



Full wwPDB EM Validation Report ⓘ

Nov 7, 2022 – 06:31 PM EST

PDB ID : 6OET
EMDB ID : EMD-20037
Title : Cryo-EM structure of mouse RAG1/2 STC complex
Authors : Chen, X.; Cui, Y.; Zhou, Z.H.; Yang, W.; Gellert, M.
Deposited on : 2019-03-27
Resolution : 3.40 Å(reported)

This is a Full wwPDB EM Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

EMDB validation analysis : 0.0.1.dev43
MolProbity : 4.02b-467
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ : 1.9.9
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.31.2

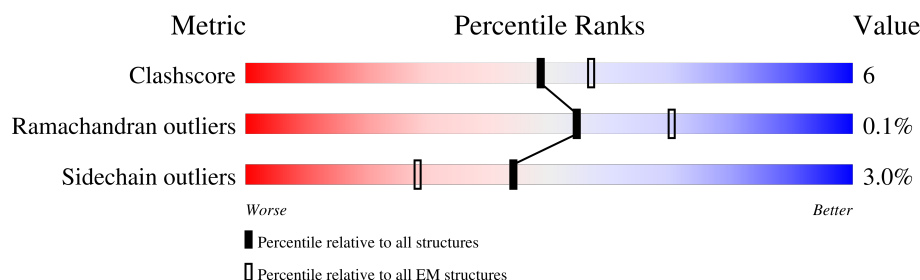
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.40 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 $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion $< 40\%$). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	1040	
1	C	1040	
2	B	527	
2	D	527	
3	F	50	
4	I	15	
4	J	15	
5	G	61	

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Mol	Chain	Length	Quality of chain
6	L	30	<div><div></div><div>20%</div><div>83%</div><div>17%</div></div>
7	M	41	<div><div></div><div>46%</div><div>66%</div><div>27%</div><div>• 5%</div></div>

2 Entry composition [i](#)

There are 9 unique types of molecules in this entry. The entry contains 19604 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 V(D)J recombination-activating protein 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	618	Total	C	N	O	S	1	0
			4969	3127	885	922	35		
1	C	623	Total	C	N	O	S	0	0
			5010	3152	898	926	34		

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	962	GLN	GLU	engineered mutation	UNP P15919
C	962	GLN	GLU	engineered mutation	UNP P15919

- Molecule 2 is a protein called V(D)J recombination-activating protein 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	342	Total	C	N	O	S	2	0
			2691	1718	460	495	18		
2	D	343	Total	C	N	O	S	1	0
			2673	1707	455	493	18		

- Molecule 3 is a DNA chain called DNA (50-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
3	F	50	Total	C	N	O	P	0	0
			1030	489	189	303	49		

- Molecule 4 is a DNA chain called DNA (5'-D(*CP*CP*TP*GP*GP*AP*TP*CP*TP*GP*GP*CP*CP*TP*G)-3').

Mol	Chain	Residues	Atoms					AltConf	Trace
4	I	15	Total	C	N	O	P	0	0
			303	145	53	91	14		
4	J	15	Total	C	N	O	P	0	0
			303	145	53	91	14		

- Molecule 5 is a DNA chain called DNA (59-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
5	G	59	Total	C	N	O	P	0	0
			1205	574	209	363	59		

- Molecule 6 is a DNA chain called DNA (30-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
6	L	30	Total	C	N	O	P	0	0
			611	290	118	173	30		

- Molecule 7 is a DNA chain called DNA (39-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
7	M	39	Total	C	N	O	P	0	0
			805	381	162	223	39		

- Molecule 8 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		AltConf
8	A	1	Total	Ca	0
			1	1	
8	C	1	Total	Ca	0
			1	1	

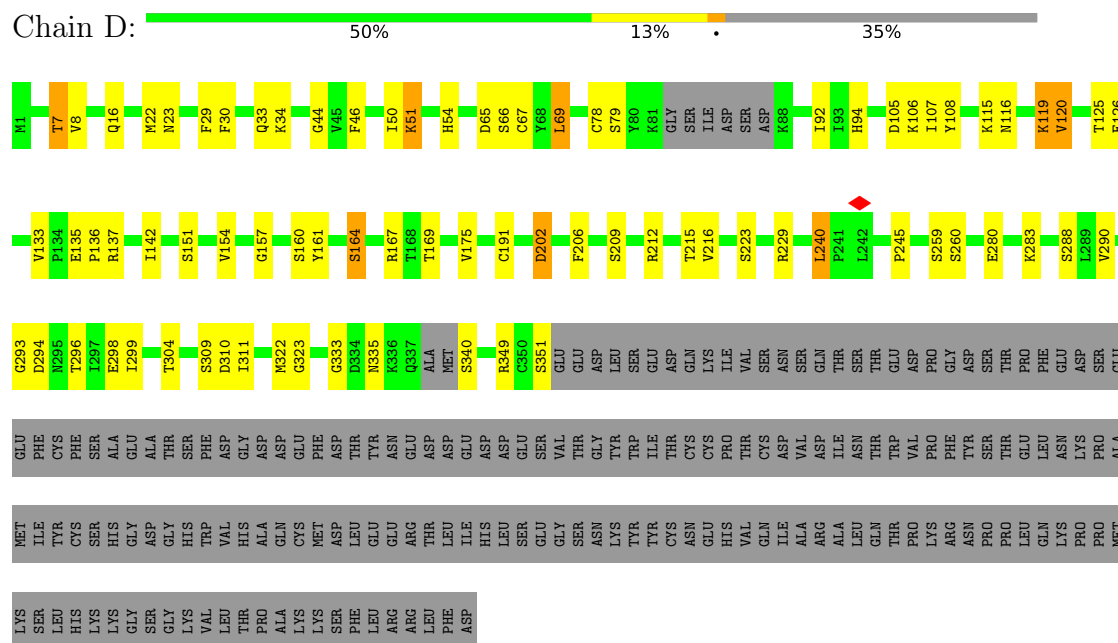
- Molecule 9 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		AltConf
9	A	1	Total	Zn	0
			1	1	
9	C	1	Total	Zn	0
			1	1	

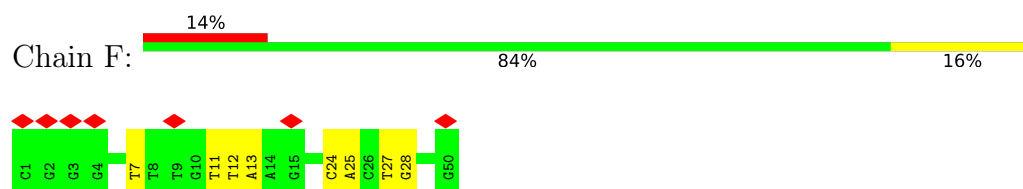


SER	THR	PRO	THR	GLU	LEU	GLN	ASN	LYS	PRO	PRO	ALA	MET	LYS	THR	TYR	CYS	SER	HIS	GLY	ASP	GLY	GLY	HIS	TRP	VAL	HIS	PRO	ALA	GLN	CYS	MET	ASP	LEU	GLU	GLU	ARG	ARG	THR	LEU	ILE	HIS	LEU	SER	GLY	GLY	ASN	LYS	TYR	TYR	CYS	ASN	ASN	GLU	HIS	VAL	GLN	ILE	ALA	ARG	ALA	ALA	GLN	THR	PRO	LYS	ARG	ASN
PRO	PRO	LEU	GLN	LYS	PRO	PRO	MET	LYS	THR	ILE	SER	HIS	LYS	LYS	GLY	SER	ASP	GLY	GLY	VAL	LEU	THR	ALA	ALA	LYS	SER	PHE	LEU	GLU	ARG	ARG	LEU	PHE	ASP																																	

