



Full wwPDB EM Validation Report ⓘ

Mar 9, 2026 – 03:47 PM UTC

PDB ID : 6NUD / pdb_00006nud
EMDB ID : EMD-0516
Title : Small conformation of ssRNA-bound CRISPR_Csm complex
Authors : Zhang, K.; Pintilie, G.; Li, S.; Zhu, Y.; Chiu, W.; Huang, Z.
Deposited on : 2019-01-31
Resolution : 3.50 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:

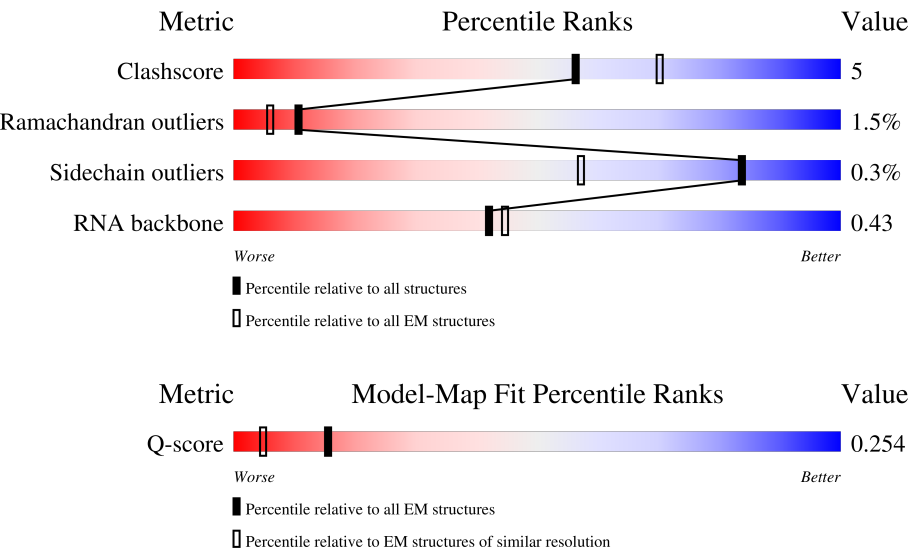
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

1 Overall quality at a glance i

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

The reported resolution of this entry is 3.50 Å.

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



Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
RNA backbone	8273	3508	-
Q-score	-	25397	13950 (3.00 - 4.00)

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	121	42% 79% 11% 11%
1	B	121	71% 80% 9% 11%
1	M	121	54% 79% 11% 11%

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Mol	Chain	Length	Quality of chain
2	H	72	
3	I	299	
4	J	758	
5	U	49	
6	C	220	
6	E	220	
6	N	220	
6	O	220	
6	P	220	

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 20831 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 CRISPR system Cms protein Csm2.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	108	Total	C	N	O	S	0	0
			894	572	154	165	3		
1	B	108	Total	C	N	O	S	0	0
			894	572	154	165	3		
1	M	108	Total	C	N	O	S	0	0
			894	572	154	165	3		

- Molecule 2 is a RNA chain called crRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	H	40	Total	C	N	O	P	0	0
			832	375	133	285	39		

- Molecule 3 is a protein called CRISPR type III-associated RAMP protein Csm4.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	I	295	Total	C	N	O	S	0	0
			2352	1503	388	456	5		

- Molecule 4 is a protein called CRISPR system single-strand-specific deoxyribonuclease Cas10/Csm1 (subtype III-A).

Mol	Chain	Residues	Atoms					AltConf	Trace
4	J	741	Total	C	N	O	S	0	0
			5997	3842	1000	1140	15		

- Molecule 5 is a RNA chain called target ssRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	U	28	Total	C	N	O	P	0	0
			602	272	119	184	27		

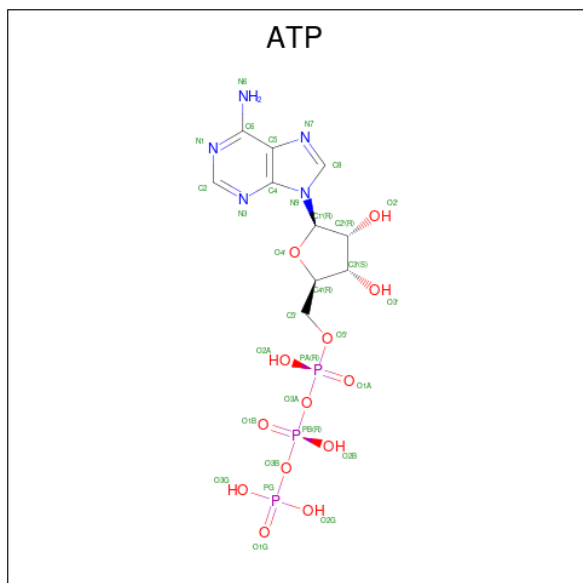
- Molecule 6 is a protein called CRISPR type III-associated RAMP protein Csm3.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	C	212	Total	C	N	O	S	0	0
			1667	1057	285	323	2		
6	E	212	Total	C	N	O	S	0	0
			1667	1057	285	323	2		
6	N	212	Total	C	N	O	S	0	0
			1667	1057	285	323	2		
6	O	212	Total	C	N	O	S	0	0
			1667	1057	285	323	2		
6	P	212	Total	C	N	O	S	0	0
			1667	1057	285	323	2		

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	33	ALA	ASP	engineered mutation	UNP A0A0A7HIF0
E	33	ALA	ASP	engineered mutation	UNP A0A0A7HIF0
N	33	ALA	ASP	engineered mutation	UNP A0A0A7HIF0
O	33	ALA	ASP	engineered mutation	UNP A0A0A7HIF0
P	33	ALA	ASP	engineered mutation	UNP A0A0A7HIF0

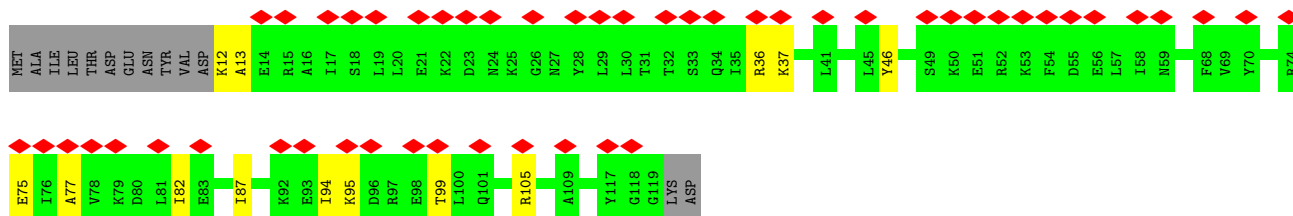
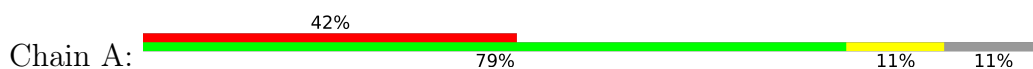
- Molecule 7 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$).



