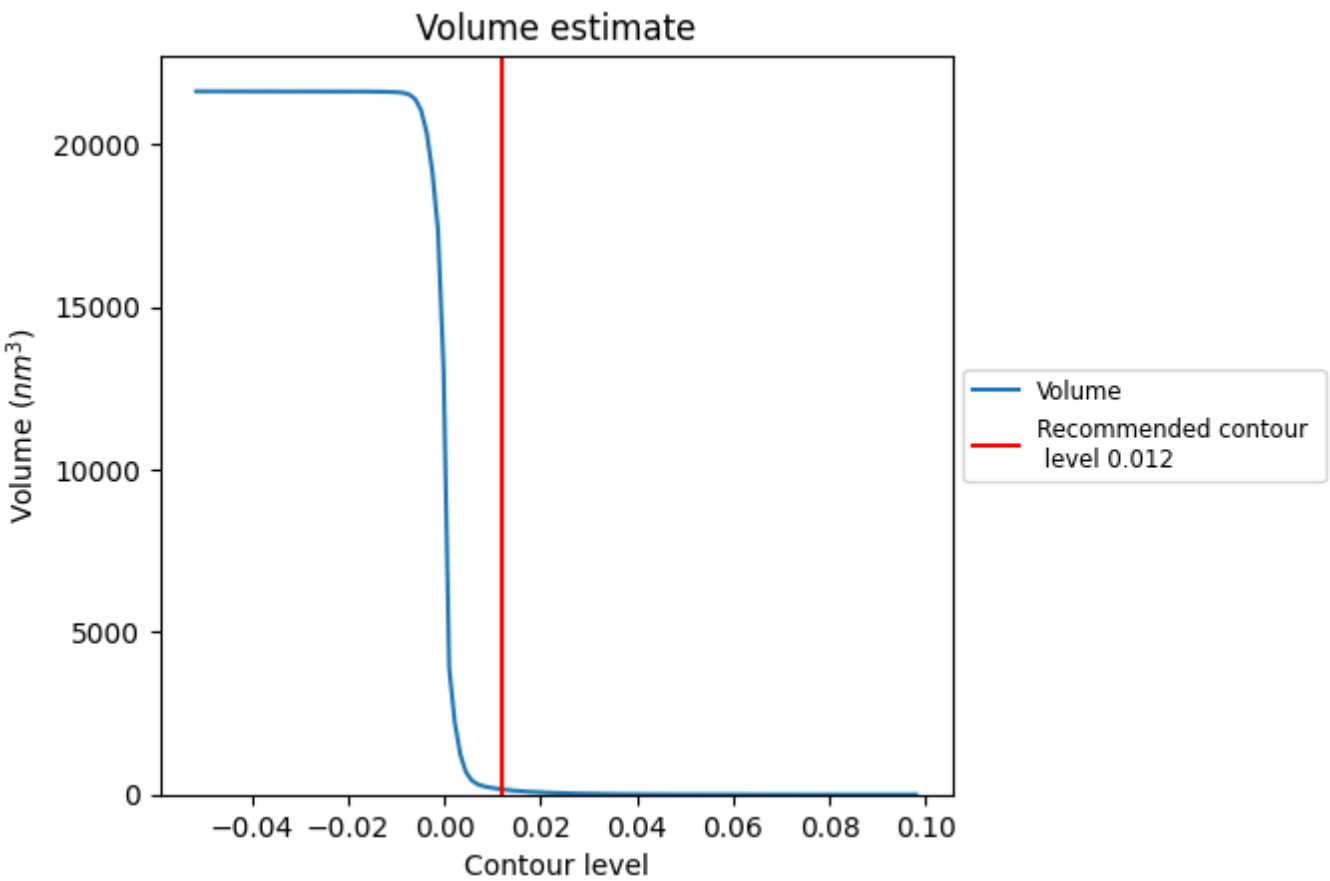
This section contains the results of statistical analysis of the map.

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.