• Molecule 2: V(D)J recombination-activating protein 2



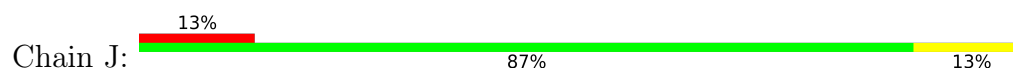
• Molecule 3: DNA (50-MER)



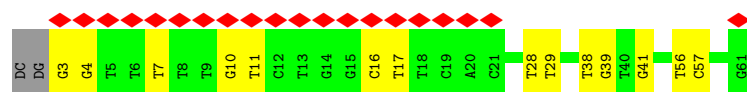
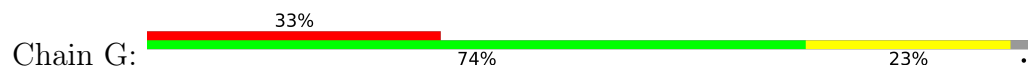
• Molecule 4: DNA (5'-D(*CP*CP*TP*GP*GP*AP*TP*CP*TP*GP*GP*CP*CP*TP*G)-3')



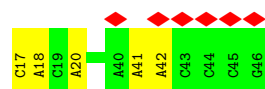
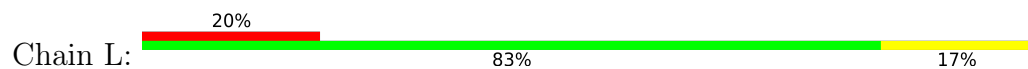
• Molecule 4: DNA (5'-D(*CP*CP*TP*GP*GP*AP*TP*CP*TP*GP*GP*CP*CP*TP*G)-3')



• Molecule 5: DNA (59-MER)



• Molecule 6: DNA (30-MER)



• Molecule 7: DNA (39-MER)



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	68085	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 ($e^-/\text{\AA}^2$)	45	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	0.216	Depositor
Minimum map value	-0.098	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.006	Depositor
Recommended contour level	0.032	Depositor
Map size (Å)	299.6, 299.6, 299.6	wwPDB
Map dimensions	280, 280, 280	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.07, 1.07, 1.07	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: CA, 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	0.52	0/5072	0.57	0/6834
1	C	0.51	0/5114	0.57	0/6887
2	B	0.51	0/2759	0.55	0/3739
2	D	0.51	0/2741	0.55	0/3719
3	F	1.09	0/1155	1.03	0/1784
4	I	0.98	0/338	1.04	0/520
4	J	0.95	0/338	1.02	0/520
5	G	0.98	0/1347	1.02	0/2077
6	L	1.07	1/686 (0.1%)	0.96	0/1052
7	M	1.04	1/907 (0.1%)	1.00	0/1395
All	All	0.67	2/20457 (0.0%)	0.71	0/28527

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	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	L	17	DC	OP3-P	-10.97	1.48	1.61
7	M	17	DC	OP3-P	-10.15	1.49	1.61

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	520	GLN	Peptide

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	4969	0	4939	62	0
1	C	5010	0	4991	74	0
2	B	2691	0	2635	52	0
2	D	2673	0	2599	43	0
3	F	1030	0	565	6	0
4	I	303	0	171	3	0
4	J	303	0	171	1	0
5	G	1205	0	668	8	0
6	L	611	0	335	4	0
7	M	805	0	435	8	0
8	A	1	0	0	0	0
8	C	1	0	0	0	0
9	A	1	0	0	0	0
9	C	1	0	0	0	0
All	All	19604	0	17509	229	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 6.