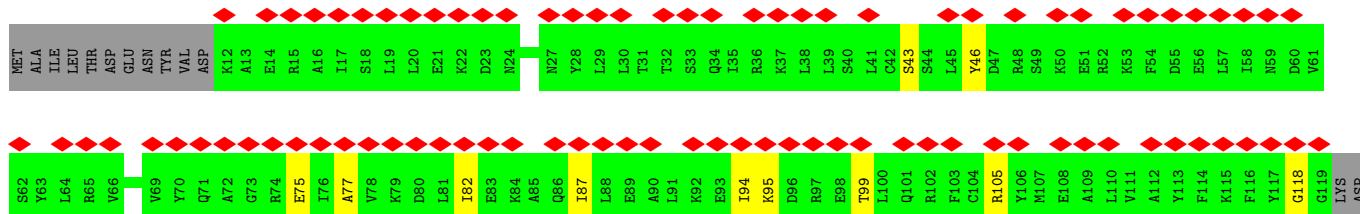
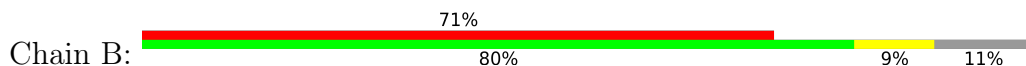
3 Residue-property plots

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.

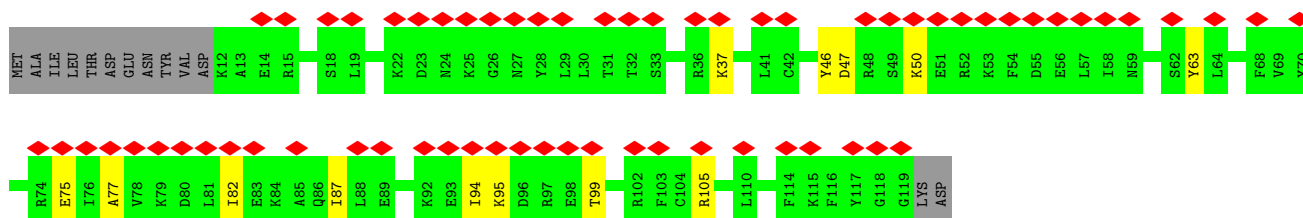
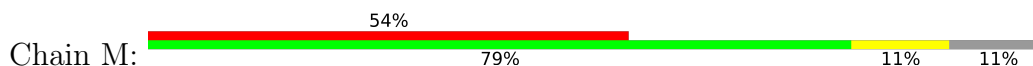
- Molecule 1: CRISPR system Cms protein Csm2



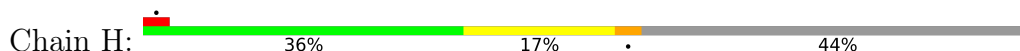
- Molecule 1: CRISPR system Cms protein Csm2

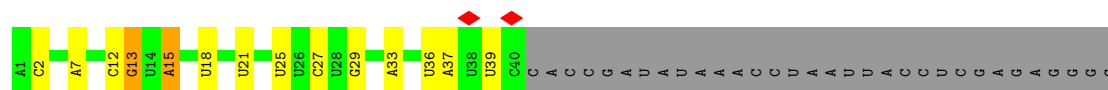


- Molecule 1: CRISPR system Cms protein Csm2

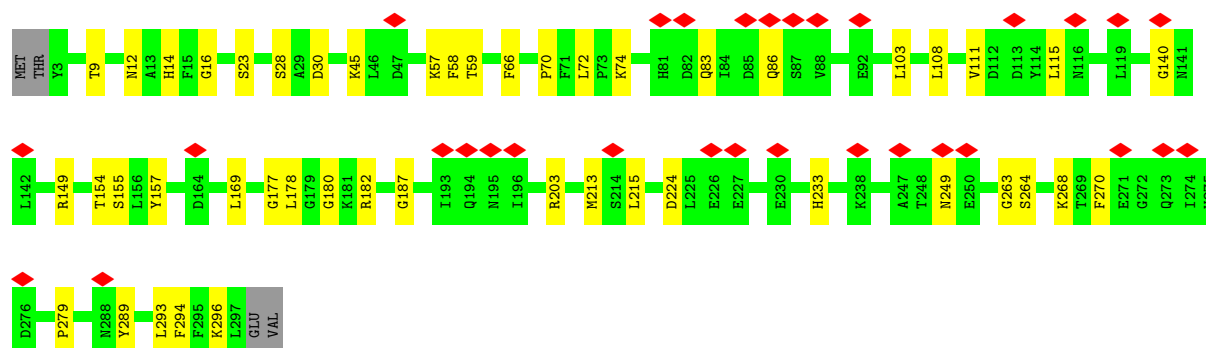
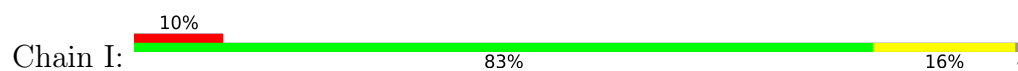


- Molecule 2: crRNA

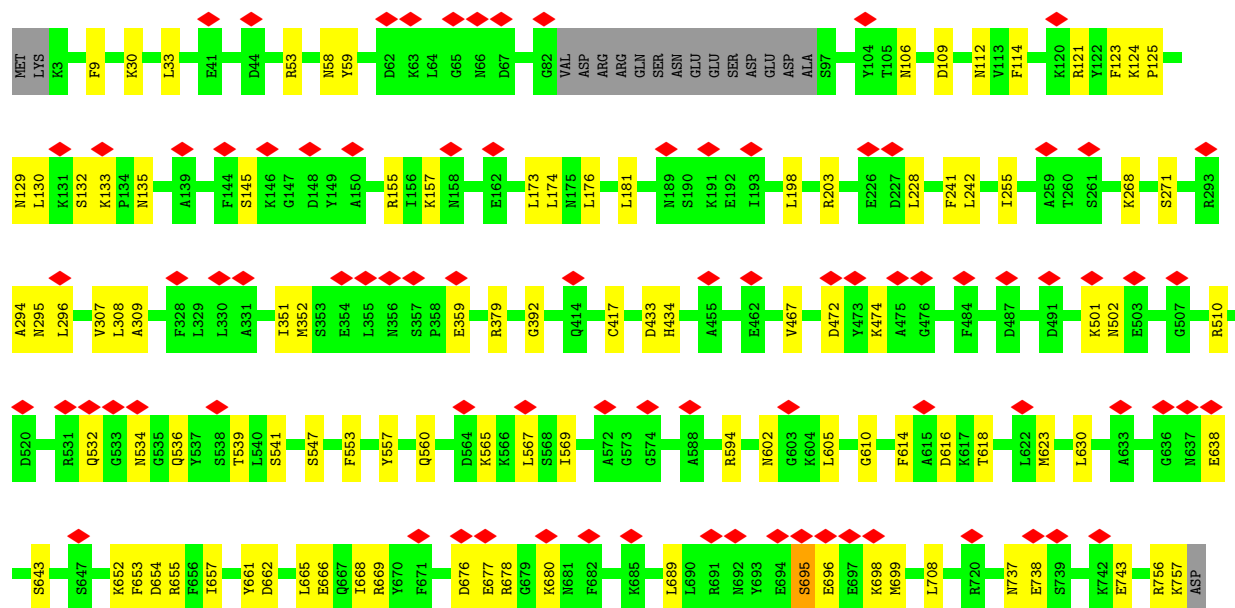
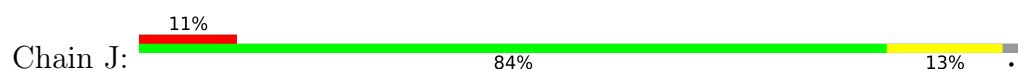