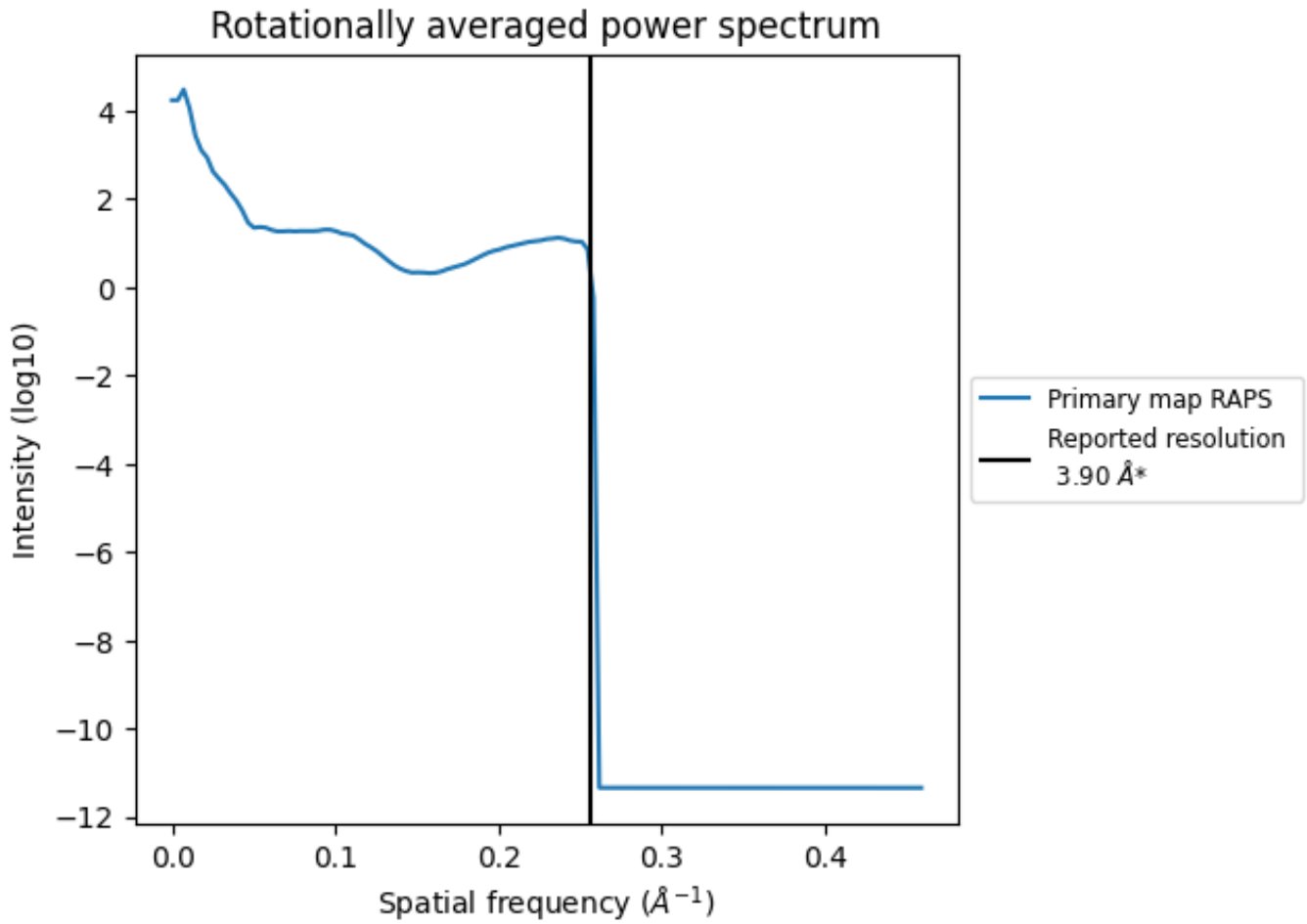
7.2 Volume estimate (i)



The volume at the recommended contour level is 160 nm^3; this corresponds to an approximate mass of 145 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 0.256 Å^{-1}

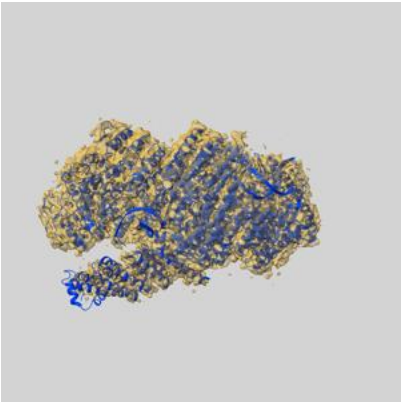
8 Fourier-Shell correlation (i)

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

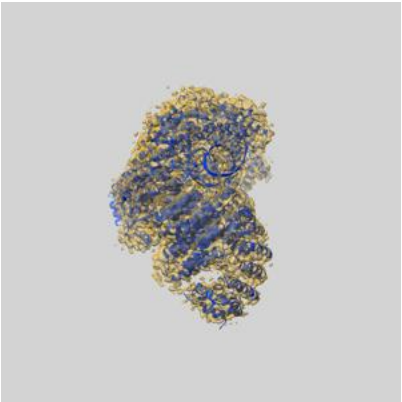
9 Map-model fit (i)

This section contains information regarding the fit between EMDB map EMD-21139 and PDB model 6VAF. Per-residue inclusion information can be found in section 3 on page 5.