All (229) 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:69:LEU:HD23	2:B:69:LEU:O	1.88	0.74
2:D:69:LEU:HD23	2:D:69:LEU:O	1.90	0.70
1:A:845:PRO:HG2	2:D:335:ASN:ND2	2.08	0.69
1:A:473:ASN:HD22	1:A:1003:MET:HG3	1.58	0.68
1:C:405:LYS:O	1:C:409:ARG:HB3	1.94	0.68
2:B:235:ARG:HD2	2:B:237[A]:ARG:HE	1.61	0.65
2:D:116:ASN:HB2	2:D:119:LYS:HB2	1.79	0.64
2:B:116:ASN:HB2	2:B:119:LYS:HB2	1.79	0.63
2:D:78:CYS:HB3	2:D:92:ILE:HB	1.81	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:23:ASN:ND2	2:B:322:MET:O	2.34	0.61
2:B:335:ASN:OD1	2:B:337:GLN:HB3	2.00	0.61
2:D:23:ASN:ND2	2:D:322:MET:O	2.34	0.61
2:B:78:CYS:HB3	2:B:92:ILE:HB	1.82	0.61
1:A:525:ASN:ND2	2:B:167:ARG:O	2.35	0.60
1:A:391:ARG:NH1	7:M:52:DA:N3	2.49	0.60
1:C:767:GLU:HB3	1:C:771:GLU:HG3	1.83	0.60
2:D:283:LYS:NZ	2:D:311:ILE:O	2.33	0.59
2:B:115:LYS:HG3	2:B:120:VAL:HG23	1.84	0.59
2:B:335:ASN:ND2	1:C:845:PRO:HG3	2.17	0.59
1:C:402:ARG:NH1	3:F:11:DT:O4	2.35	0.59
1:A:608:LYS:NZ	1:A:983:GLU:OE1	2.34	0.59
1:A:713:ARG:NH1	1:A:720:ALA:O	2.36	0.59
1:A:729:LEU:HB3	1:A:746:ILE:HG23	1.85	0.59
1:A:595:VAL:HG12	1:A:626:VAL:HA	1.85	0.58
2:D:115:LYS:HG3	2:D:120:VAL:HG23	1.84	0.58
2:B:202:ASP:OD1	2:B:202:ASP:N	2.37	0.58
2:B:283:LYS:NZ	2:B:311:ILE:O	2.33	0.58
2:B:69:LEU:CD2	2:B:69:LEU:H	2.17	0.58
2:D:202:ASP:OD1	2:D:202:ASP:N	2.37	0.58
1:C:730:CYS:O	1:C:748:ARG:NH1	2.37	0.58
1:C:534:ILE:HG13	1:C:557:PHE:HE1	1.68	0.57
1:C:915:ARG:NH1	1:C:918:GLU:OE2	2.38	0.57
1:A:723:SER:O	1:A:734:ARG:NH1	2.37	0.57
1:C:406:HIS:O	1:C:409:ARG:NH1	2.34	0.57
1:A:621:ARG:NH1	1:A:965:ASN:OD1	2.37	0.57
1:C:393:ARG:NH2	3:F:7:DT:OP1	2.38	0.57
1:A:426:ASP:OD1	1:A:426:ASP:N	2.36	0.56
1:A:560:ASP:N	1:A:560:ASP:OD1	2.37	0.56
5:G:56:DT:H1'	5:G:57:DC:H5'	1.86	0.56
1:A:872:SER:OG	1:A:873:GLU:N	2.38	0.56
1:A:471:ARG:NH1	1:A:478:CYS:SG	2.79	0.56
1:C:525:ASN:ND2	2:D:167:ARG:O	2.38	0.56
1:A:713:ARG:NH2	1:A:726:ILE:O	2.38	0.56
1:A:541:LEU:HG	2:B:173:ASN:HD22	1.71	0.55
1:C:461:GLY:O	1:C:498:GLN:NE2	2.39	0.55
1:C:471:ARG:NH1	1:C:478:CYS:SG	2.79	0.55
2:B:23:ASN:HD22	2:B:323:GLY:HA3	1.70	0.55
1:C:560:ASP:OD1	1:C:560:ASP:N	2.40	0.55
1:C:625:THR:HG22	1:C:652:CYS:HA	1.89	0.55
2:D:23:ASN:HD22	2:D:323:GLY:HA3	1.70	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:812:ILE:HA	1:C:870:ILE:HG22	1.89	0.55
2:B:310:ASP:OD2	1:C:828:ARG:NH1	2.40	0.55
1:A:391:ARG:NH2	5:G:7:DT:O2	2.41	0.54
1:A:546:ASP:OD2	2:B:259:SER:OG	2.26	0.54
2:B:94:HIS:HB2	2:B:107:ILE:HG12	1.89	0.54
1:C:406:HIS:HA	1:C:409:ARG:HD3	1.89	0.54
2:B:313:HIS:O	1:C:839:LYS:NZ	2.40	0.54
1:C:579:ARG:NH1	1:C:585:ASP:OD2	2.41	0.54
1:C:546:ASP:OD2	2:D:229:ARG:NH2	2.37	0.54
2:B:154:VAL:HG11	2:B:216:VAL:HG21	1.90	0.54
1:A:442:ARG:HG2	1:C:428:LYS:HD3	1.90	0.54
2:D:105:ASP:HB3	2:D:136:PRO:HG3	1.90	0.54
1:C:761:ARG:NH2	2:D:108:TYR:OH	2.41	0.54
2:D:310:ASP:OD1	2:D:310:ASP:N	2.41	0.54
1:A:769:VAL:HG22	2:B:68:TYR:OH	2.08	0.54
1:A:411:LEU:HD21	1:C:415:VAL:HG12	1.89	0.53
2:B:105:ASP:HB3	2:B:136:PRO:HG3	1.90	0.53
2:D:94:HIS:HB2	2:D:107:ILE:HG12	1.89	0.53
1:C:713:ARG:NH2	1:C:720:ALA:O	2.41	0.53
1:A:966:LYS:NZ	3:F:28:DG:N3	2.52	0.52
1:C:653:LYS:NZ	1:C:995:THR:O	2.41	0.52
2:D:154:VAL:HG11	2:D:216:VAL:HG21	1.90	0.52
2:B:212:ARG:NH1	2:B:293:GLY:O	2.42	0.52
2:D:212:ARG:NH1	2:D:293:GLY:O	2.42	0.52
2:D:290:VAL:HG22	2:D:299:ILE:HG12	1.92	0.52
1:A:650:LEU:HB3	1:A:1000:GLN:HE22	1.75	0.52
1:A:790:SER:OG	1:A:791:ILE:N	2.42	0.52
1:A:585:ASP:O	1:A:696:ARG:NH1	2.39	0.52
2:B:290:VAL:HG22	2:B:299:ILE:HG12	1.92	0.52
1:A:730:CYS:O	1:A:748:ARG:NH1	2.43	0.52
1:A:531:ASP:N	1:A:531:ASP:OD1	2.43	0.52
1:A:654:PRO:HG3	1:A:991:HIS:HB3	1.91	0.52
1:C:691:MET:HB3	1:C:696:ARG:HE	1.75	0.52
2:D:105:ASP:OD1	2:D:105:ASP:N	2.39	0.52
2:B:105:ASP:OD1	2:B:105:ASP:N	2.39	0.51
2:D:66:SER:OG	2:D:67:CYS:N	2.43	0.51
1:C:708:ASP:N	1:C:708:ASP:OD1	2.37	0.51
2:B:66:SER:OG	2:B:67:CYS:N	2.43	0.51
1:C:729:LEU:HB3	1:C:746:ILE:HG23	1.92	0.51
1:A:477:SER:OG	1:A:478:CYS:N	2.43	0.51
1:A:645:LYS:NZ	6:L:20:DA:OP2	2.44	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:678:GLU:HG3	2:D:169:THR:HG21	1.92	0.51
2:D:349:ARG:NH1	2:D:351:SER:O	2.43	0.51
2:B:310:ASP:OD1	2:B:310:ASP:N	2.41	0.50
2:D:160:SER:OG	2:D:161:TYR:N	2.44	0.50
1:C:551:ASP:OD1	1:C:551:ASP:N	2.43	0.50
2:B:160:SER:OG	2:B:161:TYR:N	2.44	0.50
1:C:805:TYR:OH	1:C:834:ASP:OD1	2.30	0.50
7:M:38:DT:H2'	7:M:39:DG:C8	2.48	0.49
1:C:666:GLU:OE1	2:D:34:LYS:NZ	2.41	0.49
1:A:588:ASN:HA	1:A:696:ARG:HG2	1.94	0.49
1:C:608:LYS:NZ	1:C:978:GLN:OE1	2.45	0.49
2:B:65:ASP:OD1	2:B:65:ASP:N	2.45	0.48
1:A:541:LEU:HD11	1:A:670:ALA:HB1	1.93	0.48
2:D:16:GLN:NE2	2:D:33:GLN:O	2.47	0.48
1:C:596:LYS:NZ	1:C:957:ALA:O	2.39	0.48
2:B:16:GLN:NE2	2:B:33:GLN:O	2.47	0.48
1:C:606:SER:O	1:C:972:ARG:NH2	2.46	0.48
1:C:740:ASN:HD21	1:C:743:PHE:HB2	1.78	0.48
1:C:462:LEU:O	1:C:498:GLN:NE2	2.38	0.47
2:B:237[A]:ARG:NH2	2:B:295:ASN:OD1	2.48	0.47
2:B:240:LEU:HD22	2:B:245:PRO:HB3	1.96	0.47
1:A:477:SER:HB3	1:A:480:GLN:HG2	1.95	0.47
1:A:430:VAL:HG12	1:C:396:LEU:HD23	1.95	0.47
1:C:723:SER:OG	1:C:725:TYR:O	2.32	0.47
1:A:414:GLN:HA	1:A:417:GLU:HG2	1.96	0.47
2:B:51:LYS:HB3	2:B:51:LYS:HE3	1.61	0.47
