Full wwPDB EM Validation Report ⓘ

Sep 7, 2023 – 05:19 pm BST

PDB ID	:	8QH5
EMDB ID	:	EMD-18398
Title	:	CryoEM structure of UVSSA(VHS)-CSA-DDB1-DDA1
Deposited on	:	2023-09-06
Resolution	:	3.40 Å(reported)
Based on initial models	:	7OO3, 6ZX9, ., 6UD7

This wwPDB validation report is for manuscript review

This is a Full wwPDB EM Validation Report.

This report is produced by the wwPDB biocuration pipeline after annotation of the structure.

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:

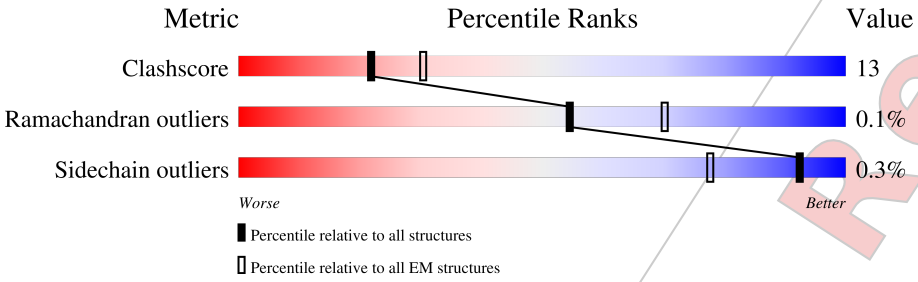
EMDB validation analysis	:	0.0.1.dev50
MolProbity	:	4.02b-467
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ	:	1.9.9
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

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 3.40 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 <=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	408	59% 28% 12%
2	B	1160	9% 68% 28%
3	C	152	5% 21% 11% 68%
4	D	729	16% 18% 79%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 13214 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 excision repair protein ERCC-8.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	358	2787	1737	496	535	19	0	0

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	397	GLY	-	expression tag	UNP Q13216
A	398	THR	-	expression tag	UNP Q13216
A	399	SER	-	expression tag	UNP Q13216
A	400	ALA	-	expression tag	UNP Q13216
A	401	TRP	-	expression tag	UNP Q13216
A	402	SER	-	expression tag	UNP Q13216
A	403	HIS	-	expression tag	UNP Q13216
A	404	PRO	-	expression tag	UNP Q13216
A	405	GLN	-	expression tag	UNP Q13216
A	406	PHE	-	expression tag	UNP Q13216
A	407	GLU	-	expression tag	UNP Q13216
A	408	LYS	-	expression tag	UNP Q13216

- Molecule 2 is a protein called DNA damage-binding protein 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
2	B	1120	8779	5562	1479	1689	49	0	0

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-19	MET	-	initiating methionine	UNP Q16531
B	-18	ALA	-	expression tag	UNP Q16531
B	-17	HIS	-	expression tag	UNP Q16531
B	-16	HIS	-	expression tag	UNP Q16531
B	-15	HIS	-	expression tag	UNP Q16531

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-14	HIS	-	expression tag	UNP Q16531
B	-13	HIS	-	expression tag	UNP Q16531
B	-12	HIS	-	expression tag	UNP Q16531
B	-11	SER	-	expression tag	UNP Q16531
B	-10	ALA	-	expression tag	UNP Q16531
B	-9	ALA	-	expression tag	UNP Q16531
B	-8	LEU	-	expression tag	UNP Q16531
B	-7	GLU	-	expression tag	UNP Q16531
B	-6	VAL	-	expression tag	UNP Q16531
B	-5	LEU	-	expression tag	UNP Q16531
B	-4	PHE	-	expression tag	UNP Q16531
B	-3	GLN	-	expression tag	UNP Q16531
B	-2	GLY	-	expression tag	UNP Q16531
B	-1	PRO	-	expression tag	UNP Q16531
B	0	GLY	-	expression tag	UNP Q16531

- Molecule 3 is a protein called DET1- and DDB1-associated protein 1.

Mol	Chain	Residues	Atoms				AltConf	Trace
			Total	C	N	O		
3	C	48	418	276	70	72	1	0

There are 50 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	103	ASP	-	expression tag	UNP Q9BW61
C	104	VAL	-	expression tag	UNP Q9BW61
C	105	LEU	-	expression tag	UNP Q9BW61
C	106	PHE	-	expression tag	UNP Q9BW61
C	107	GLN	-	expression tag	UNP Q9BW61
C	108	GLY	-	expression tag	UNP Q9BW61
C	109	PRO	-	expression tag	UNP Q9BW61
C	110	GLY	-	expression tag	UNP Q9BW61
C	111	ALA	-	expression tag	UNP Q9BW61
C	112	TRP	-	expression tag	UNP Q9BW61
C	113	SER	-	expression tag	UNP Q9BW61
C	114	HIS	-	expression tag	UNP Q9BW61
C	115	PRO	-	expression tag	UNP Q9BW61
C	116	GLN	-	expression tag	UNP Q9BW61
C	117	PHE	-	expression tag	UNP Q9BW61
C	118	GLU	-	expression tag	UNP Q9BW61
C	119	LYS	-	expression tag	UNP Q9BW61

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Chain	Residue	Modelled	Actual	Comment	Reference
C	120	GLY	-	expression tag	UNP Q9BW61
C	121	GLY	-	expression tag	UNP Q9BW61
C	122	GLY	-	expression tag	UNP Q9BW61
C	123	SER	-	expression tag	UNP Q9BW61
C	124	GLY	-	expression tag	UNP Q9BW61
C	125	GLY	-	expression tag	UNP Q9BW61
C	126	GLY	-	expression tag	UNP Q9BW61
C	127	SER	-	expression tag	UNP Q9BW61
C	128	GLY	-	expression tag	UNP Q9BW61
C	129	GLY	-	expression tag	UNP Q9BW61
C	130	GLY	-	expression tag	UNP Q9BW61
C	131	SER	-	expression tag	UNP Q9BW61
C	132	TRP	-	expression tag	UNP Q9BW61
C	133	SER	-	expression tag	UNP Q9BW61
C	134	HIS	-	expression tag	UNP Q9BW61
C	135	PRO	-	expression tag	UNP Q9BW61
C	136	GLN	-	expression tag	UNP Q9BW61
C	137	PHE	-	expression tag	UNP Q9BW61
C	138	GLU	-	expression tag	UNP Q9BW61
C	139	LYS	-	expression tag	UNP Q9BW61
C	140	GLY	-	expression tag	UNP Q9BW61
C	141	ALA	-	expression tag	UNP Q9BW61
C	142	SER	-	expression tag	UNP Q9BW61
C	143	GLY	-	expression tag	UNP Q9BW61
C	144	GLU	-	expression tag	UNP Q9BW61
C	145	ASP	-	expression tag	UNP Q9BW61
C	146	TYR	-	expression tag	UNP Q9BW61
C	147	LYS	-	expression tag	UNP Q9BW61
C	148	ASP	-	expression tag	UNP Q9BW61
C	149	ASP	-	expression tag	UNP Q9BW61
C	150	ASP	-	expression tag	UNP Q9BW61
C	151	ASP	-	expression tag	UNP Q9BW61
C	152	LYS	-	expression tag	UNP Q9BW61

- Molecule 4 is a protein called UV-stimulated scaffold protein A.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
4	D	150	1230	779	220	227	4	0	0

There are 20 discrepancies between the modelled and reference sequences:

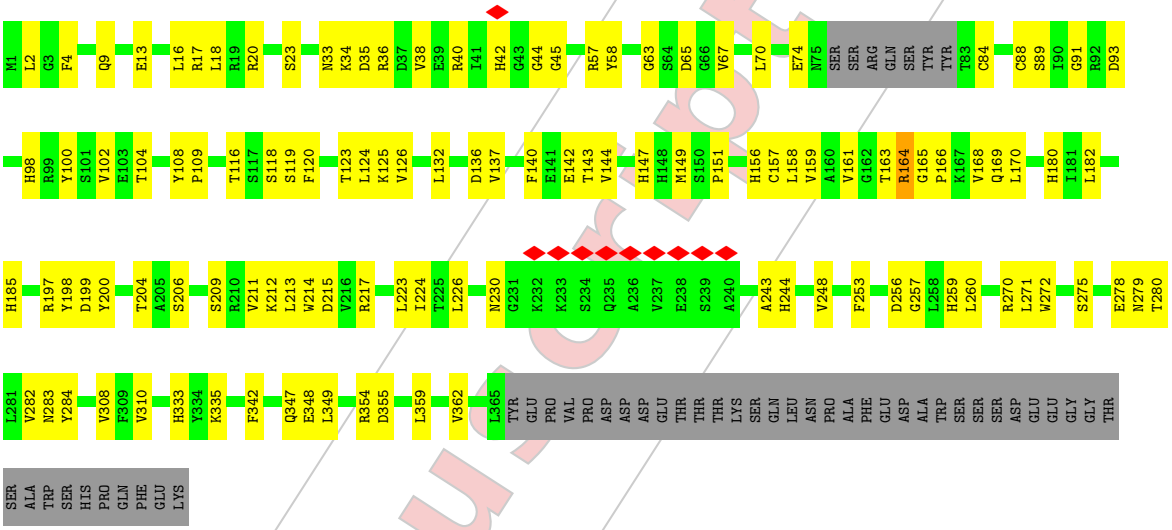
Chain	Residue	Modelled	Actual	Comment	Reference
D	-19	MET	-	initiating methionine	UNP Q2YD98
D	-18	ALA	-	expression tag	UNP Q2YD98
D	-17	HIS	-	expression tag	UNP Q2YD98
D	-16	HIS	-	expression tag	UNP Q2YD98
D	-15	HIS	-	expression tag	UNP Q2YD98
D	-14	HIS	-	expression tag	UNP Q2YD98
D	-13	HIS	-	expression tag	UNP Q2YD98
D	-12	HIS	-	expression tag	UNP Q2YD98
D	-11	SER	-	expression tag	UNP Q2YD98
D	-10	ALA	-	expression tag	UNP Q2YD98
D	-9	ALA	-	expression tag	UNP Q2YD98
D	-8	LEU	-	expression tag	UNP Q2YD98
D	-7	GLU	-	expression tag	UNP Q2YD98
D	-6	VAL	-	expression tag	UNP Q2YD98
D	-5	LEU	-	expression tag	UNP Q2YD98
D	-4	PHE	-	expression tag	UNP Q2YD98
D	-3	GLN	-	expression tag	UNP Q2YD98
D	-2	GLY	-	expression tag	UNP Q2YD98
D	-1	PRO	-	expression tag	UNP Q2YD98
D	0	GLY	-	expression tag	UNP Q2YD98

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.