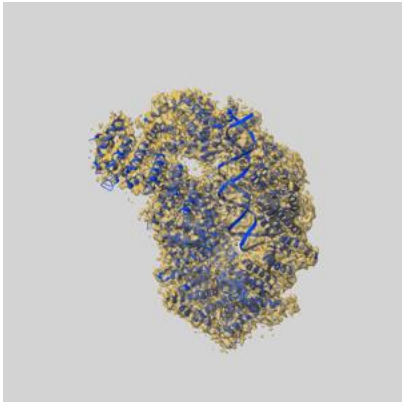
9.1 Map-model overlay (i)



X



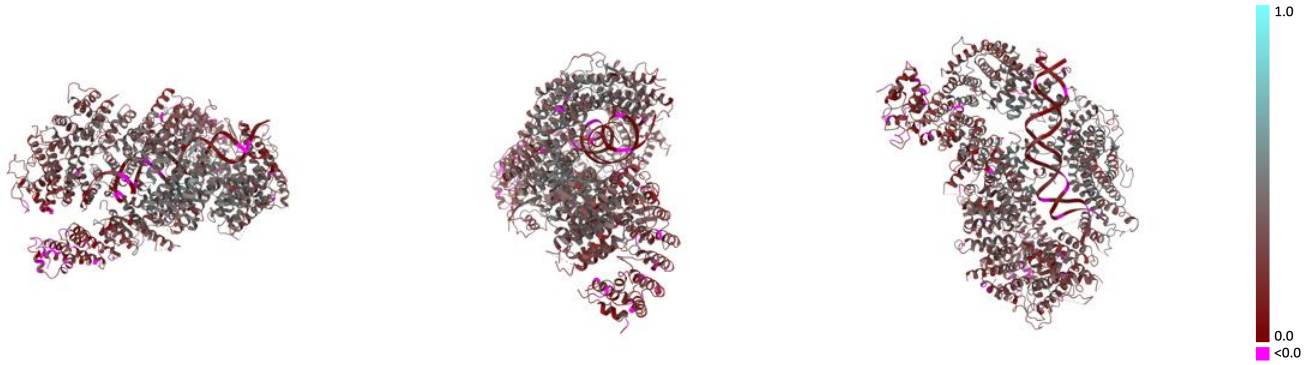
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.012 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.

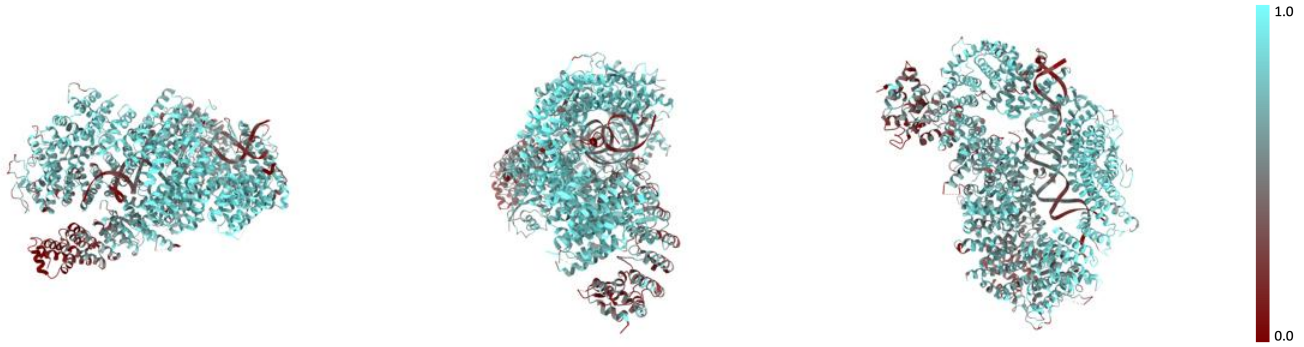
9.2 Q-score mapped to coordinate model (i)



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.