2:D:240:LEU:HD22	2:D:245:PRO:HB3	1.96	0.47
7:M:44:DC:H2'	7:M:45:DA:C8	2.50	0.47
1:A:933:THR:HG22	4:I:14:DC:H5''	1.96	0.47
2:D:33:GLN:HE21	2:D:44:GLY:HA2	1.80	0.47
1:C:713:ARG:NH2	1:C:719:GLU:O	2.48	0.47
2:B:135:GLU:H	2:B:137:ARG:HH21	1.63	0.46
1:A:405:LYS:O	1:A:409:ARG:HB2	2.14	0.46
2:B:33:GLN:HE21	2:B:44:GLY:HA2	1.81	0.46
1:A:539:SER:HB3	2:B:172:TRP:H	1.81	0.46
1:C:748:ARG:NH2	1:C:777:LYS:O	2.49	0.46
1:A:727:CYS:HB2	1:A:937:HIS:CD2	2.50	0.46
2:B:164:SER:O	2:B:164:SER:OG	2.33	0.46
1:A:740:ASN:ND2	1:A:743:PHE:O	2.47	0.46
2:D:22:MET:HG3	2:D:29:PHE:HB2	1.98	0.46
5:G:3:DG:H2''	5:G:4:DG:H5''	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:433:THR:HA	1:C:436:LEU:HD12	1.96	0.46
1:C:598:SER:HB2	1:C:706:GLY:HA3	1.98	0.46
1:C:853:PHE:O	1:C:857:LEU:HB2	2.15	0.46
2:D:135:GLU:H	2:D:137:ARG:HH21	1.63	0.46
2:B:22:MET:HG3	2:B:29:PHE:HB2	1.98	0.46
1:A:719:GLU:OE2	1:A:773:ARG:NE	2.43	0.45
1:C:443:ASN:HD21	7:M:42:DG:H21	1.62	0.45
1:C:894:ARG:HA	1:C:894:ARG:HD2	1.68	0.45
1:C:989:LYS:HB3	1:C:989:LYS:HE3	1.81	0.45
7:M:21:DG:H2'	7:M:22:DT:H71	1.98	0.45
1:C:391:ARG:HB2	6:L:41:DA:H2	1.80	0.45
1:C:962:GLN:HB2	5:G:41:DG:H21	1.82	0.45
2:D:137:ARG:HH11	2:D:157:GLY:HA2	1.82	0.45
1:A:558:ARG:NH2	1:A:560:ASP:OD2	2.40	0.45
1:C:863:VAL:HA	1:C:866:VAL:HG12	1.98	0.45
2:B:137:ARG:HH11	2:B:157:GLY:HA2	1.82	0.45
1:C:410:GLU:HA	1:C:413:ILE:HD12	1.98	0.45
1:C:588:ASN:HA	1:C:696:ARG:HG2	2.00	0.45
1:A:404:GLN:HG3	1:C:437:LEU:HD11	1.99	0.44
1:A:703:ARG:NH2	1:A:950:ASP:OD2	2.50	0.44
1:A:576:GLU:OE2	1:A:579:ARG:NE	2.45	0.44
1:C:752:GLU:OE2	1:C:756:ARG:NH2	2.38	0.44
1:A:482:HIS:HB2	1:A:500:LEU:HD11	2.00	0.44
1:A:754:LEU:HD23	1:A:783:PRO:HD2	1.98	0.44
1:C:833:LEU:O	1:C:837:LEU:HB2	2.18	0.44
1:A:792:ASP:OD2	1:A:795:HIS:ND1	2.37	0.44
5:G:28:DT:H2'	5:G:29:DT:C6	2.53	0.44
1:C:639:LYS:HE2	1:C:639:LYS:HB3	1.78	0.44
2:D:51:LYS:HB3	2:D:51:LYS:HE3	1.61	0.44
5:G:10:DG:H3'	5:G:11:DT:H71	2.00	0.44
2:B:7:THR:OG1	2:B:8:VAL:N	2.51	0.44
1:C:517:PHE:HB2	1:C:687:LEU:HD11	1.98	0.44
2:B:69:LEU:CD2	2:B:69:LEU:N	2.79	0.43
1:C:752:GLU:HG2	1:C:756:ARG:HE	1.82	0.43
2:B:106:LYS:NZ	2:B:126:GLU:OE2	2.43	0.43
1:C:727:CYS:HB2	1:C:937:HIS:CD2	2.54	0.43
1:C:952:SER:OG	1:C:953:ILE:N	2.51	0.43
2:B:50:ILE:HA	2:B:54:HIS:O	2.19	0.43
1:C:680:GLU:HA	1:C:683:LYS:HD3	2.01	0.43
3:F:12:DT:H2'	3:F:13:DA:C8	2.53	0.43
4:I:6:DG:H2''	4:I:7:DA:C8	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:50:ILE:HA	2:D:54:HIS:O	2.19	0.43
1:C:891:PRO:HB3	7:M:17:DC:N3	2.33	0.43
7:M:18:DA:C8	7:M:18:DA:H5'	2.54	0.42
2:B:206:PHE:HB3	2:B:260:SER:HB3	2.01	0.42
2:D:206:PHE:HB3	2:D:260:SER:HB3	2.02	0.42
2:D:78:CYS:HB2	2:D:142:ILE:HG23	2.01	0.42
6:L:18:DA:H5'	6:L:18:DA:C8	2.54	0.42
1:C:823:LYS:HE3	1:C:823:LYS:HB3	1.87	0.42
2:B:212:ARG:HH12	2:B:294:ASP:HA	1.85	0.42
1:C:707:TYR:HD2	1:C:711:LEU:HD23	1.84	0.42
7:M:54:DC:H2'	7:M:55:DC:C6	2.54	0.42
1:A:797:ASP:OD2	1:A:894:ARG:NE	2.43	0.42
1:A:853:PHE:O	1:A:857:LEU:HB2	2.19	0.42
2:D:7:THR:OG1	2:D:8:VAL:N	2.51	0.42
4:J:6:DG:H2''	4:J:7:DA:C8	2.54	0.42
5:G:38:DT:H2'	5:G:39:DG:C8	2.54	0.42
2:B:30:PHE:HB3	2:B:46:PHE:HB2	2.02	0.42
2:B:69:LEU:H	2:B:69:LEU:HD22	1.85	0.42
2:B:69:LEU:HD23	2:B:69:LEU:N	2.35	0.41
2:B:78:CYS:HB2	2:B:142:ILE:HG23	2.01	0.41
2:D:212:ARG:HH12	2:D:294:ASP:HA	1.85	0.41
4:I:14:DC:H2'	4:I:15:DT:C6	2.55	0.41
1:A:629:ILE:HB	1:A:641:PHE:HB3	2.01	0.41
2:B:179:LEU:HA	2:B:180:PRO:HD3	1.84	0.41
1:C:790:SER:OG	1:C:791:ILE:N	2.50	0.41
5:G:16:DC:H2'	5:G:17:DT:C6	2.55	0.41
1:A:415:VAL:HA	1:C:411:LEU:HD11	2.02	0.41
1:A:727:CYS:O	1:A:748:ARG:NH1	2.53	0.41
1:C:789:PRO:O	1:C:949:ARG:NH1	2.41	0.41
2:D:164:SER:O	2:D:164:SER:OG	2.33	0.41
1:A:648:SER:OG	1:A:649:GLU:N	2.53	0.41
1:A:543:SER:O	1:A:543:SER:OG	2.31	0.41
2:B:14:LEU:HG	2:B:33:GLN:HB2	2.03	0.41
2:B:219:LEU:HD12	2:B:256:ILE:HG12	2.03	0.41
1:C:555:LYS:HD3	1:C:555:LYS:HA	1.84	0.41
1:C:653:LYS:HE3	1:C:653:LYS:HB3	1.76	0.41
1:A:684:SER:O	1:A:684:SER:OG	2.39	0.41
1:A:1006:HIS:CE1	1:C:486:ARG:HD2	2.55	0.41
2:B:117:ASN:O	2:B:119:LYS:NZ	2.35	0.41
1:C:672:LEU:HD12	1:C:675:LEU:HD12	2.03	0.41
2:D:92:ILE:HD13	2:D:92:ILE:HA	1.94	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:106:LYS:NZ	2:D:126:GLU:OE2	2.43	0.41
1:A:803:GLU:OE2	1:A:928:TYR:OH	2.26	0.40
1:A:870:ILE:HG21	1:A:875:ARG:HB2	2.03	0.40
1:C:909:TYR:O	1:C:913:SER:OG	2.34	0.40
2:D:30:PHE:HB3	2:D:46:PHE:HB2	2.02	0.40
1:A:653:LYS:HZ3	1:A:653:LYS:HG2	1.77	0.40
1:A:768:SER:OG	1:A:771:GLU:OE1	2.33	0.40
1:C:1003:MET:O	1:C:1004:ASN:ND2	2.54	0.40
3:F:24:DC:H2'	3:F:25:DA:C8	2.56	0.40
6:L:41:DA:H2''	6:L:42:DA:C8	2.57	0.40
1:C:545:VAL:N	2:D:280:GLU:OE2	2.54	0.40
2:D:65:ASP:OD1	2:D:65:ASP:N	2.45	0.40
2:D:116:ASN:HB2	2:D:119:LYS:HD2	2.04	0.40
3:F:27:DT:H2'	3:F:28:DG:C8	2.56	0.40
1:A:927:ARG:HA	1:A:927:ARG:HD2	1.82	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	617/1040 (59%)	586 (95%)	31 (5%)	0	100	100
1	C	621/1040 (60%)	586 (94%)	35 (6%)	0	100	100
2	B	338/527 (64%)	316 (94%)	21 (6%)	1 (0%)	41	72
2	D	338/527 (64%)	315 (93%)	22 (6%)	1 (0%)	41	72
All	All	1914/3134 (61%)	1803 (94%)	109 (6%)	2 (0%)	54	82