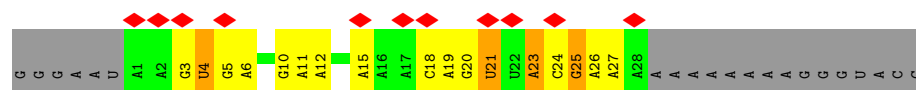
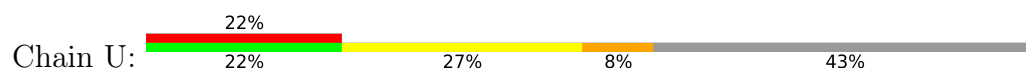
- Molecule 3: CRISPR type III-associated RAMP protein Csm4



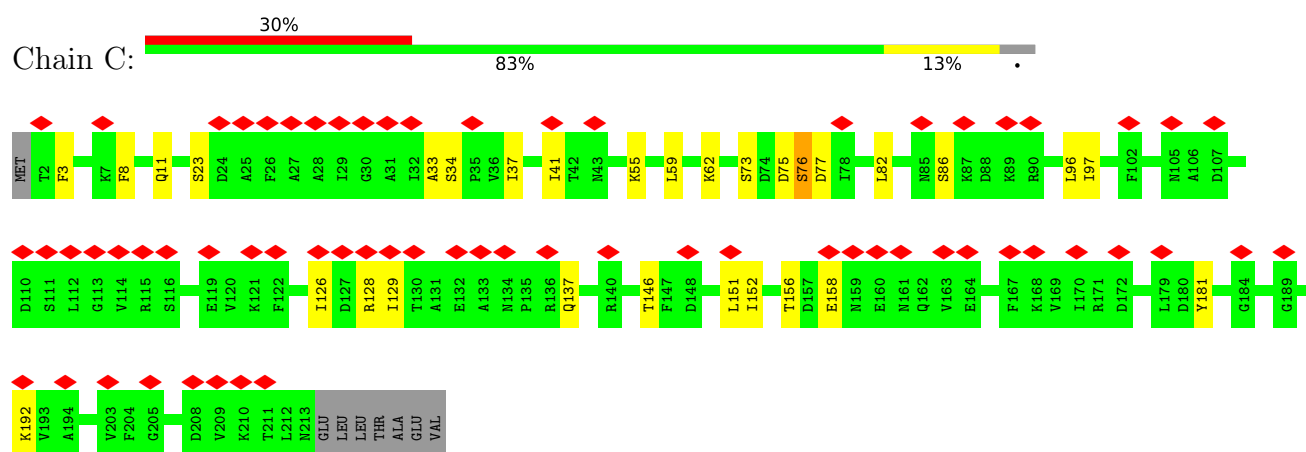
- Molecule 4: CRISPR system single-strand-specific deoxyribonuclease Cas10/Csm1 (subtype III-A)



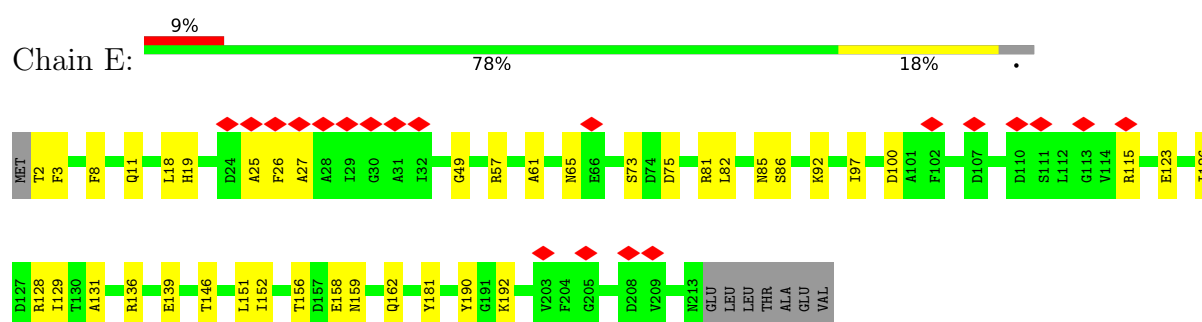
- Molecule 5: target ssRNA



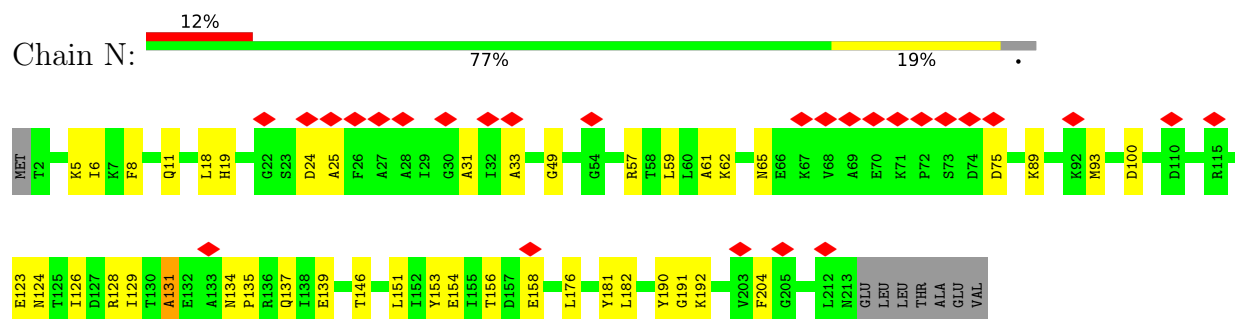
- Molecule 6: CRISPR type III-associated RAMP protein Csm3



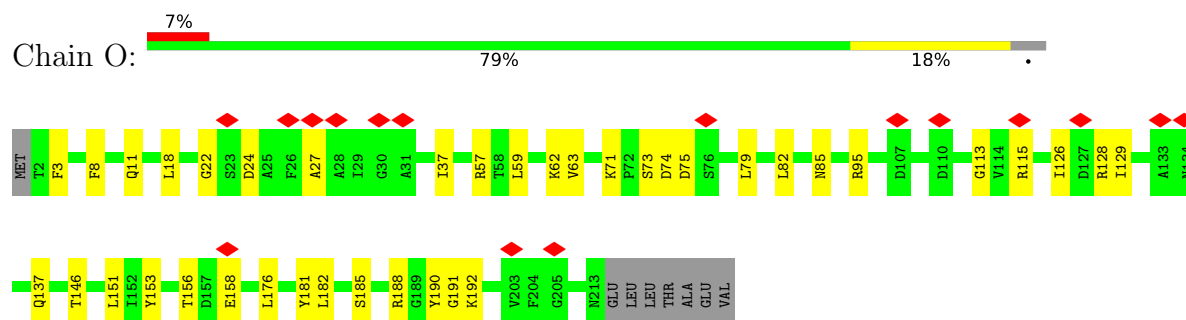
- Molecule 6: CRISPR type III-associated RAMP protein Csm3



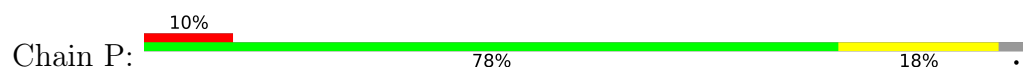
- Molecule 6: CRISPR type III-associated RAMP protein Csm3