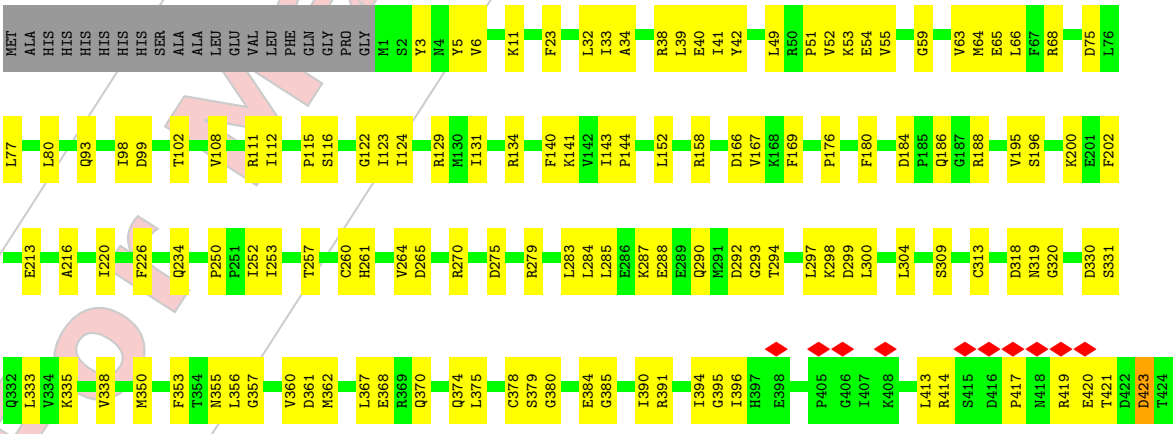
• Molecule 1: DNA excision repair protein ERCC-8

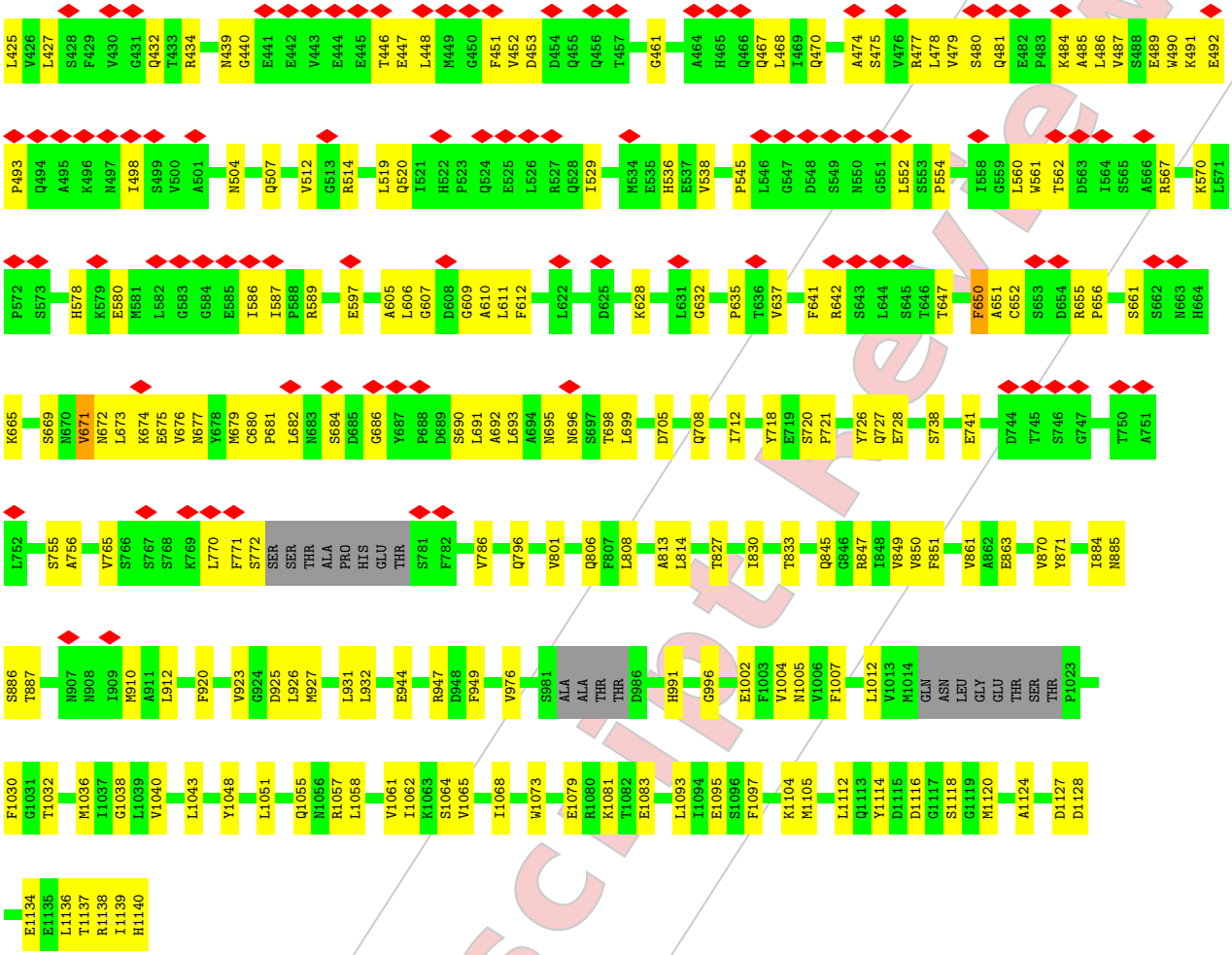
Chain A:



• Molecule 2: DNA damage-binding protein 1

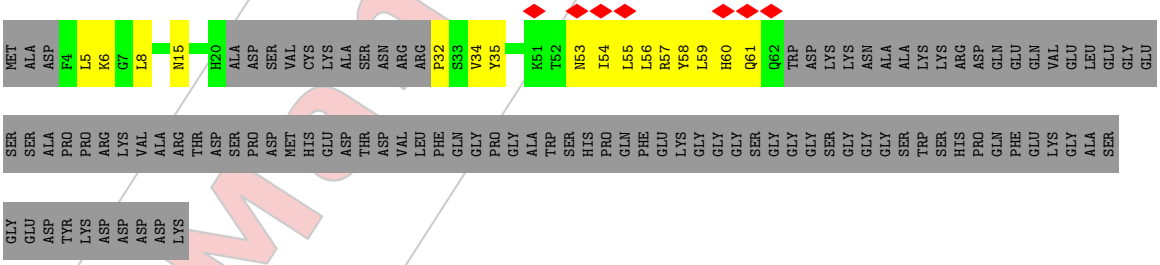
Chain B:





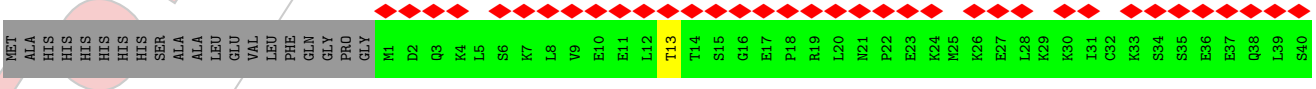
• Molecule 3: DET1- and DDB1-associated protein 1

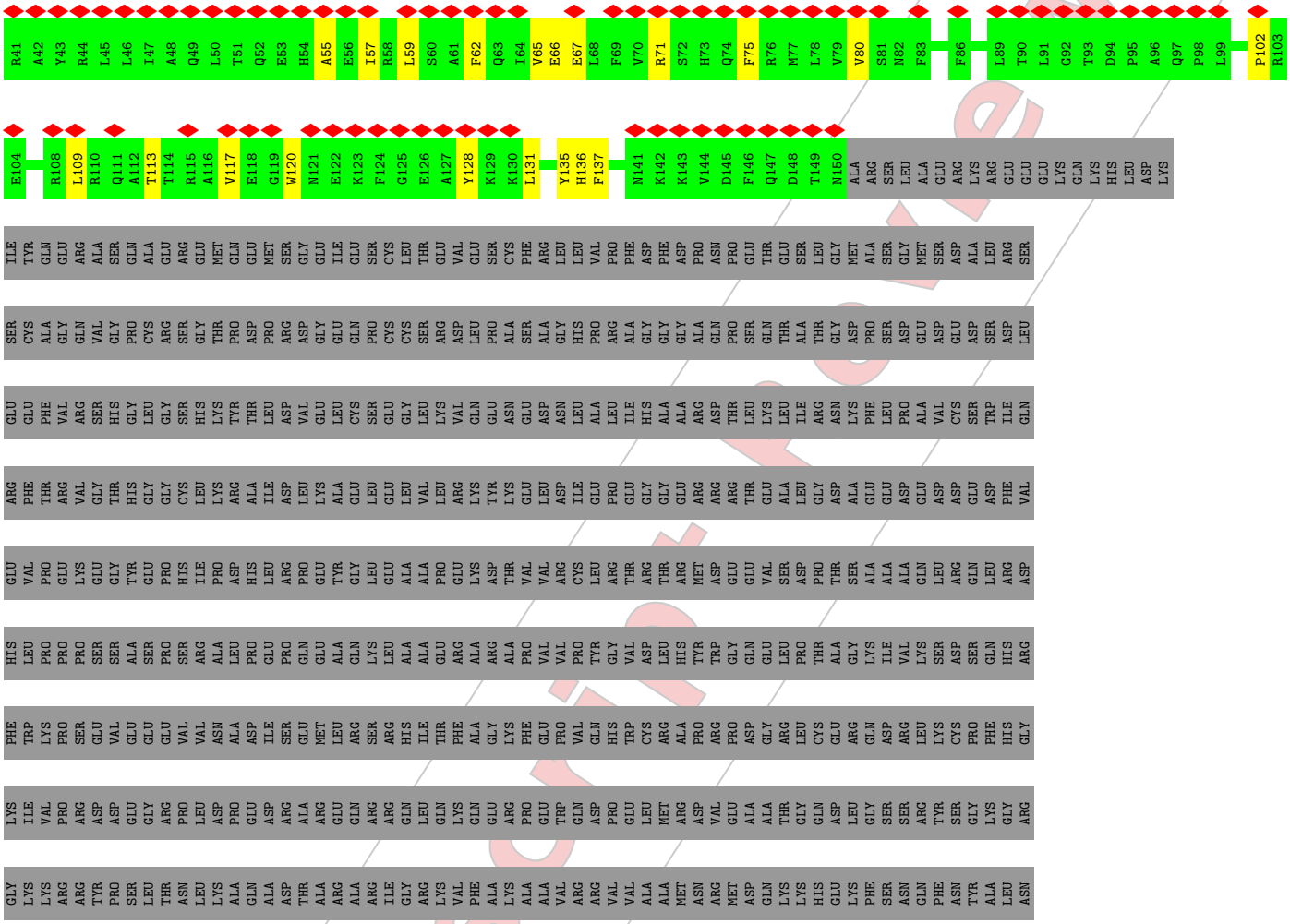
Chain C:



• Molecule 4: UV-stimulated scaffold protein A

Chain D:





4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	294142	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 ($e^-/Å^2$)	60	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	3000	Depositor
Magnification	81000	Depositor
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	85.662	Depositor
Minimum map value	-1.899	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.948	Depositor
Recommended contour level	0.828	Depositor
Map size (Å)	338.24, 338.24, 338.24	wwPDB
Map dimensions	320, 320, 320	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.0569999, 1.0569999, 1.0569999	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	0.27	0/2843	0.56	0/3850
2	B	0.29	0/8939	0.56	0/12101
3	C	0.29	0/433	0.50	0/585
4	D	0.24	0/1253	0.45	0/1685
All	All	0.28	0/13468	0.55	0/18221

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2787	0	2723	85	0
2	B	8779	0	8753	246	0
3	C	418	0	421	19	0
4	D	1230	0	1248	13	0
All	All	13214	0	13145	341	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 13.

All (341) 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:335:LYS:NZ	4:D:137:PHE:HE1	1.50	1.07
1:A:335:LYS:NZ	4:D:137:PHE:CE1	2.31	0.93
2:B:675:GLU:HB3	2:B:695:ASN:HB2	1.59	0.83
2:B:467:GLN:HG3	2:B:478:LEU:HD11	1.61	0.81
1:A:215:ASP:OD1	1:A:217:ARG:NH2	2.16	0.78
2:B:394:ILE:HG23	2:B:705:ASP:HB2	1.65	0.77
3:C:53:ASN:HB2	3:C:56:LEU:HD13	1.68	0.76
2:B:368:GLU:HG3	2:B:674:LYS:HE3	1.68	0.74
2:B:671:VAL:HG12	2:B:672:ASN:H	1.53	0.74
2:B:676:VAL:HG22	2:B:693:LEU:HD23	1.70	0.73
2:B:656:PRO:CD	2:B:674:LYS:HA	2.18	0.73
2:B:606:LEU:HD11	2:B:612:PHE:HE1	1.53	0.73
1:A:118:SER:HB2	1:A:144:VAL:HG11	1.72	0.71
3:C:57:ARG:O	3:C:60:HIS:ND1	2.23	0.71
2:B:55:VAL:HG23	3:C:35:TYR:HD2	1.55	0.70
1:A:123:THR:OG1	1:A:125:LYS:NZ	2.24	0.69
2:B:656:PRO:HG3	2:B:676:VAL:HB	1.74	0.69
2:B:771:PHE:CE1	2:B:845:GLN:HB3	2.27	0.69
2:B:675:GLU:HB3	2:B:695:ASN:CB	2.23	0.68
2:B:391:ARG:NH2	2:B:672:ASN:OD1	2.25	0.68
2:B:367:LEU:HD23	2:B:368:GLU:HG2	1.75	0.68
3:C:60:HIS:CE1	3:C:61:GLN:HG3	2.29	0.68
1:A:144:VAL:HA	1:A:163:THR:HG22	1.76	0.68
1:A:259:HIS:HB3	1:A:271:LEU:HD11	1.76	0.68
2:B:391:ARG:NH2	2:B:672:ASN:CG	2.46	0.68
2:B:53:LYS:HE2	3:C:35:TYR:HB2	1.77	0.67
2:B:102:THR:OG1	2:B:1065:VAL:O	2.12	0.67
2:B:419:ARG:HG3	2:B:420:GLU:N	2.09	0.67
2:B:367:LEU:HG	2:B:796:GLN:HE21	1.59	0.67
2:B:765:VAL:HG22	2:B:806:GLN:HB3	1.77	0.67
2:B:338:VAL:HA	3:C:6:LYS:HG3	1.75	0.66
1:A:283:ASN:OD1	1:A:284:TYR:N	2.27	0.66
1:A:197:ARG:NH2	2:B:112:ILE:O	2.28	0.66
1:A:109:PRO:HG2	1:A:151:PRO:HA	1.77	0.66
2:B:1058:LEU:HA	2:B:1061:VAL:HG12	1.79	0.65
2:B:656:PRO:HD2	2:B:674:LYS:HA	1.78	0.65
2:B:677:ASN:ND2	2:B:696:ASN:HB2	2.12	0.65
2:B:284:LEU:HD11	3:C:5:LEU:HG	1.79	0.64
2:B:391:ARG:HH22	2:B:672:ASN:HA	1.62	0.64
1:A:244:HIS:HD2	1:A:270:ARG:HE	1.46	0.64
2:B:650:PHE:CE1	2:B:676:VAL:HG11	2.34	0.63
2:B:419:ARG:HG3	2:B:420:GLU:H	1.64	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:54:GLU:HB2	3:C:32:PRO:HB2	1.81	0.61
2:B:250:PRO:HG2	2:B:253:ILE:HG12	1.82	0.61
2:B:578:HIS:NE2	2:B:580:GLU:HG2	2.15	0.61
2:B:391:ARG:HH22	2:B:672:ASN:CG	2.04	0.61
2:B:375:LEU:HG	2:B:1012:LEU:HD21	1.83	0.61
2:B:413:LEU:HD11	2:B:468:LEU:HD22	1.81	0.61
2:B:830:ILE:HG12	2:B:850:VAL:HG22	1.83	0.61
2:B:498:ILE:HA	2:B:512:VAL:HG12	1.82	0.61
1:A:67:VAL:HG23	1:A:91:GLY:HA2	1.83	0.61
2:B:252:ILE:HD11	2:B:304:LEU:HB2	1.82	0.60
1:A:58:TYR:HB3	1:A:70:LEU:HD11	1.83	0.60
2:B:39:LEU:HB3	2:B:55:VAL:CG1	2.31	0.60
1:A:230:ASN:HD21	1:A:279:ASN:H	1.50	0.60
2:B:196:SER:O	2:B:200:LYS:N	2.35	0.60
2:B:432:GLN:HB2	2:B:453:ASP:O	2.02	0.60
2:B:665:LYS:HD3	2:B:1138:ARG:HH12	1.67	0.59
1:A:36:ARG:NH1	1:A:348:GLU:OE1	2.35	0.59
1:A:158:LEU:HD13	1:A:170:LEU:HD23	1.84	0.59
1:A:74:GLU:OE2	1:A:74:GLU:N	2.31	0.59
2:B:925:ASP:OD1	2:B:927:MET:N	2.31	0.59
2:B:108:VAL:O	2:B:141:LYS:NZ	2.28	0.59
1:A:199:ASP:OD2	2:B:111:ARG:NH1	2.35	0.59
2:B:265:ASP:OD1	2:B:270:ARG:NH1	2.35	0.58
2:B:283:LEU:HD11	2:B:300:LEU:HD12	1.85	0.58
2:B:673:LEU:HG	2:B:675:GLU:H	1.68	0.58
4:D:13:THR:HB	4:D:57:ILE:HG23	1.85	0.58
2:B:275:ASP:OD1	2:B:279:ARG:N	2.37	0.58
2:B:318:ASP:OD1	2:B:319:ASN:N	2.37	0.58
2:B:1064:SER:OG	2:B:1068:ILE:N	2.35	0.58
2:B:53:LYS:HD3	2:B:98:ILE:HD11	1.84	0.58
2:B:1002:GLU:HG3	2:B:1036:MET:SD	2.44	0.58
2:B:641:PHE:HB3	2:B:681:PRO:HG3	1.86	0.58
1:A:198:TYR:OH	3:C:54:ILE:HG12	2.04	0.57
1:A:215:ASP:HB2	1:A:223:LEU:HD21	1.87	0.57
2:B:213:GLU:OE1	2:B:213:GLU:N	2.31	0.57
1:A:88:CYS:SG	1:A:89:SER:N	2.77	0.57
2:B:370:GLN:N	2:B:370:GLN:OE1	2.37	0.57
2:B:489:GLU:OE2	2:B:491:LYS:HE3	2.04	0.57
2:B:65:GLU:N	2:B:65:GLU:OE1	2.37	0.57
2:B:370:GLN:NE2	2:B:674:LYS:NZ	2.52	0.57
3:C:54:ILE:HG13	3:C:55:LEU:N	2.20	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:156:HIS:CD2	1:A:200:TYR:HE1	2.23	0.57
1:A:142:GLU:HB3	1:A:164:ARG:HE	1.69	0.56
2:B:370:GLN:HG2	2:B:655:ARG:NH1	2.20	0.56
2:B:432:GLN:HG2	2:B:434:ARG:NH1	2.20	0.56
2:B:756:ALA:HB1	2:B:801:VAL:HG21	1.86	0.56
1:A:280:THR:HG23	1:A:282:VAL:HG23	1.87	0.56
2:B:589:ARG:NH1	2:B:637:VAL:HG22	2.20	0.56
1:A:136:ASP:OD1	1:A:137:VAL:N	2.37	0.56
2:B:166:ASP:OD1	2:B:167:VAL:N	2.39	0.56
1:A:33:ASN:OD1	1:A:34:LYS:N	2.39	0.56
1:A:169:GLN:OE1	1:A:169:GLN:N	2.39	0.56
2:B:425:LEU:HD21	2:B:427:LEU:HD21	1.88	0.56
2:B:771:PHE:HE1	2:B:845:GLN:HB3	1.71	0.56
2:B:949:PHE:HE2	2:B:991:HIS:CD2	2.24	0.56
2:B:32:LEU:HD23	2:B:39:LEU:HD21	1.88	0.56
2:B:485:ALA:O	2:B:487:VAL:HG23	2.06	0.56