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	333	GLY
2	D	333	GLY

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	543/933 (58%)	541 (100%)	2 (0%)	91	95
1	C	547/933 (59%)	547 (100%)	0	100	100
2	B	299/469 (64%)	273 (91%)	26 (9%)	10	34
2	D	295/469 (63%)	271 (92%)	24 (8%)	11	38
All	All	1684/2804 (60%)	1632 (97%)	52 (3%)	44	68

All (52) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	412	LYS
1	A	761	ARG
2	B	7	THR
2	B	51	LYS
2	B	69	LEU
2	B	79	SER
2	B	119	LYS
2	B	120	VAL
2	B	125	THR
2	B	133	VAL
2	B	151	SER
2	B	164	SER
2	B	175	VAL
2	B	191	CYS
2	B	202	ASP
2	B	209	SER
2	B	215	THR
2	B	223	SER
2	B	237[A]	ARG
2	B	237[B]	ARG

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Mol	Chain	Res	Type
2	B	240	LEU
2	B	259	SER
2	B	288	SER
2	B	296	THR
2	B	298	GLU
2	B	304	THR
2	B	309	SER
2	B	340	SER
2	D	7	THR
2	D	51	LYS
2	D	69	LEU
2	D	79	SER
2	D	119	LYS
2	D	120	VAL
2	D	125	THR
2	D	133	VAL
2	D	151	SER
2	D	164	SER
2	D	175	VAL
2	D	191	CYS
2	D	202	ASP
2	D	209	SER
2	D	215	THR
2	D	223	SER
2	D	240	LEU
2	D	259	SER
2	D	288	SER
2	D	296	THR
2	D	298	GLU
2	D	304	THR
2	D	309	SER
2	D	340	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	404	GLN
1	A	473	ASN
1	A	753	ASN
1	A	830	GLN
1	A	978	GLN
1	A	1000	GLN

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Mol	Chain	Res	Type
1	A	1006	HIS
1	A	1007	ASN
2	B	33	GLN
2	B	53	ASN
2	B	101	ASN
2	B	173	ASN
2	B	181	HIS
2	B	337	GLN
1	C	965	ASN
1	C	1004	ASN
2	D	10	HIS
2	D	33	GLN
2	D	181	HIS
2	D	337	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 4 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.

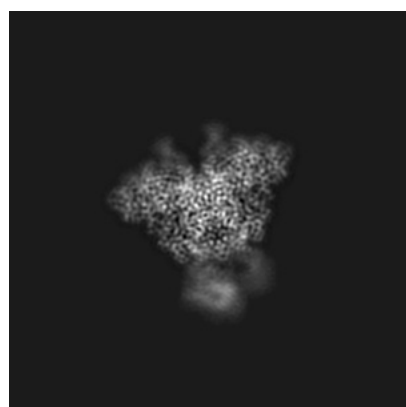
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-20037. These allow visual inspection of the internal detail of the map and identification of artifacts.

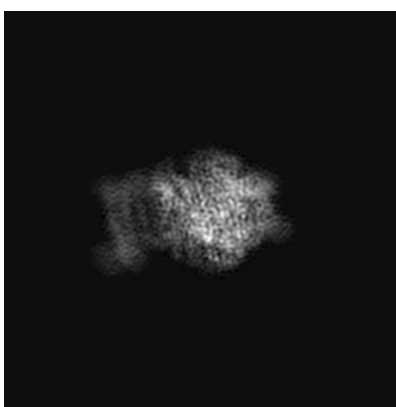
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

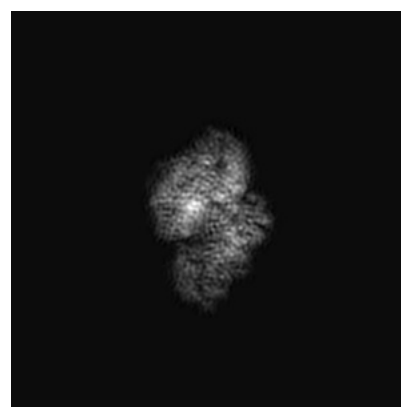
6.1.1 Primary map



X



Y

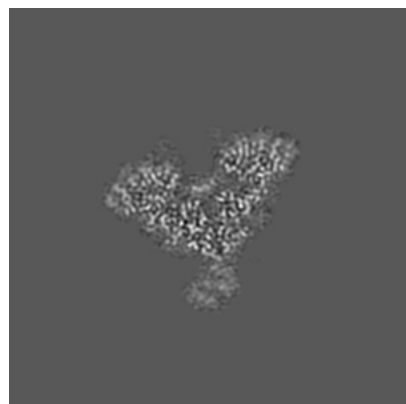


Z

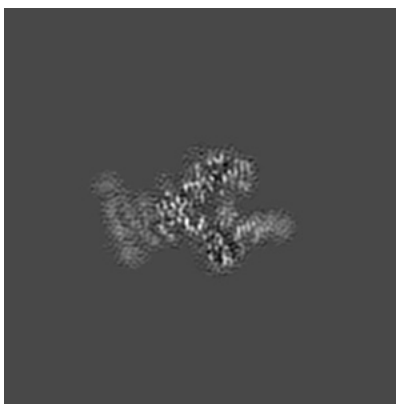
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

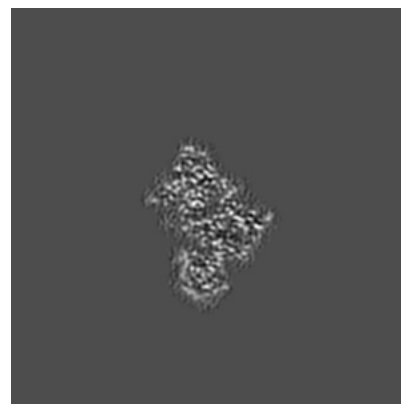
6.2.1 Primary map



X Index: 140



Y Index: 140

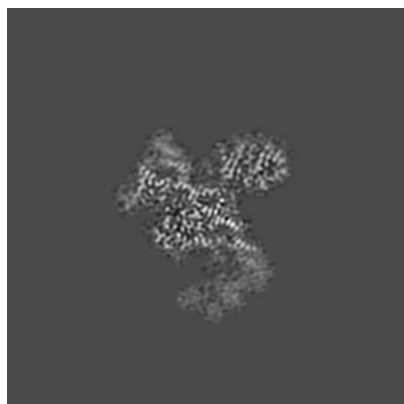


Z Index: 140

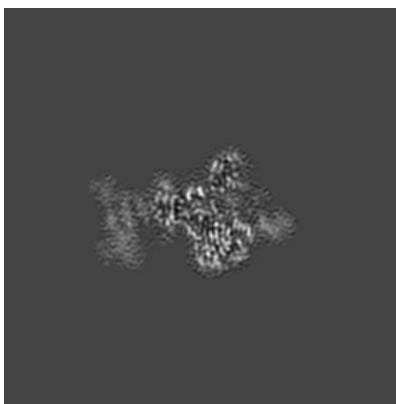
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