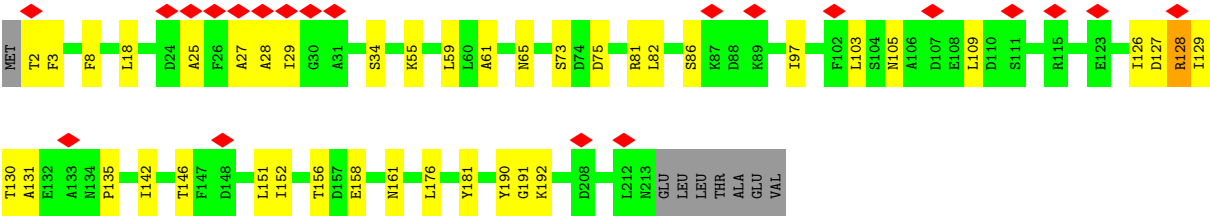


- Molecule 6: CRISPR type III-associated RAMP protein Csm3



- Molecule 6: CRISPR type III-associated RAMP protein Csm3





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	50092	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 ($e^-/\text{\AA}^2$)	7	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	3.348	Depositor
Minimum map value	-1.954	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.099	Depositor
Recommended contour level	0.5	Depositor
Map size (Å)	407.03998, 407.03998, 407.03998	wwPDB
Map dimensions	384, 384, 384	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.06, 1.06, 1.06	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: ATP

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	0.08	0/905	0.25	0/1209
1	B	0.07	0/905	0.21	0/1209
1	M	0.07	0/905	0.22	0/1209
2	H	0.07	0/926	0.19	0/1437
3	I	0.07	0/2404	0.22	0/3247
4	J	0.07	0/6123	0.22	0/8248
5	U	0.06	0/677	0.18	0/1054
6	C	0.09	0/1692	0.27	0/2275
6	E	0.09	0/1692	0.23	0/2275
6	N	0.09	0/1692	0.24	0/2275
6	O	0.09	0/1692	0.26	1/2275 (0.0%)
6	P	0.08	0/1692	0.25	0/2275
All	All	0.08	0/21305	0.23	1/28988 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	O	63	VAL	N-CA-C	-5.02	108.94	113.71

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	894	0	923	8	0
1	B	894	0	923	5	0
1	M	894	0	923	7	0
2	H	832	0	425	9	0
3	I	2352	0	2298	26	0
4	J	5997	0	5865	52	0
5	U	602	0	305	10	0
6	C	1667	0	1685	17	0
6	E	1667	0	1685	26	0
6	N	1667	0	1685	26	0
6	O	1667	0	1685	22	0
6	P	1667	0	1685	26	0
7	J	31	0	12	1	0
All	All	20831	0	20099	201	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 5.