3:C:54:ILE:HG13	3:C:55:LEU:H	1.71	0.56
2:B:1104:LYS:HE3	2:B:1104:LYS:HA	1.88	0.56
2:B:771:PHE:CE2	2:B:847:ARG:HD3	2.42	0.55
4:D:67:GLU:OE2	4:D:71:ARG:NH1	2.39	0.55
2:B:23:PHE:HB2	2:B:66:LEU:HD21	1.88	0.55
2:B:642:ARG:NH1	2:B:647:THR:HB	2.22	0.55
4:D:62:PHE:CZ	4:D:113:THR:HA	2.42	0.55
2:B:446:THR:HG22	2:B:447:GLU:H	1.72	0.55
2:B:1136:LEU:O	2:B:1139:ILE:HG12	2.06	0.54
2:B:330:ASP:HA	2:B:355:ASN:HB3	1.89	0.54
1:A:244:HIS:ND1	1:A:248:VAL:HG12	2.23	0.54
2:B:361:ASP:OD1	2:B:362:MET:N	2.40	0.54
1:A:98:HIS:ND1	1:A:119:SER:OG	2.36	0.54
2:B:370:GLN:HG2	2:B:655:ARG:HH11	1.73	0.54
1:A:116:THR:HG22	1:A:126:VAL:HG23	1.90	0.54
2:B:920:PHE:HB3	2:B:932:LEU:HD11	1.90	0.53
2:B:234:GLN:HE22	2:B:257:THR:HG22	1.72	0.53
2:B:561:TRP:HB3	2:B:562:THR:HG23	1.90	0.53
2:B:448:LEU:HA	2:B:484:LYS:NZ	2.23	0.53
2:B:378:CYS:HB3	2:B:721:PRO:HB2	1.91	0.53
4:D:117:VAL:HG12	4:D:135:TYR:HD1	1.74	0.53
1:A:224:ILE:HG12	3:C:59:LEU:HD21	1.91	0.53
2:B:771:PHE:O	2:B:772:SER:C	2.47	0.53
2:B:167:VAL:HG23	2:B:180:PHE:HB3	1.91	0.53
1:A:342:PHE:HD1	1:A:349:LEU:HD13	1.74	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:695:ASN:OD1	2:B:698:THR:N	2.37	0.53
2:B:1030:PHE:CZ	2:B:1038:GLY:HA3	2.44	0.52
1:A:224:ILE:HG22	1:A:226:LEU:HD12	1.91	0.52
1:A:230:ASN:ND2	1:A:278:GLU:HA	2.25	0.52
1:A:333:HIS:CE1	1:A:359:LEU:HD13	2.45	0.52
4:D:59:LEU:HA	4:D:109:LEU:HD13	1.92	0.52
2:B:34:ALA:HB2	2:B:64:MET:HE1	1.91	0.52
2:B:690:SER:OG	2:B:691:LEU:N	2.42	0.52
2:B:589:ARG:HH12	2:B:637:VAL:HG22	1.73	0.52
1:A:42:HIS:CD2	1:A:63:GLY:HA3	2.45	0.51
2:B:285:LEU:HD22	2:B:297:LEU:HD11	1.91	0.51
2:B:1114:TYR:HE2	2:B:1116:ASP:HB3	1.75	0.51
2:B:507:GLN:NE2	2:B:552:LEU:HG	2.26	0.51
1:A:13:GLU:OE1	1:A:17:ARG:NH1	2.43	0.51
1:A:226:LEU:HD23	1:A:272:TRP:CG	2.46	0.51
1:A:185:HIS:ND1	1:A:206:SER:OG	2.39	0.51
2:B:49:LEU:HD13	2:B:333:LEU:HD11	1.93	0.51
2:B:671:VAL:HG12	2:B:672:ASN:N	2.24	0.51
2:B:944:GLU:HG2	2:B:947:ARG:HH12	1.76	0.51
1:A:244:HIS:CD2	1:A:270:ARG:HE	2.26	0.51
2:B:5:TYR:HB2	2:B:1043:LEU:HD11	1.93	0.51
2:B:195:VAL:HG12	2:B:202:PHE:CE1	2.46	0.51
2:B:63:VAL:HG21	2:B:122:GLY:HA3	1.93	0.50
2:B:453:ASP:N	2:B:453:ASP:OD1	2.44	0.50
2:B:53:LYS:CE	3:C:35:TYR:HB2	2.41	0.50
2:B:490:TRP:CG	2:B:519:LEU:HD21	2.47	0.50
1:A:45:GLY:HA3	1:A:354:ARG:HA	1.93	0.50
1:A:120:PHE:HA	1:A:143:THR:OG1	2.11	0.50
2:B:129:ARG:HE	2:B:176:PRO:HG3	1.76	0.50
1:A:149:MET:HG3	1:A:159:VAL:HG12	1.92	0.50
2:B:390:ILE:HG12	2:B:712:ILE:HG12	1.92	0.50
1:A:4:PHE:CD1	1:A:18:LEU:HB3	2.47	0.50
2:B:396:ILE:HD12	2:B:396:ILE:O	2.11	0.50
1:A:335:LYS:HG2	1:A:355:ASP:OD1	2.11	0.50
2:B:504:ASN:HD22	2:B:545:PRO:HD3	1.76	0.50
1:A:98:HIS:CE1	1:A:125:LYS:HE2	2.47	0.49
2:B:188:ARG:NH1	2:B:216:ALA:O	2.44	0.49
2:B:467:GLN:HE21	2:B:478:LEU:HD21	1.77	0.49
2:B:650:PHE:CZ	2:B:676:VAL:HG11	2.47	0.49
2:B:11:LYS:HE2	2:B:38:ARG:HD2	1.93	0.49
2:B:467:GLN:NE2	2:B:478:LEU:HD21	2.27	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:607:GLY:HA2	2:B:635:PRO:HB3	1.95	0.49
2:B:380:GLY:HA3	2:B:385:GLY:HA2	1.94	0.49
2:B:538:VAL:HA	2:B:560:LEU:HD23	1.94	0.49
2:B:597:GLU:OE2	2:B:661:SER:OG	2.28	0.49
2:B:813:ALA:HA	2:B:833:THR:HG22	1.94	0.49
2:B:1127:ASP:OD1	2:B:1128:ASP:N	2.46	0.49
1:A:4:PHE:HD1	1:A:18:LEU:HB3	1.78	0.49
2:B:41:ILE:O	2:B:52:VAL:HG22	2.12	0.48
2:B:374:GLN:HG3	2:B:391:ARG:HG2	1.94	0.48
2:B:421:THR:OG1	2:B:423:ASP:OD1	2.28	0.48
2:B:490:TRP:CD2	2:B:519:LEU:HD21	2.49	0.48
2:B:871:TYR:HE2	2:B:885:ASN:HA	1.78	0.48
2:B:116:SER:OG	2:B:134:ARG:NH2	2.47	0.48
2:B:319:ASN:ND2	3:C:15:ASN:O	2.47	0.48
2:B:677:ASN:HD22	2:B:696:ASN:HB2	1.77	0.48
2:B:726:TYR:OH	2:B:728:GLU:OE1	2.32	0.48
2:B:1055:GLN:HG3	2:B:1093:LEU:HD23	1.96	0.48
2:B:63:VAL:HG12	2:B:80:LEU:HB3	1.95	0.48
1:A:200:TYR:HD2	1:A:217:ARG:NH1	2.12	0.48
1:A:230:ASN:HD21	1:A:278:GLU:HA	1.79	0.48
2:B:1114:TYR:HB2	2:B:1124:ALA:HB2	1.96	0.48
2:B:1118:SER:OG	2:B:1120:MET:SD	2.60	0.48
1:A:33:ASN:HB2	1:A:362:VAL:HG12	1.96	0.48
2:B:665:LYS:CD	2:B:1138:ARG:HH12	2.27	0.48
2:B:578:HIS:CD2	2:B:580:GLU:HG2	2.49	0.47
1:A:40:ARG:HH12	4:D:136:HIS:CG	2.31	0.47
2:B:606:LEU:HD11	2:B:612:PHE:CE1	2.42	0.47
2:B:480:SER:O	2:B:484:LYS:HA	2.14	0.47
1:A:209:SER:O	1:A:243:ALA:N	2.47	0.47
1:A:16:LEU:HD11	2:B:1079:GLU:OE2	2.14	0.47
1:A:102:VAL:HG12	1:A:119:SER:HB3	1.97	0.47
2:B:356:LEU:HD21	2:B:390:ILE:HD11	1.95	0.47
1:A:42:HIS:HD2	1:A:63:GLY:HA3	1.80	0.47
2:B:370:GLN:CB	2:B:655:ARG:NH1	2.78	0.47
2:B:910:MET:HE2	2:B:910:MET:HB3	1.59	0.46
2:B:1114:TYR:CE2	2:B:1116:ASP:HB3	2.50	0.46
2:B:54:GLU:HB3	3:C:34:VAL:HA	1.98	0.46
2:B:195:VAL:HG12	2:B:202:PHE:HE1	1.79	0.46
1:A:108:TYR:OH	1:A:157:CYS:SG	2.60	0.46
1:A:256:ASP:OD1	1:A:257:GLY:N	2.49	0.46
1:A:198:TYR:CZ	3:C:54:ILE:HG12	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:6:VAL:HG22	2:B:1040:VAL:HG22	1.98	0.46
2:B:474:ALA:O	2:B:475:SER:HB2	2.16	0.46
2:B:1051:LEU:HD11	2:B:1136:LEU:HD11	1.97	0.46
2:B:808:LEU:HD23	2:B:847:ARG:HH12	1.78	0.46
2:B:42:TYR:CD1	2:B:51:PRO:HA	2.51	0.46
1:A:23:SER:HB3	1:A:347:GLN:HE21	1.81	0.46
2:B:220:ILE:HG12	2:B:261:HIS:CE1	2.51	0.46
2:B:440:GLY:O	2:B:686:GLY:HA3	2.16	0.46
1:A:253:PHE:CE1	1:A:260:LEU:HD13	2.51	0.46
2:B:417:PRO:HG3	2:B:481:GLN:OE1	2.16	0.46
2:B:1116:ASP:N	2:B:1116:ASP:OD1	2.49	0.46
1:A:104:THR:OG1	1:A:147:HIS:ND1	2.41	0.45
2:B:468:LEU:HD21	2:B:481:GLN:NE2	2.31	0.45
2:B:260:CYS:SG	2:B:313:CYS:HA	2.57	0.45
2:B:850:VAL:O	2:B:861:VAL:HG12	2.15	0.45
1:A:118:SER:HB3	1:A:124:LEU:HD12	1.97	0.45
2:B:40:GLU:OE1	2:B:40:GLU:HA	2.16	0.45