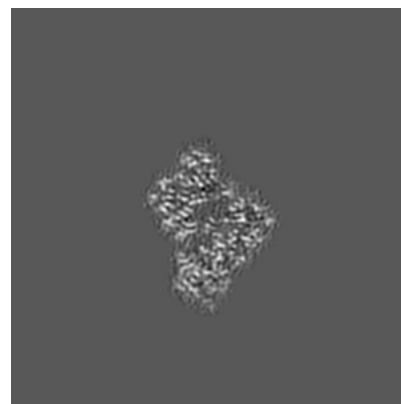
6.3.1 Primary map



X Index: 151



Y Index: 146



Z Index: 144

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

6.4 Orthogonal surface views [i](#)

6.4.1 Primary map



X



Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.032. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

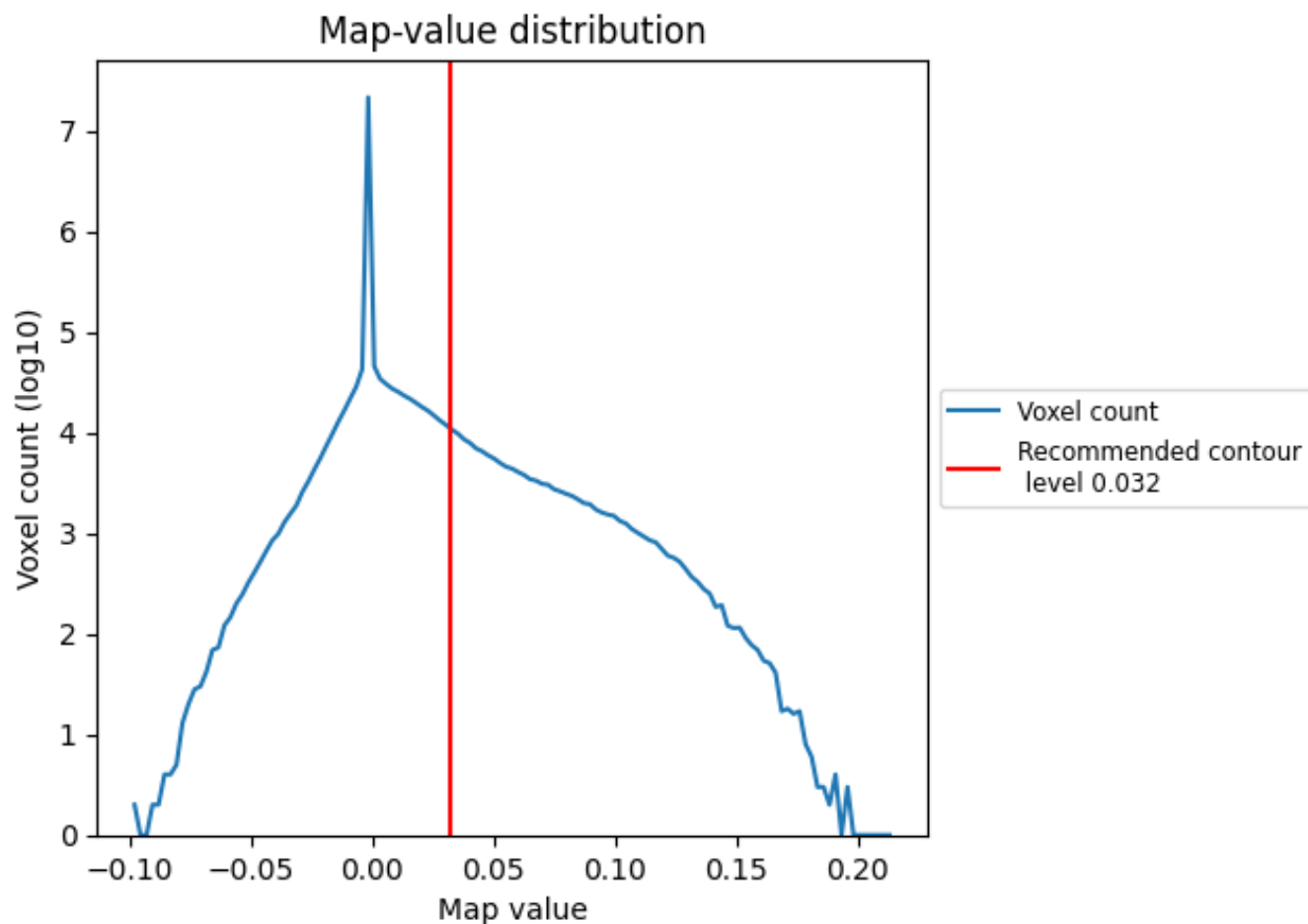
6.5 Mask visualisation

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

7 Map analysis [i](#)

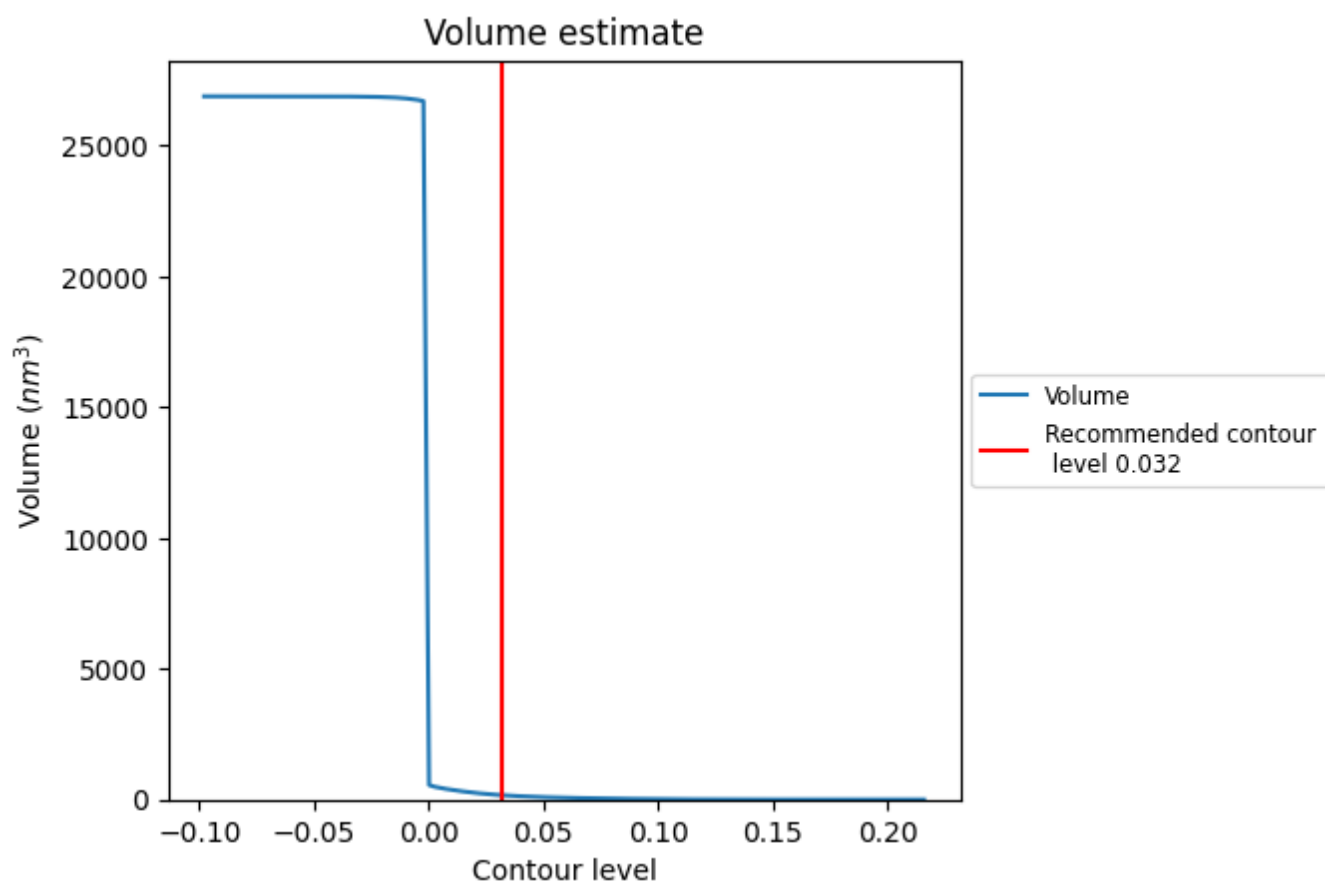
This section contains the results of statistical analysis of the map.