All (201) 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)
6:C:8:PHE:HB2	6:C:151:LEU:HB3	1.74	0.70
5:U:23:A:H1'	6:N:135:PRO:HD2	1.73	0.69
6:E:92:LYS:HE2	6:N:131:ALA:HB2	1.76	0.68
6:P:8:PHE:HB2	6:P:151:LEU:HB3	1.77	0.67
6:E:181:TYR:HA	6:E:192:LYS:H	1.60	0.66
3:I:249:ASN:HA	6:N:89:LYS:HE3	1.78	0.65
6:N:126:ILE:HG22	6:N:128:ARG:H	1.63	0.64
6:O:126:ILE:HG22	6:O:128:ARG:H	1.63	0.64
6:P:181:TYR:HA	6:P:192:LYS:H	1.63	0.64
6:O:8:PHE:HB2	6:O:151:LEU:HB3	1.79	0.63
6:N:8:PHE:HB2	6:N:151:LEU:HB3	1.79	0.63
6:E:8:PHE:HB2	6:E:151:LEU:HB3	1.81	0.63
6:O:181:TYR:HA	6:O:192:LYS:H	1.64	0.62
6:E:126:ILE:HG22	6:E:128:ARG:H	1.64	0.62
1:M:37:LYS:HB2	6:P:25:ALA:HB1	1.84	0.60
6:P:61:ALA:O	6:P:65:ASN:ND2	2.35	0.59
6:E:18:LEU:HA	6:E:190:TYR:HB3	1.85	0.59
4:J:657:ILE:HA	4:J:661:TYR:HB3	1.85	0.59
6:N:181:TYR:HA	6:N:192:LYS:H	1.68	0.59
2:H:15:A:N6	6:O:24:ASP:OD2	2.36	0.58
3:I:59:THR:HG21	3:I:293:LEU:HD21	1.86	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:C:23:SER:HB3	6:C:37:ILE:HD13	1.85	0.58
2:H:15:A:O2'	6:N:124:ASN:ND2	2.38	0.57
6:C:126:ILE:HG22	6:C:128:ARG:H	1.70	0.57
6:P:18:LEU:HA	6:P:190:TYR:HB3	1.87	0.57
6:O:182:LEU:H	6:O:191:GLY:HA3	1.69	0.56
6:O:18:LEU:HA	6:O:190:TYR:HB3	1.87	0.56
6:N:18:LEU:HA	6:N:190:TYR:HB3	1.88	0.56
6:E:26:PHE:O	6:O:115:ARG:NH2	2.39	0.55
5:U:4:U:H2'	6:P:135:PRO:HG3	1.88	0.55
1:A:75:GLU:HG2	1:A:77:ALA:H	1.72	0.55
6:E:19:HIS:ND1	6:E:139:GLU:O	2.38	0.55
3:I:70:PRO:HD2	3:I:108:LEU:HD13	1.89	0.55
4:J:501:LYS:NZ	4:J:757:LYS:O	2.40	0.54
1:A:37:LYS:HB2	6:E:25:ALA:HB1	1.88	0.54
6:N:61:ALA:O	6:N:65:ASN:ND2	2.41	0.54
6:P:127:ASP:O	6:P:129:ILE:N	2.41	0.54
2:H:7:A:OP1	6:N:57:ARG:NH1	2.41	0.54
1:A:12:LYS:NZ	4:J:699:MET:SD	2.81	0.54
2:H:13:G:OP2	6:O:57:ARG:NH1	2.41	0.54
6:E:73:SER:O	6:E:85:ASN:ND2	2.41	0.53
4:J:618:THR:OG1	4:J:623:MET:SD	2.67	0.53
6:O:11:GLN:NE2	6:O:146:THR:OG1	2.42	0.53
4:J:565:LYS:HD2	4:J:567:LEU:HD23	1.90	0.53
4:J:678:ARG:NH1	5:U:20:G:OP2	2.42	0.53
1:B:75:GLU:HG2	1:B:77:ALA:H	1.74	0.53
1:A:36:ARG:NH1	6:O:137:GLN:OE1	2.42	0.53
4:J:614:PHE:HB2	4:J:623:MET:HE3	1.91	0.53
1:M:47:ASP:HA	1:M:50:LYS:HE3	1.91	0.52
6:O:156:THR:O	6:O:158:GLU:N	2.41	0.52
2:H:13:G:N1	5:U:25:G:O6	2.42	0.52
3:I:45:LYS:HD2	3:I:169:LEU:HD21	1.90	0.52
4:J:121:ARG:HD2	4:J:602:ASN:HB2	1.91	0.52
4:J:652:LYS:O	4:J:654:ASP:N	2.43	0.52
1:A:82:ILE:HG23	1:A:87:ILE:HB	1.90	0.52
6:N:25:ALA:HB2	6:N:31:ALA:HB2	1.91	0.52
4:J:124:LYS:O	4:J:135:ASN:ND2	2.42	0.52
6:P:126:ILE:HG22	6:P:128:ARG:H	1.75	0.52
6:E:115:ARG:NH2	6:P:27:ALA:O	2.43	0.52
6:E:156:THR:O	6:E:158:GLU:N	2.43	0.51
1:A:46:TYR:OH	1:A:105:ARG:NH2	2.44	0.51
3:I:177:GLY:HA3	3:I:182:ARG:HG3	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:I:213:MET:HE3	3:I:215:LEU:HD11	1.92	0.51
4:J:124:LYS:HD2	4:J:125:PRO:HD2	1.93	0.51
3:I:16:GLY:HA2	3:I:23:SER:HA	1.92	0.51
3:I:66:PHE:HB2	3:I:157:TYR:HD1	1.76	0.51
3:I:182:ARG:NH1	6:N:93:MET:SD	2.83	0.51
1:M:46:TYR:OH	1:M:105:ARG:NH2	2.44	0.51
6:P:156:THR:O	6:P:158:GLU:N	2.43	0.51
5:U:20:G:N2	6:E:86:SER:OG	2.44	0.50
6:C:181:TYR:HA	6:C:192:LYS:H	1.77	0.50
6:E:97:ILE:HB	6:E:152:ILE:HB	1.92	0.50
3:I:233:HIS:O	3:I:264:SER:OG	2.29	0.50
1:M:75:GLU:HG2	1:M:77:ALA:H	1.75	0.50
6:C:73:SER:HB2	6:C:86:SER:H	1.76	0.50
6:P:81:ARG:HG2	6:P:82:LEU:HD12	1.94	0.50
3:I:270:PHE:H	3:I:294:PHE:HZ	1.58	0.50
4:J:594:ARG:NH1	4:J:654:ASP:OD1	2.45	0.50
6:E:11:GLN:NE2	6:E:146:THR:OG1	2.44	0.50
3:I:178:LEU:H	3:I:187:GLY:HA3	1.77	0.49
6:P:73:SER:HB2	6:P:86:SER:H	1.76	0.49
1:B:82:ILE:HG23	1:B:87:ILE:HB	1.94	0.49
3:I:66:PHE:HB3	3:I:155:SER:HB3	1.93	0.49
6:N:11:GLN:NE2	6:N:146:THR:OG1	2.44	0.49
6:O:59:LEU:HA	6:O:62:LYS:HE3	1.94	0.49
4:J:630:LEU:HB3	4:J:643:SER:HB3	1.94	0.49
1:B:46:TYR:OH	1:B:105:ARG:NH2	2.46	0.49
4:J:106:ASN:HB2	4:J:145:SER:HB2	1.95	0.49
4:J:680:LYS:HB3	5:U:23:A:H62	1.78	0.49
6:C:11:GLN:NE2	6:C:146:THR:OG1	2.45	0.49
6:E:3:PHE:HE2	6:O:176:LEU:HD11	1.78	0.49
6:C:3:PHE:HE2	6:P:176:LEU:HD11	1.78	0.49
6:E:159:ASN:ND2	6:E:162:GLN:OE1	2.46	0.49
4:J:129:ASN:HD22	4:J:130:LEU:H	1.61	0.48
4:J:666:GLU:HG3	4:J:669:ARG:HH21	1.78	0.48
6:C:41:ILE:HD12	6:P:109:LEU:HD11	1.95	0.48
4:J:737:ASN:HD22	4:J:738:GLU:H	1.61	0.48
6:N:6:ILE:HB	6:N:153:TYR:HB3	1.96	0.48
6:C:82:LEU:HD23	6:C:96:LEU:HD11	1.96	0.47
3:I:86:GLN:NE2	4:J:359:GLU:OE2	2.47	0.47
6:O:185:SER:HB2	6:O:188:ARG:HH12	1.79	0.47
4:J:676:ASP:O	4:J:678:ARG:N	2.48	0.47
6:C:76:SER:OG	6:C:77:ASP:N	2.48	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:E:159:ASN:HD21	6:E:162:GLN:HB2	1.79	0.47
6:N:156:THR:O	6:N:158:GLU:N	2.42	0.47
6:O:22:GLY:HA3	6:O:37:ILE:HB	1.96	0.47
6:P:127:ASP:HB3	6:P:130:THR:HG22	1.96	0.47
6:C:156:THR:O	6:C:158:GLU:N	2.42	0.47
1:B:43:SER:HA	6:P:29:ILE:HD13	1.97	0.47
3:I:203:ARG:HD3	3:I:296:LYS:HB2	1.97	0.47
6:E:61:ALA:O	6:E:65:ASN:ND2	2.48	0.47
4:J:109:ASP:HA	4:J:123:PHE:HE2	1.79	0.47
4:J:547:SER:HB2	7:J:801:ATP:HN62	1.80	0.46
6:E:49:GLY:N	6:E:100:ASP:OD1	2.47	0.46
4:J:532:GLN:HG2	4:J:534:ASN:H	1.79	0.46
6:C:33:ALA:HB2	6:C:137:GLN:HG3	1.97	0.46
6:E:81:ARG:HG2	6:E:82:LEU:HD12	1.98	0.46
6:O:73:SER:O	6:O:85:ASN:ND2	2.48	0.46
1:A:13:ALA:HB3	4:J:698:LYS:HB3	1.98	0.46
3:I:12:ASN:H	3:I:149:ARG:HA	1.81	0.46
3:I:12:ASN:HB2	3:I:149:ARG:HH11	1.79	0.46
3:I:72:LEU:HD12	3:I:111:VAL:HG11	1.97	0.46
4:J:242:LEU:HB3	4:J:308:LEU:HB2	1.98	0.46
4:J:510:ARG:NH1	6:N:24:ASP:OD2	2.49	0.46
6:N:176:LEU:HD11	6:O:3:PHE:HE2	1.79	0.46
6:P:75:ASP:N	6:P:75:ASP:OD1	2.48	0.45
6:P:97:ILE:HB	6:P:152:ILE:HB	1.98	0.45
4:J:132:SER:O	4:J:560:GLN:NE2	2.49	0.45
6:N:49:GLY:N	6:N:100:ASP:OD1	2.47	0.45
6:N:123:GLU:OE2	6:N:190:TYR:OH	2.31	0.45
6:E:75:ASP:N	6:E:75:ASP:OD1	2.49	0.45
4:J:30:LYS:HD2	4:J:33:LEU:HD23	1.97	0.45
4:J:174:LEU:HB3	4:J:203:ARG:HH12	1.81	0.45
1:B:94:ILE:HG22	1:B:99:THR:HG22	1.98	0.45
4:J:203:ARG:HH21	4:J:296:LEU:HB3	1.82	0.45
4:J:434:HIS:HA	4:J:467:VAL:HG12	1.98	0.45
6:N:75:ASP:OD1	6:N:75:ASP:N	2.49	0.45
4:J:9:PHE:HE1	4:J:294:ALA:HB2	1.80	0.45
6:O:75:ASP:OD1	6:O:75:ASP:N	2.48	0.45
1:A:94:ILE:HG22	1:A:99:THR:HG22	1.99	0.45
2:H:25:U:O2	6:E:136:ARG:NE	2.50	0.45
1:M:82:ILE:HG23	1:M:87:ILE:HB	1.97	0.45
4:J:30:LYS:NZ	4:J:58:ASN:OD1	2.46	0.45
6:P:55:LYS:HE3	6:P:59:LEU:HD11	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:94:ILE:HG22	1:M:99:THR:HG22	1.98	0.45
6:O:71:LYS:HZ2	6:O:74:ASP:HB2	1.81	0.45
2:H:36:U:O4	2:H:37:A:N6	2.51	0.44
4:J:680:LYS:HZ1	4:J:756:ARG:HA	1.82	0.44
3:I:83:GLN:NE2	3:I:115:LEU:O	2.43	0.44
6:P:105:ASN:HB2	6:P:146:THR:HG22	2.00	0.44
4:J:155:ARG:NH1	4:J:472:ASP:OD1	2.51	0.43
6:C:96:LEU:HD22	6:C:151:LEU:HD11	1.99	0.43
3:I:28:SER:OG	3:I:30:ASP:OD1	2.36	0.43
5:U:20:G:O2'	5:U:21:U:O2	2.31	0.43
6:N:19:HIS:ND1	6:N:139:GLU:O	2.49	0.43
6:N:182:LEU:H	6:N:191:GLY:HA3	1.83	0.43
6:C:59:LEU:HA	6:C:62:LYS:HE3	2.01	0.43
6:E:123:GLU:OE2	6:E:190:TYR:OH	2.35	0.43
3:I:215:LEU:HG	3:I:263:GLY:H	1.84	0.43
5:U:19:A:N6	5:U:20:G:O6	2.52	0.43
6:C:97:ILE:HB	6:C:152:ILE:HB	1.99	0.43
6:O:95:ARG:NH1	6:O:153:TYR:OH	2.51	0.43
6:P:103:LEU:HD11	6:P:142:ILE:HD11	2.00	0.43
4:J:133:LYS:HA	4:J:560:GLN:HE22	1.83	0.43
4:J:553:PHE:HA	4:J:557:TYR:HB2	2.00	0.43
4:J:352:MET:HA	4:J:536:GLN:HE22	1.84	0.43
4:J:295:ASN:ND2	4:J:307:VAL:O	2.51	0.43
3:I:14:HIS:HB3	3:I:180:GLY:HA3	2.00	0.43
6:P:2:THR:OG1	6:P:3:PHE:N	2.50	0.42
4:J:241:PHE:HA	4:J:309:ALA:HA	2.01	0.42
4:J:268:LYS:NZ	4:J:569:ILE:O	2.40	0.42
4:J:114:PHE:HE2	4:J:605:LEU:HD21	1.84	0.42
5:U:23:A:O2'	6:N:134:ASN:OD1	2.32	0.42
6:O:18:LEU:HB3	6:O:191:GLY:HA2	2.01	0.42
4:J:181:LEU:HD23	4:J:198:LEU:HD23	2.02	0.42
6:P:161:ASN:OD1	6:P:161:ASN:N	2.51	0.42
3:I:9:THR:OG1	3:I:154:THR:O	2.37	0.42
4:J:433:ASP:OD2	4:J:474:LYS:NZ	2.46	0.42
4:J:695:SER:OG	4:J:696:GLU:N	2.50	0.42
1:M:63:TYR:OH	6:P:28:ALA:O	2.34	0.42
2:H:18:U:O2'	6:E:57:ARG:NH1	2.52	0.42
4:J:255:ILE:HD13	4:J:271:SER:HA	2.01	0.42
6:C:41:ILE:HD11	6:P:142:ILE:HB	2.00	0.42
6:C:55:LYS:HE3	6:C:59:LEU:HD11	2.00	0.42
4:J:510:ARG:HH21	4:J:616:ASP:H	1.66	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:18:U:HI'	6:E:73:SER:HB3	2.01	0.41
4:J:539:THR:HG22	4:J:541:SER:H	1.84	0.41
5:U:10:G:N2	5:U:11:A:N7	2.68	0.41
3:I:58:PHE:O	3:I:289:TYR:OH	2.37	0.41
4:J:351:ILE:O	4:J:536:GLN:NE2	2.53	0.41
6:P:18:LEU:HB3	6:P:191:GLY:HA2	2.03	0.41
4:J:9:PHE:HZ	4:J:174:LEU:HD11	1.85	0.41
4:J:109:ASP:OD2	4:J:112:ASN:N	2.54	0.41
3:I:224:ASP:OD1	3:I:268:LYS:NZ	2.52	0.41
3:I:74:LYS:HD3	3:I:103:LEU:HB3	2.03	0.41
6:N:5:LYS:HD3	6:N:154:GLU:HB3	2.03	0.41
6:N:33:ALA:HA	6:N:137:GLN:HG3	2.02	0.41
6:N:59:LEU:HA	6:N:62:LYS:HE3	2.03	0.41
4:J:655:ARG:NE	4:J:743:GLU:OE2	2.54	0.41
6:O:79:LEU:HA	6:O:82:LEU:HD13	2.04	0.40
6:E:2:THR:OG1	6:E:3:PHE:N	2.52	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	106/121 (88%)	97 (92%)	8 (8%)	1 (1%)	14	47
1	B	106/121 (88%)	98 (92%)	6 (6%)	2 (2%)	6	33
1	M	106/121 (88%)	98 (92%)	7 (7%)	1 (1%)	14	47
3	I	293/299 (98%)	240 (82%)	51 (17%)	2 (1%)	18	51
4	J	737/758 (97%)	608 (82%)	114 (16%)	15 (2%)	6	32
6	C	210/220 (96%)	171 (81%)	35 (17%)	4 (2%)	6	33
6	E	210/220 (96%)	173 (82%)	34 (16%)	3 (1%)	9	38