2:B:299:ASP:OD1	2:B:300:LEU:N	2.50	0.45
2:B:912:LEU:HD11	2:B:926:LEU:HD12	1.98	0.45
2:B:320:GLY:HA2	3:C:8:LEU:HD13	1.98	0.45
2:B:520:GLN:HG3	2:B:529:ILE:HG13	1.99	0.45
1:A:9:GLN:O	2:B:1005:ASN:ND2	2.46	0.45
1:A:211:VAL:HG12	1:A:243:ALA:HB2	1.98	0.45
2:B:370:GLN:HB3	2:B:655:ARG:NH1	2.32	0.45
2:B:718:TYR:HB2	2:B:755:SER:HA	1.99	0.45
1:A:65:ASP:OD1	1:A:67:VAL:HG12	2.17	0.45
2:B:33:ILE:O	2:B:39:LEU:HD12	2.17	0.45
1:A:185:HIS:CD2	1:A:212:LYS:HD3	2.51	0.45
2:B:560:LEU:HD12	2:B:567:ARG:HH11	1.81	0.45
1:A:44:GLY:N	1:A:65:ASP:OD2	2.50	0.45
2:B:357:GLY:N	2:B:379:SER:OG	2.50	0.45
1:A:182:LEU:HD21	1:A:214:TRP:CE2	2.52	0.45
1:A:213:LEU:HD12	1:A:224:ILE:HB	1.98	0.44
2:B:770:LEU:CD1	2:B:863:GLU:HG3	2.47	0.44
1:A:33:ASN:OD1	1:A:35:ASP:N	2.51	0.44
2:B:560:LEU:CD1	2:B:567:ARG:HH11	2.30	0.44
2:B:586:ILE:HD12	2:B:587:ILE:H	1.82	0.44
1:A:38:VAL:HG22	1:A:84:CYS:HB2	1.98	0.44
4:D:55:ALA:HB3	4:D:102:PRO:HD2	1.99	0.44
2:B:184:ASP:OD2	2:B:186:GLN:NE2	2.51	0.44
2:B:196:SER:O	2:B:200:LYS:CA	2.65	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:170:LEU:HB2	1:A:180:HIS:HB3	1.98	0.44
2:B:3:TYR:HD1	2:B:1048:TYR:CG	2.35	0.44
2:B:414:ARG:NH2	2:B:419:ARG:O	2.49	0.44
2:B:492:GLU:HG3	2:B:512:VAL:HG21	1.99	0.44
2:B:93:GLN:HG3	2:B:98:ILE:HG22	2.00	0.44
2:B:425:LEU:HD12	2:B:682:LEU:HD12	2.00	0.44
2:B:474:ALA:O	2:B:491:LYS:HE2	2.18	0.44
2:B:610:ALA:HB1	2:B:628:LYS:HE3	2.00	0.44
2:B:656:PRO:HD3	2:B:674:LYS:HA	1.97	0.44
3:C:56:LEU:HA	3:C:59:LEU:HD12	2.00	0.44
1:A:100:TYR:HB2	1:A:120:PHE:HB2	2.00	0.43
2:B:692:ALA:C	2:B:693:LEU:HD12	2.38	0.43
4:D:80:VAL:HG22	4:D:131:LEU:HD23	1.99	0.43
2:B:68:ARG:HB2	2:B:75:ASP:OD1	2.18	0.43
2:B:395:GLY:HA2	2:B:672:ASN:HB2	1.99	0.43
2:B:923:VAL:HB	2:B:931:LEU:HB3	2.00	0.43
2:B:1058:LEU:HD12	2:B:1062:ILE:HD12	2.00	0.43
1:A:132:LEU:HD23	1:A:132:LEU:HA	1.87	0.43
1:A:168:VAL:HG21	1:A:204:THR:HG21	1.99	0.43
1:A:275:SER:HB2	2:B:158:ARG:NH1	2.32	0.43
2:B:39:LEU:HB3	2:B:55:VAL:HG12	1.99	0.43
2:B:770:LEU:HD13	2:B:863:GLU:HG3	2.00	0.43
2:B:885:ASN:HB3	2:B:886:SER:H	1.59	0.43
2:B:41:ILE:HD12	2:B:53:LYS:HB2	2.01	0.43
2:B:637:VAL:HB	2:B:652:CYS:HB2	2.00	0.43
2:B:741:GLU:HB2	2:B:786:VAL:HG13	1.98	0.43
1:A:140:PHE:CZ	1:A:161:VAL:HG21	2.54	0.43
1:A:226:LEU:HD23	1:A:272:TRP:CD2	2.52	0.43
2:B:439:ASN:OD1	2:B:439:ASN:N	2.51	0.43
4:D:80:VAL:HG21	4:D:128:TYR:HD2	1.83	0.43
1:A:93:ASP:OD1	1:A:93:ASP:N	2.47	0.43
2:B:477:ARG:NH2	2:B:486:LEU:HD22	2.32	0.43
2:B:554:PRO:O	2:B:570:LYS:HD2	2.19	0.43
2:B:331:SER:OG	2:B:353:PHE:HB2	2.18	0.43
1:A:20:ARG:NH2	2:B:115:PRO:HG2	2.34	0.42
2:B:335:LYS:HB2	2:B:350:MET:SD	2.58	0.42
2:B:609:GLY:HA3	2:B:632:GLY:O	2.20	0.42
2:B:1004:VAL:HA	2:B:1032:THR:HG22	2.01	0.42
4:D:65:VAL:HG13	4:D:75:PHE:HE2	1.84	0.42
2:B:53:LYS:HD3	2:B:98:ILE:CD1	2.49	0.42
2:B:360:VAL:HG21	2:B:721:PRO:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:884:ILE:O	2:B:887:THR:OG1	2.35	0.42
2:B:309:SER:HA	2:B:384:GLU:OE2	2.19	0.42
2:B:394:ILE:HG12	2:B:708:GLN:HA	2.02	0.42
2:B:99:ASP:OD1	2:B:99:ASP:N	2.53	0.42
2:B:290:GLN:OE1	2:B:294:THR:OG1	2.37	0.42
2:B:669:SER:OG	2:B:708:GLN:HB3	2.20	0.42
4:D:66:GLU:HA	4:D:120:TRP:HZ2	1.84	0.42
2:B:124:ILE:CD1	2:B:131:ILE:HG12	2.50	0.42
2:B:123:ILE:HD12	2:B:169:PHE:CD2	2.55	0.42
2:B:976:VAL:HG22	2:B:996:GLY:HA3	2.02	0.42
2:B:144:PRO:O	2:B:152:LEU:HD23	2.20	0.42
2:B:478:LEU:HD12	2:B:479:VAL:H	1.84	0.42
2:B:1095:GLU:OE2	2:B:1140:HIS:NE2	2.44	0.42
2:B:1097:PHE:CE1	2:B:1105:MET:HG2	2.55	0.42
1:A:57:ARG:HG2	1:A:57:ARG:HH11	1.85	0.41
2:B:375:LEU:HB2	2:B:390:ILE:HB	2.01	0.41
2:B:605:ALA:HB2	2:B:611:LEU:HD23	2.01	0.41
2:B:770:LEU:HD22	2:B:863:GLU:OE1	2.20	0.41
2:B:870:VAL:HG22	2:B:884:ILE:HD12	2.01	0.41
2:B:1064:SER:HG	2:B:1068:ILE:N	2.16	0.41
2:B:64:MET:HG3	2:B:77:LEU:HD11	2.02	0.41
2:B:413:LEU:HD13	2:B:461:GLY:HA2	2.02	0.41
1:A:2:LEU:HD22	2:B:814:LEU:HD21	2.01	0.41
1:A:23:SER:CB	1:A:347:GLN:HE21	2.34	0.41
2:B:849:VAL:HG11	2:B:851:PHE:CZ	2.54	0.41
2:B:451:PHE:HA	2:B:470:GLN:OE1	2.21	0.41
2:B:514:ARG:HD2	2:B:536:HIS:HA	2.03	0.41
2:B:368:GLU:CG	2:B:674:LYS:HE3	2.45	0.41
1:A:165:GLY:HA2	1:A:166:PRO:HD3	1.94	0.41
2:B:131:ILE:HB	2:B:143:ILE:HB	2.03	0.41
2:B:226:PHE:CZ	2:B:287:LYS:HG2	2.55	0.41
2:B:650:PHE:CD2	2:B:679:MET:HG3	2.56	0.41
2:B:720:SER:OG	2:B:738:SER:HB3	2.20	0.41
1:A:308:VAL:HG12	1:A:310:VAL:HG13	2.02	0.41
2:B:59:GLY:HA2	2:B:1073:TRP:CZ3	2.56	0.41
2:B:492:GLU:HG2	2:B:493:PRO:HD2	2.02	0.41
2:B:650:PHE:HD1	2:B:651:ALA:N	2.19	0.41
2:B:196:SER:O	2:B:200:LYS:HA	2.21	0.41
2:B:1057:ARG:NH1	2:B:1112:LEU:HG	2.36	0.41
2:B:288:GLU:HB2	2:B:298:LYS:HB2	2.02	0.40
2:B:292:ASP:OD1	2:B:293:GLY:N	2.54	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:452:VAL:HG12	2:B:470:GLN:NE2	2.37	0.40
2:B:512:VAL:HG23	2:B:512:VAL:O	2.21	0.40
2:B:1007:PHE:CD1	2:B:1030:PHE:HB3	2.56	0.40
2:B:1081:LYS:HG2	2:B:1083:GLU:OE1	2.21	0.40
2:B:1134:GLU:HA	2:B:1137:THR:HG22	2.02	0.40
2:B:264:VAL:HB	2:B:270:ARG:HB2	2.03	0.40
2:B:446:THR:HG22	2:B:447:GLU:N	2.35	0.40
2:B:727:GLN:HE21	2:B:827:THR:HG22	1.87	0.40
2:B:448:LEU:HA	2:B:484:LYS:HZ3	1.86	0.40
3:C:58:TYR:HA	3:C:61:GLN:OE1	2.22	0.40
2:B:123:ILE:HD11	2:B:140:PHE:HZ	1.87	0.40
2:B:682:LEU:HD22	2:B:684:SER:HB2	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	354/408 (87%)	332 (94%)	22 (6%)	0	100	100
2	B	1112/1160 (96%)	1053 (95%)	58 (5%)	1 (0%)	51	82
3	C	45/152 (30%)	44 (98%)	1 (2%)	0	100	100
4	D	148/729 (20%)	148 (100%)	0	0	100	100
All	All	1659/2449 (68%)	1577 (95%)	81 (5%)	1 (0%)	54	82