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.

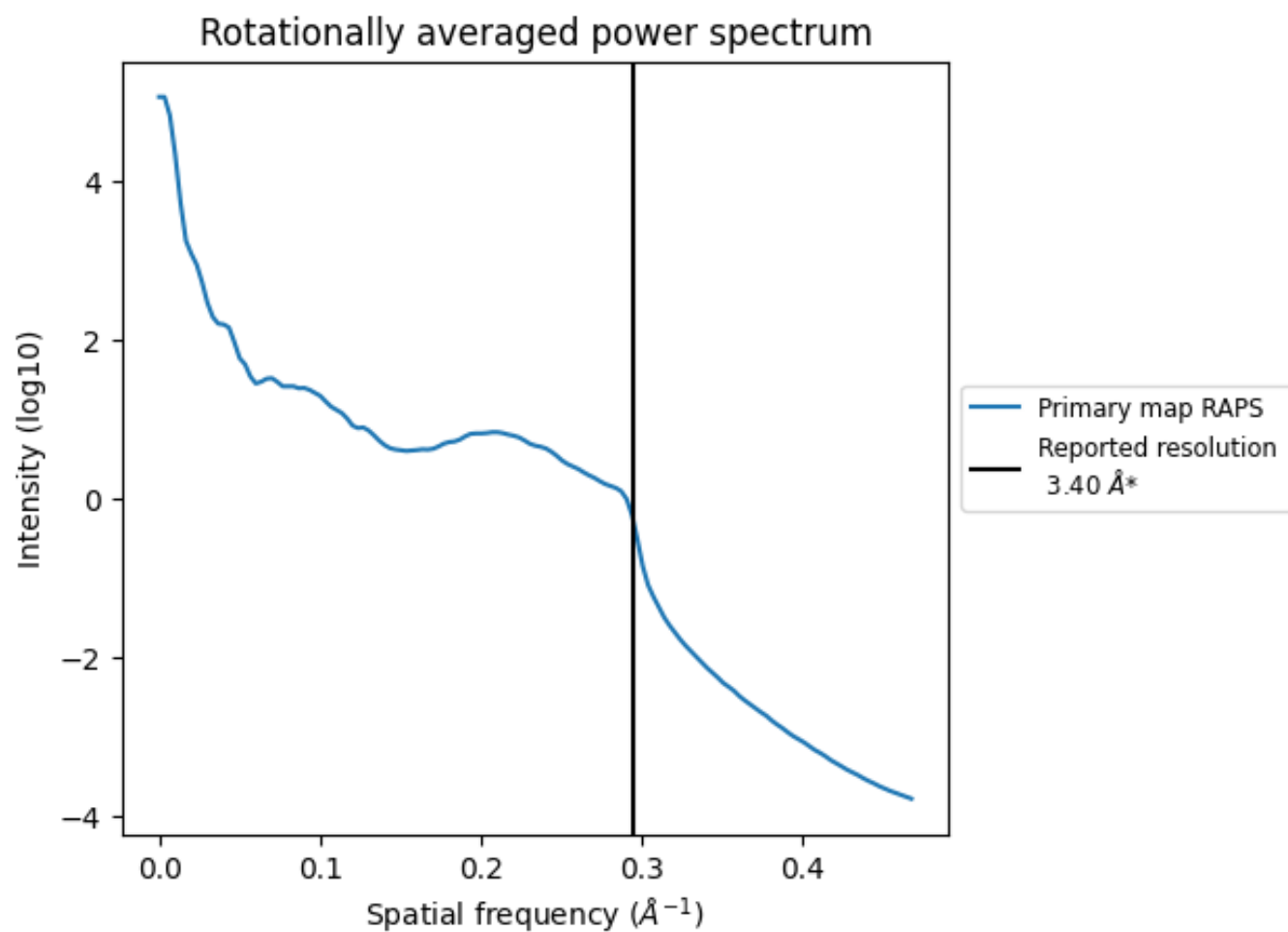
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 167 nm³; this corresponds to an approximate mass of 151 kDa.

The volume estimate graph shows how the enclosed volume varies with the contour level. The recommended contour level is shown as a vertical line and the intersection between the line and the curve gives the volume of the enclosed surface at the given level.

7.3 Rotationally averaged power spectrum ⓘ



*Reported resolution corresponds to spatial frequency of 0.294 Å⁻¹

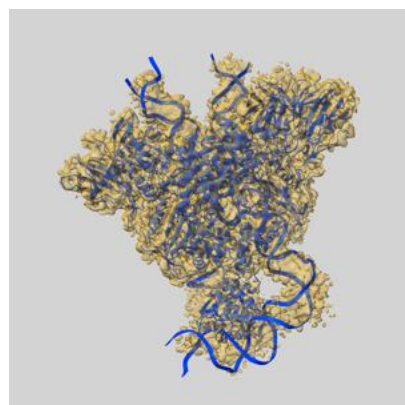
8 Fourier-Shell correlation

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

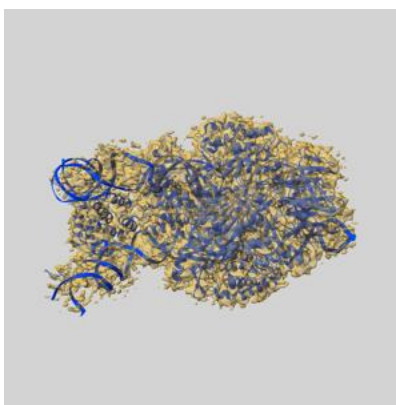
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-20037 and PDB model 6OET. Per-residue inclusion information can be found in [section 3](#) on [page 6](#).

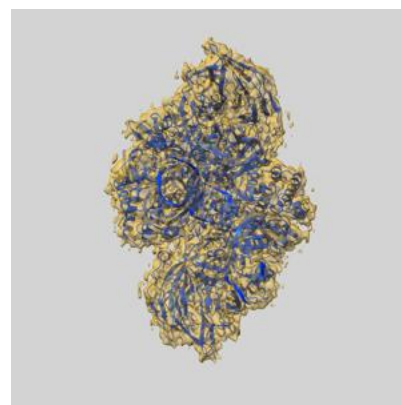
9.1 Map-model overlay [i](#)