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
6	N	210/220 (96%)	173 (82%)	34 (16%)	3 (1%)	9	38
6	O	210/220 (96%)	171 (81%)	36 (17%)	3 (1%)	9	38
6	P	210/220 (96%)	173 (82%)	34 (16%)	3 (1%)	9	38
All	All	2398/2520 (95%)	2002 (84%)	359 (15%)	37 (2%)	11	37

All (37) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
4	J	668	ILE
3	I	279	PRO
4	J	653	PHE
4	J	662	ASP
4	J	677	GLU
6	C	75	ASP
6	N	129	ILE
6	O	27	ALA
6	O	129	ILE
6	P	131	ALA
3	I	140	GLY
4	J	53	ARG
4	J	59	TYR
4	J	228	LEU
4	J	417	CYS
4	J	502	ASN
4	J	695	SER
6	C	76	SER
6	C	129	ILE
6	E	129	ILE
6	E	131	ALA
6	N	131	ALA
1	B	95	LYS
4	J	379	ARG
6	N	204	PHE
1	A	95	LYS
4	J	157	LYS
4	J	392	GLY
4	J	638	GLU
1	M	95	LYS
6	C	34	SER
6	E	27	ALA
6	P	128	ARG

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Mol	Chain	Res	Type
1	B	118	GLY
6	O	113	GLY
4	J	610	GLY
6	P	34	SER

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	96/108 (89%)	96 (100%)	0	100	100
1	B	96/108 (89%)	96 (100%)	0	100	100
1	M	96/108 (89%)	96 (100%)	0	100	100
3	I	259/263 (98%)	258 (100%)	1 (0%)	84	81
4	J	633/649 (98%)	628 (99%)	5 (1%)	73	77
6	C	180/187 (96%)	180 (100%)	0	100	100
6	E	180/187 (96%)	180 (100%)	0	100	100
6	N	180/187 (96%)	180 (100%)	0	100	100
6	O	180/187 (96%)	180 (100%)	0	100	100
6	P	180/187 (96%)	180 (100%)	0	100	100
All	All	2080/2171 (96%)	2074 (100%)	6 (0%)	84	83