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	671	VAL

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	313/358 (87%)	312 (100%)	1 (0%)	92	97
2	B	983/1014 (97%)	979 (100%)	4 (0%)	91	95
3	C	48/128 (38%)	48 (100%)	0	100	100
4	D	134/623 (22%)	134 (100%)	0	100	100
All	All	1478/2123 (70%)	1473 (100%)	5 (0%)	92	97

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	164	ARG
2	B	423	ASP
2	B	650	PHE
2	B	680	CYS
2	B	699	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	42	HIS
1	A	47	ASN
1	A	230	ASN
1	A	244	HIS
2	B	370	GLN
2	B	467	GLN
2	B	677	ASN
2	B	796	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)

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.

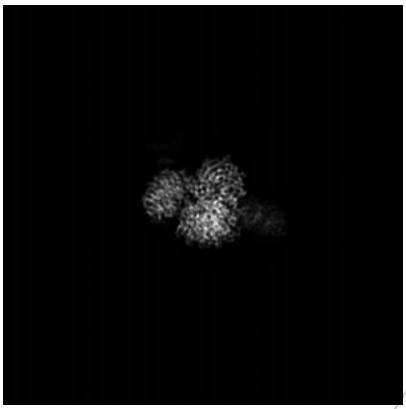
6 Map visualisation

This section contains visualisations of the EMDB entry EMD-18398. These allow visual inspection of the internal detail of the map and identification of artifacts.

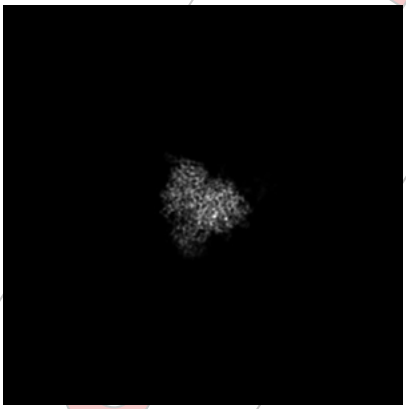
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

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: 160

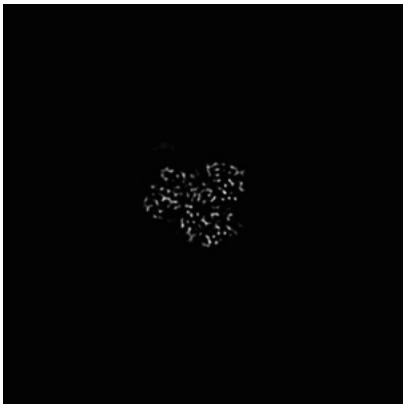
Y Index: 160

Z Index: 160

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: 166

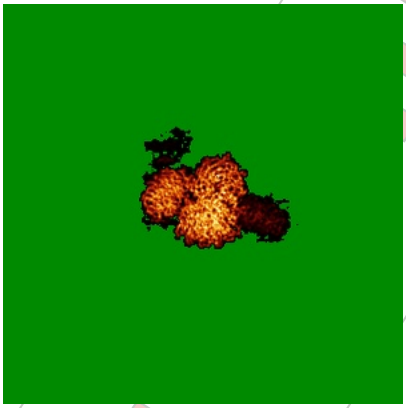
Y Index: 168

Z Index: 157

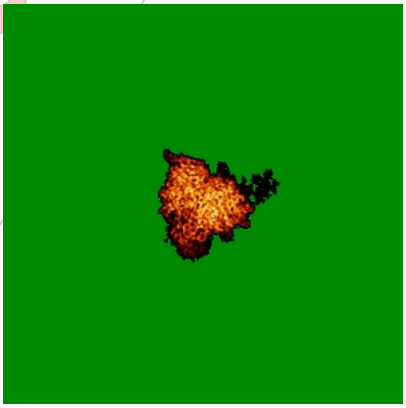
The images above show the largest variance slices of the map in three orthogonal directions.