X



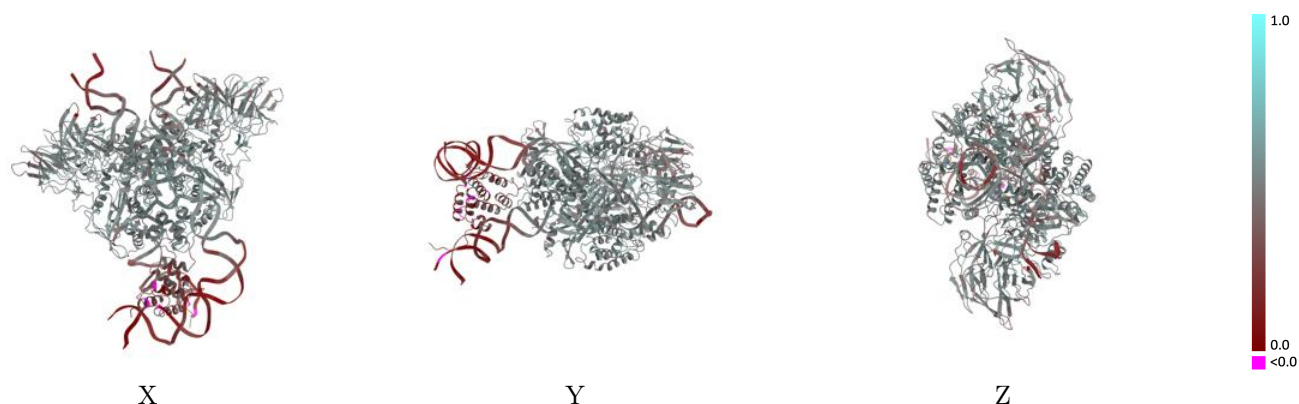
Y



Z

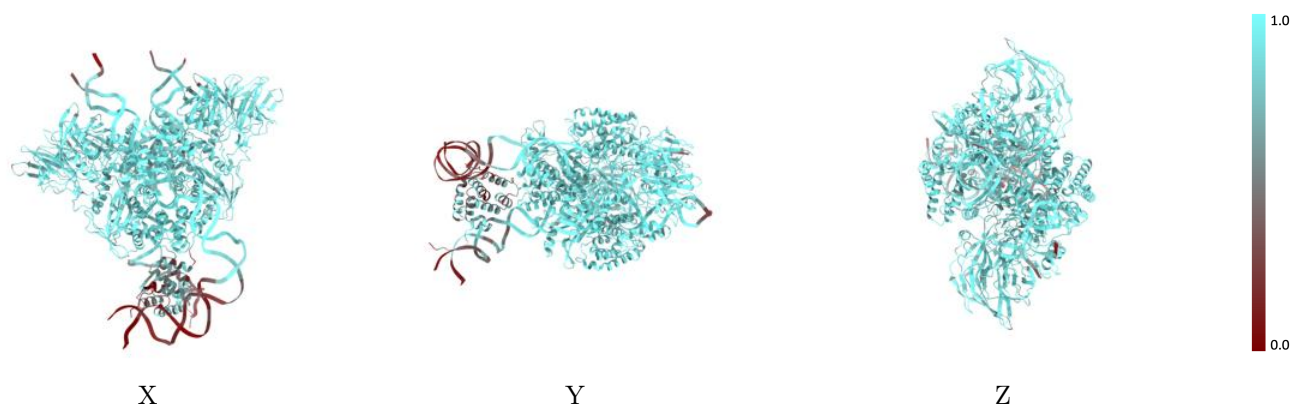
The images above show the 3D surface view of the map at the recommended contour level 0.032 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



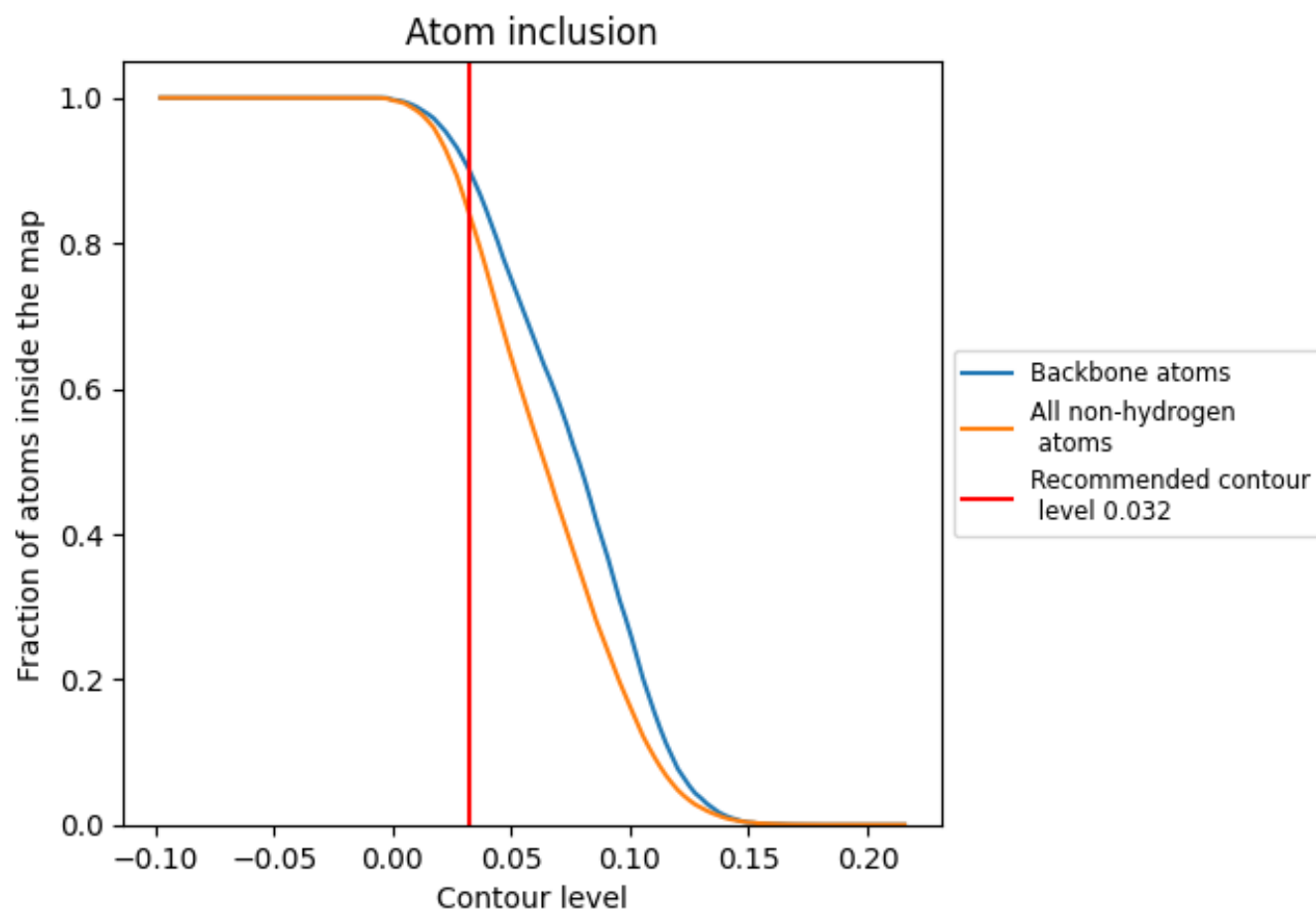
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions at which atoms can be resolved. Low Q-score values may therefore be expected for many entries.

9.3 Atom inclusion mapped to coordinate model [i](#)



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.032).

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 (0.032) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div></div> 0.8437	<div></div> 0.4590
A	<div></div> 0.8743	<div></div> 0.4920
B	<div></div> 0.8915	<div></div> 0.4950
C	<div></div> 0.8731	<div></div> 0.4890
D	<div></div> 0.8988	<div></div> 0.4940
F	<div></div> 0.7893	<div></div> 0.3670
G	<div></div> 0.6598	<div></div> 0.3370
I	<div></div> 0.8680	<div></div> 0.3810
J	<div></div> 0.8218	<div></div> 0.3580
L	<div></div> 0.7349	<div></div> 0.3390
M	<div></div> 0.5727	<div></div> 0.2940

1.0

0.0

<0.0