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	I	57	LYS
4	J	173	LEU
4	J	176	LEU
4	J	665	LEU
4	J	689	LEU
4	J	708	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (52) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	27	ASN
1	B	27	ASN
1	B	34	GLN
3	I	12	ASN
3	I	135	GLN
3	I	137	HIS
3	I	207	ASN
3	I	255	GLN
4	J	68	HIS
4	J	106	ASN
4	J	129	ASN
4	J	169	GLN
4	J	189	ASN
4	J	292	ASN
4	J	356	ASN
4	J	403	HIS
4	J	414	GLN
4	J	489	GLN
4	J	502	ASN
4	J	534	ASN
4	J	536	GLN
4	J	559	ASN
4	J	560	GLN
4	J	602	ASN
4	J	667	GLN
4	J	700	ASN
4	J	725	GLN
4	J	737	ASN
1	M	27	ASN
1	M	34	GLN
6	C	11	GLN
6	C	85	ASN
6	E	11	GLN
6	E	85	ASN
6	E	159	ASN
6	E	206	ASN
6	E	213	ASN
6	N	11	GLN
6	N	85	ASN
6	N	124	ASN
6	N	159	ASN
6	N	213	ASN
6	O	11	GLN

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Mol	Chain	Res	Type
6	O	85	ASN
6	O	105	ASN
6	O	206	ASN
6	O	213	ASN
6	P	85	ASN
6	P	105	ASN
6	P	159	ASN
6	P	206	ASN
6	P	213	ASN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	H	39/72 (54%)	8 (20%)	1 (2%)
5	U	27/49 (55%)	12 (44%)	1 (3%)
All	All	66/121 (54%)	20 (30%)	2 (3%)

All (20) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	H	2	C
2	H	13	G
2	H	15	A
2	H	21	U
2	H	27	C
2	H	29	G
2	H	33	A
2	H	39	U
5	U	3	G
5	U	4	U
5	U	5	G
5	U	6	A
5	U	12	A
5	U	15	A
5	U	18	C
5	U	21	U
5	U	23	A
5	U	24	C
5	U	26	A
5	U	27	A

All (2) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
2	H	12	C
5	U	25	G

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

1 ligand is modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
7	ATP	J	801	-	32,33,33	1.42	4 (12%)	48,52,52	1.76	9 (18%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
7	ATP	J	801	-	-	0/22/38/38	0/3/3/3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	J	801	ATP	C5-C4	4.74	1.47	1.39
7	J	801	ATP	C5-C6	2.75	1.48	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	J	801	ATP	C8-N7	2.37	1.36	1.31
7	J	801	ATP	C5-N7	-2.25	1.35	1.39

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	J	801	ATP	C5-C4-N3	-5.87	118.63	126.72
7	J	801	ATP	N3-C4-N9	4.64	135.06	127.17
7	J	801	ATP	C2-N3-C4	3.70	120.88	111.83
7	J	801	ATP	C4-C5-N7	-3.49	106.59	110.58
7	J	801	ATP	N3-C2-N1	-3.21	123.73	128.58
7	J	801	ATP	C4-N9-C8	2.64	108.51	105.74
7	J	801	ATP	C5-N7-C8	2.56	107.48	103.45
7	J	801	ATP	C3'-C2'-C1'	2.33	105.87	101.46
7	J	801	ATP	C6-C5-N7	2.07	136.07	132.09

There are no chirality outliers.

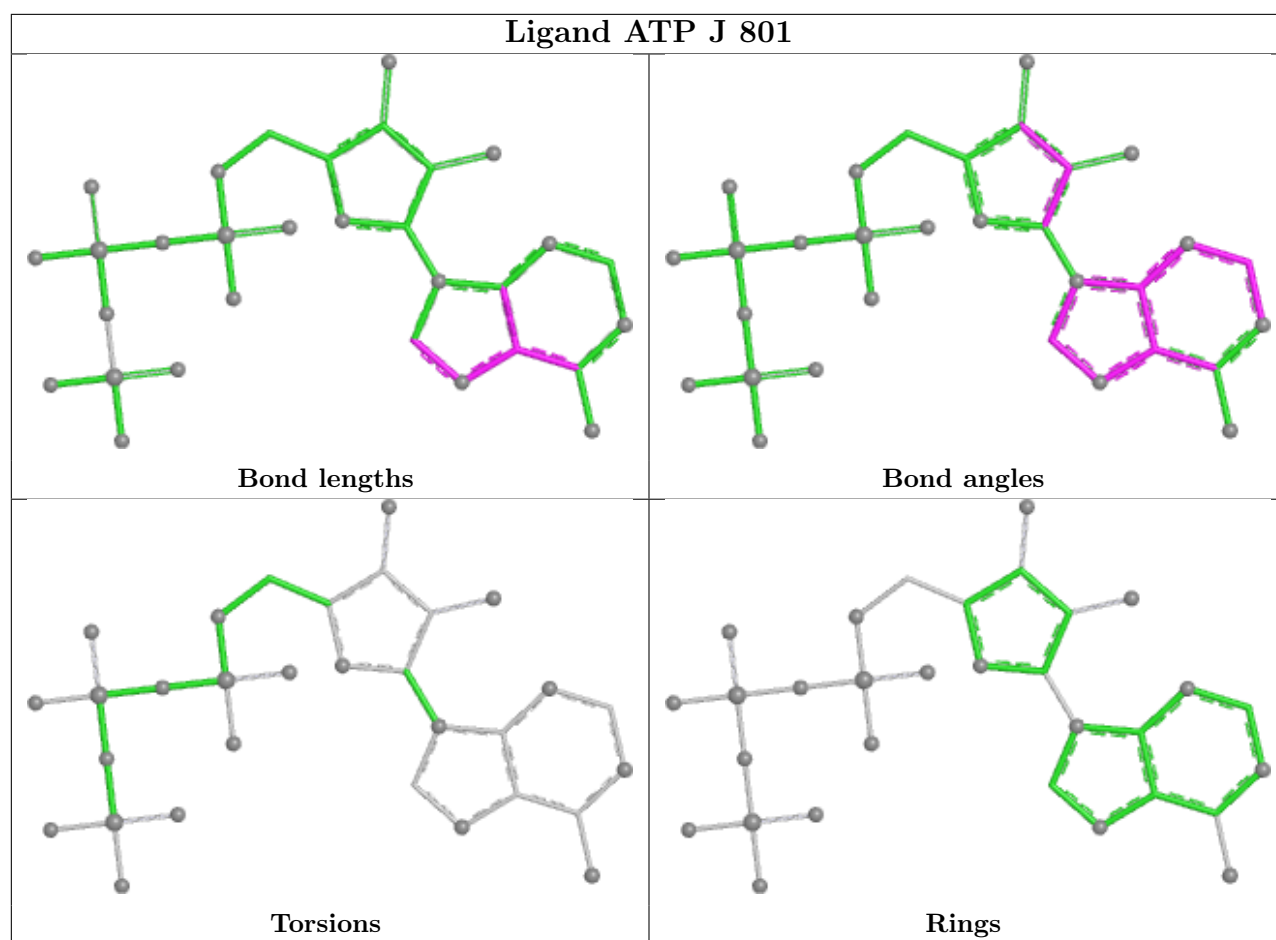
There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	J	801	ATP	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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.