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

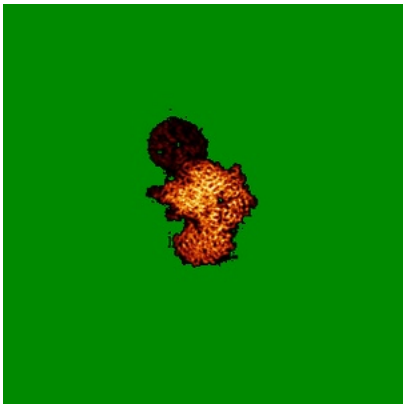
6.4.1 Primary map



X



Y

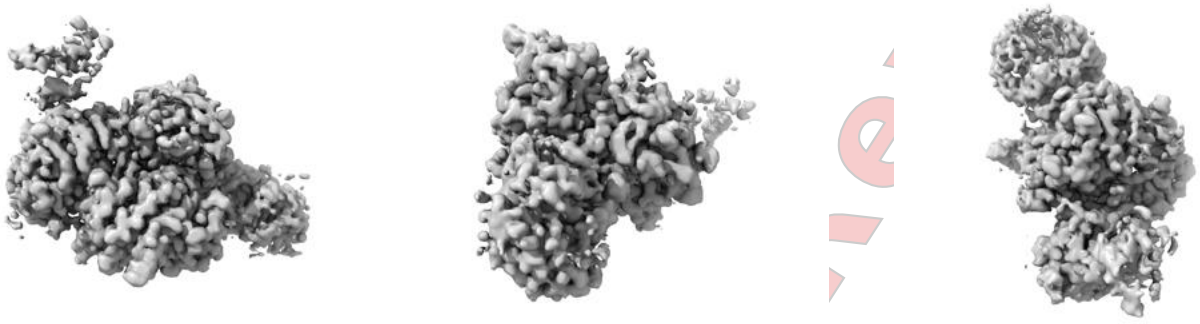


Z

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

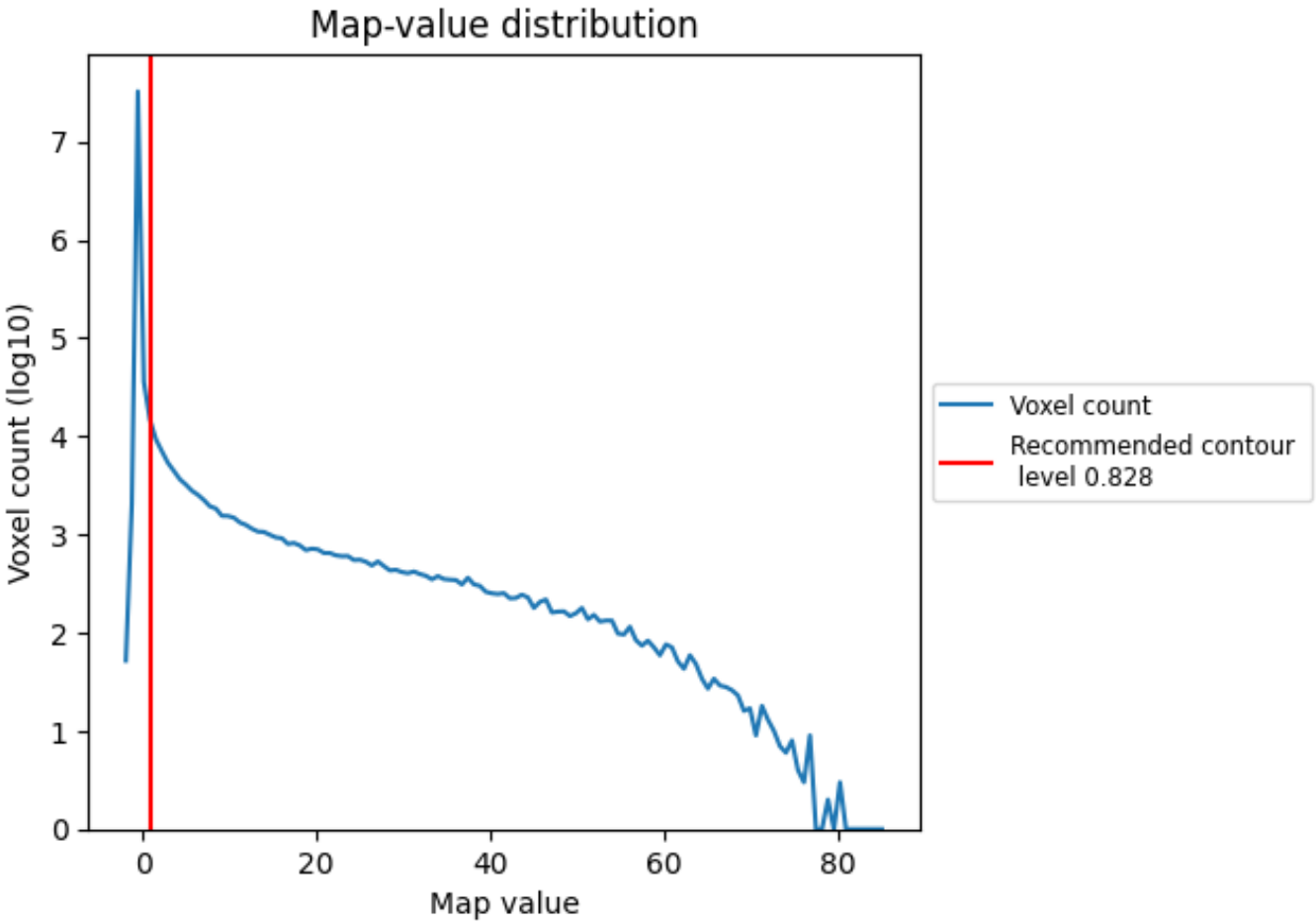
6.6 Mask visualisation (i)

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

7 Map analysis

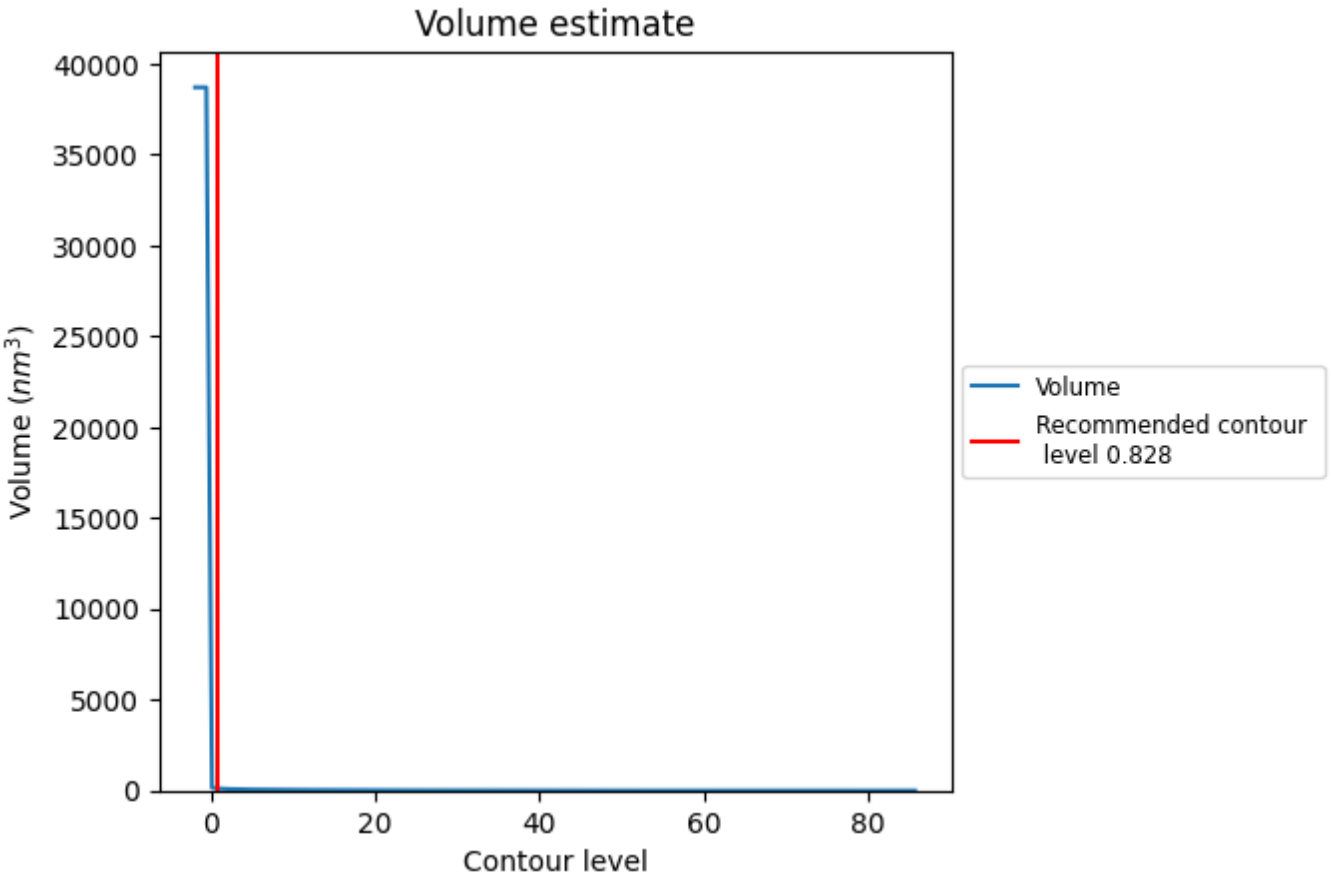
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

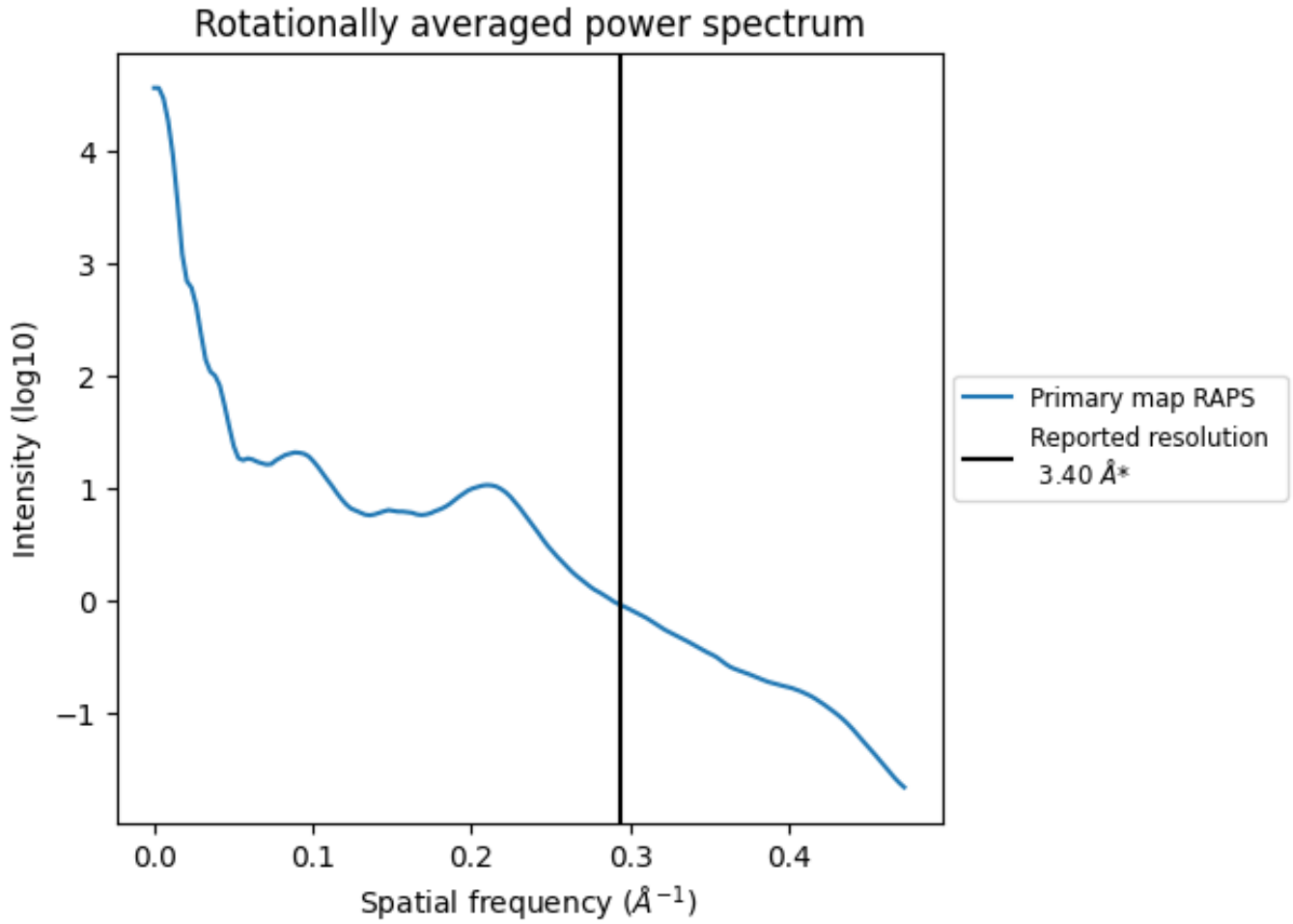
7.2 Volume estimate (i)