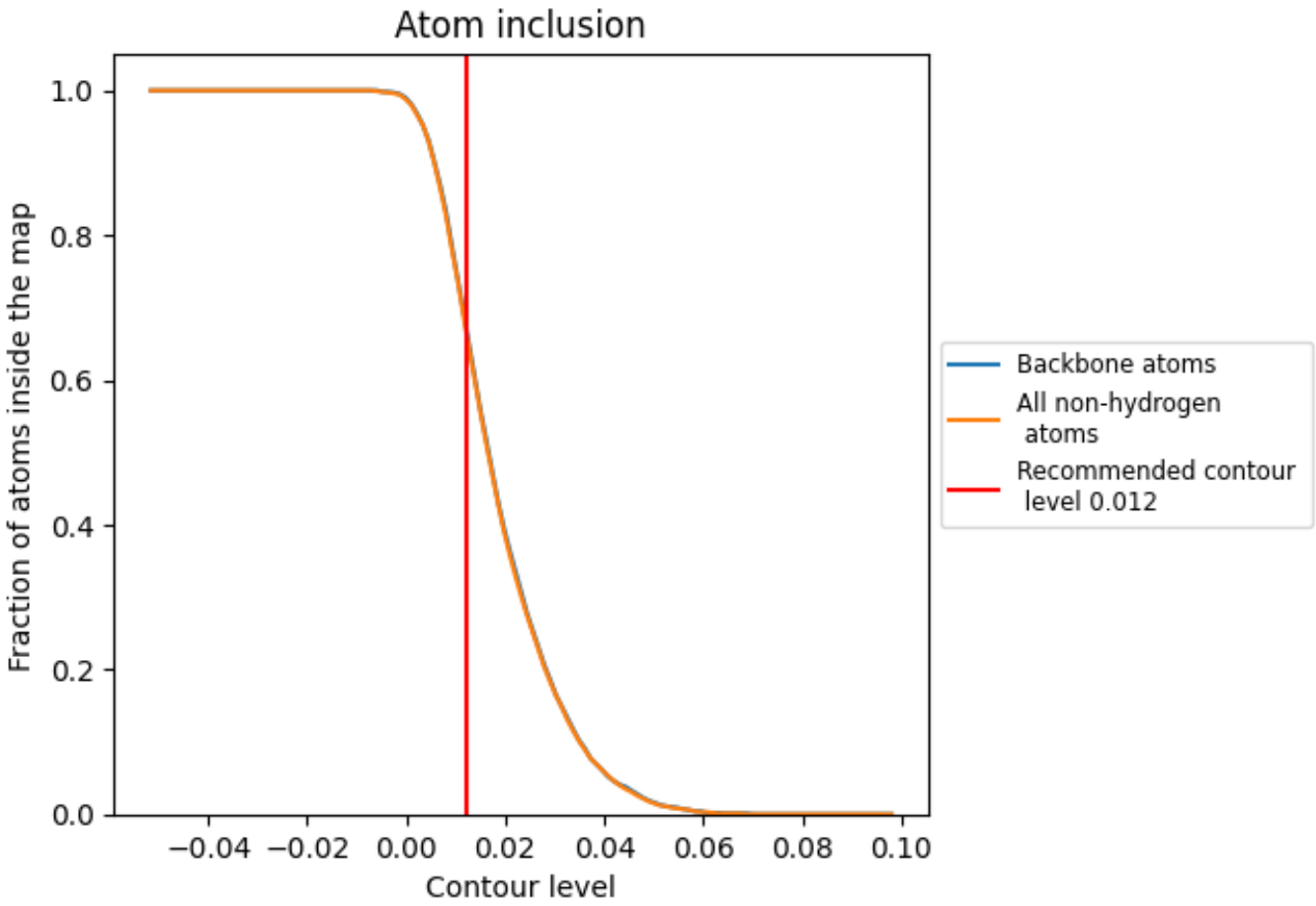
9.3 Atom inclusion mapped to coordinate model (i)



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 (0.012).

9.4 Atom inclusion (i)



At the recommended contour level, 68% of all backbone atoms, 67% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary (i)

The table lists the average atom inclusion at the recommended contour level (0.012) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.6710	0.3030
A	0.7170	0.3400
B	0.6750	0.2920
D	0.6280	0.2800
S	0.3350	0.0880
T	0.3970	0.1100