The volume at the recommended contour level is 113 nm^3; this corresponds to an approximate mass of 102 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 0.294 $Å^{-1}$

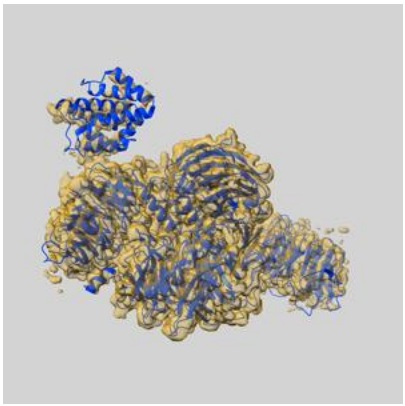
8 Fourier-Shell correlation (i)

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

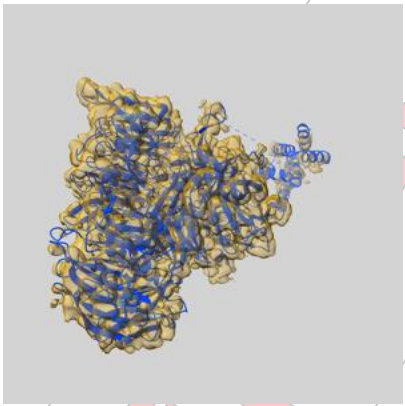
9 Map-model fit

This section contains information regarding the fit between EMDB map EMD-18398 and PDB model 8QH5. Per-residue inclusion information can be found in section 3 on page 7.

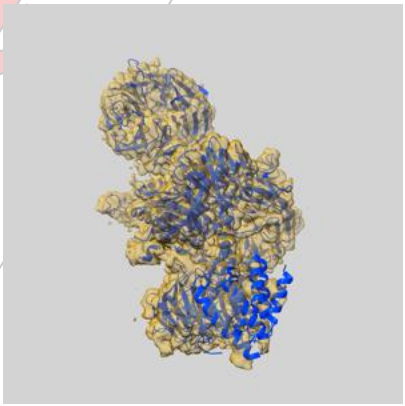
9.1 Map-model overlay



X



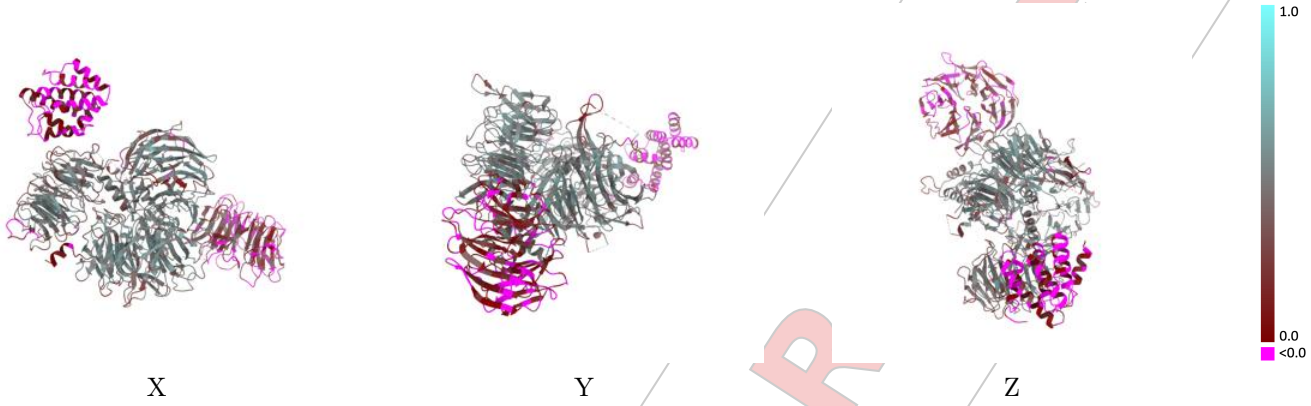
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.828 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.

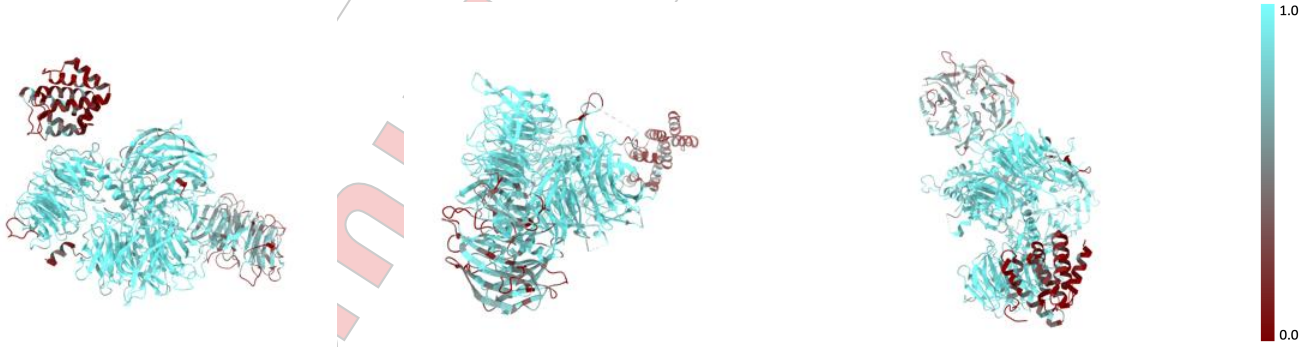
9.2 Q-score mapped to coordinate model



X Y Z

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

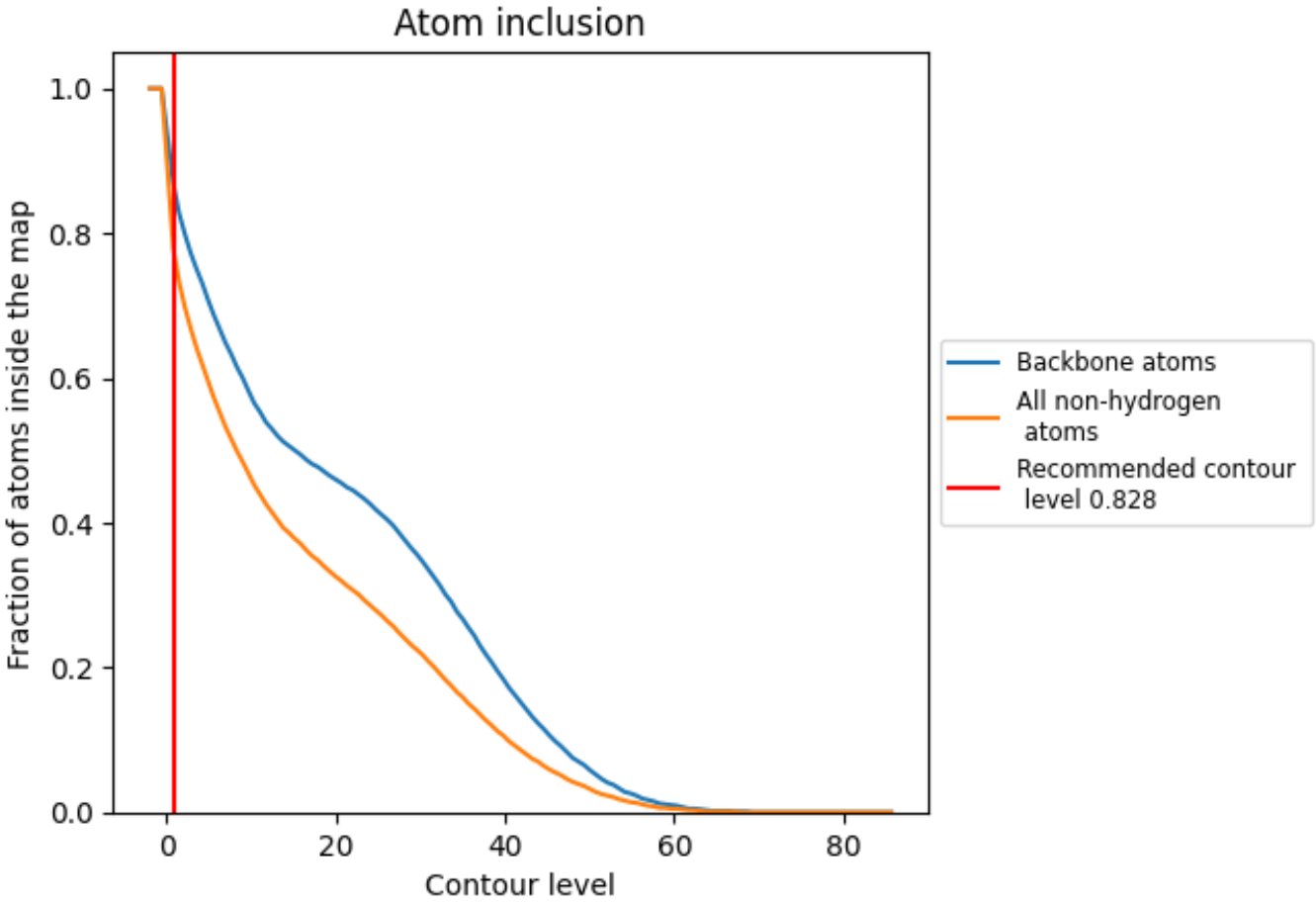
9.3 Atom inclusion mapped to coordinate model



X Y Z

The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.828).

9.4 Atom inclusion (i)



At the recommended contour level, 87% of all backbone atoms, 78% 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 (0.828) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7800	0.3310
A	0.8930	0.3990
B	0.8250	0.3570
C	0.7620	0.3280
D	0.2000	-0.0030