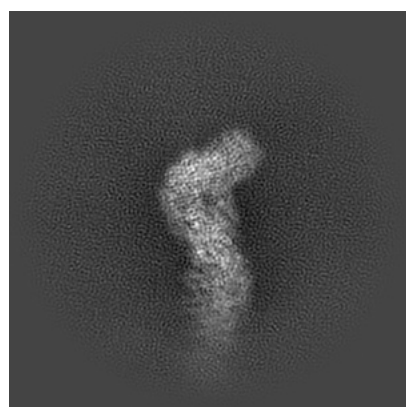
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-0516. These allow visual inspection of the internal detail of the map and identification of artifacts.

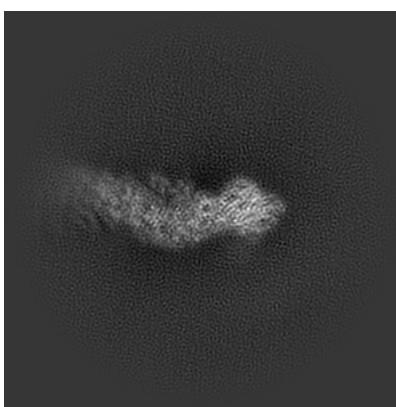
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

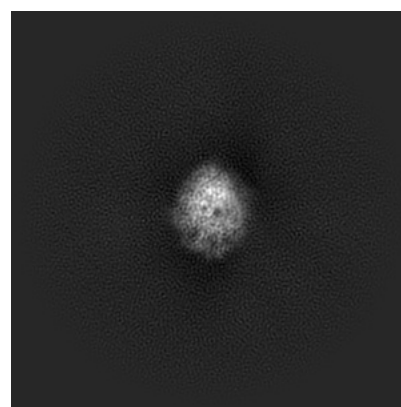
6.1.1 Primary map



X



Y

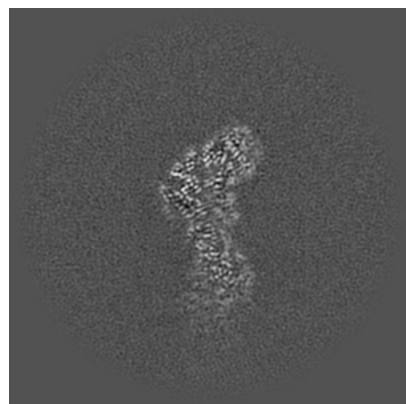


Z

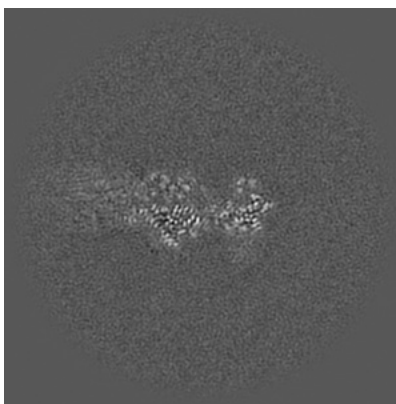
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

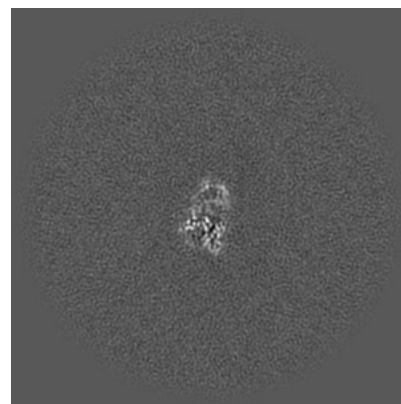
6.2.1 Primary map



X Index: 192



Y Index: 192

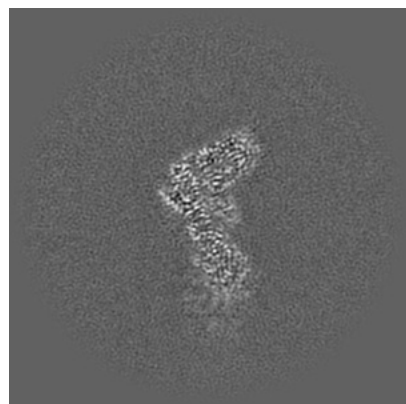


Z Index: 192

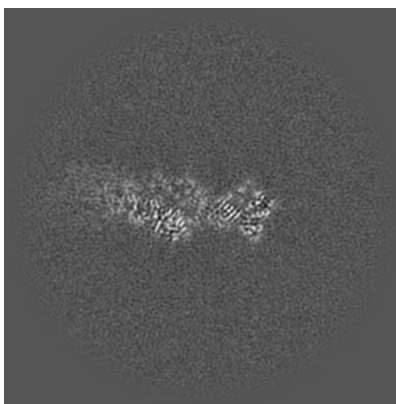
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

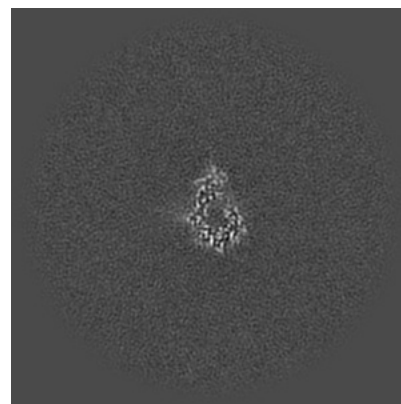
6.3.1 Primary map



X Index: 188



Y Index: 199

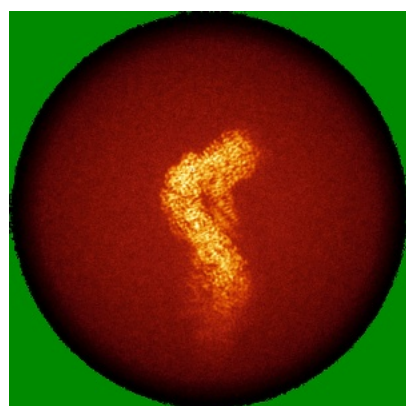


Z Index: 224

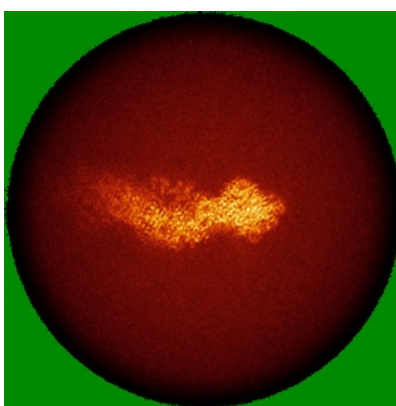
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) [i](#)

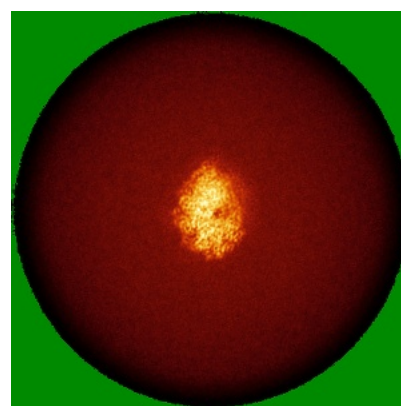
6.4.1 Primary map



X



Y

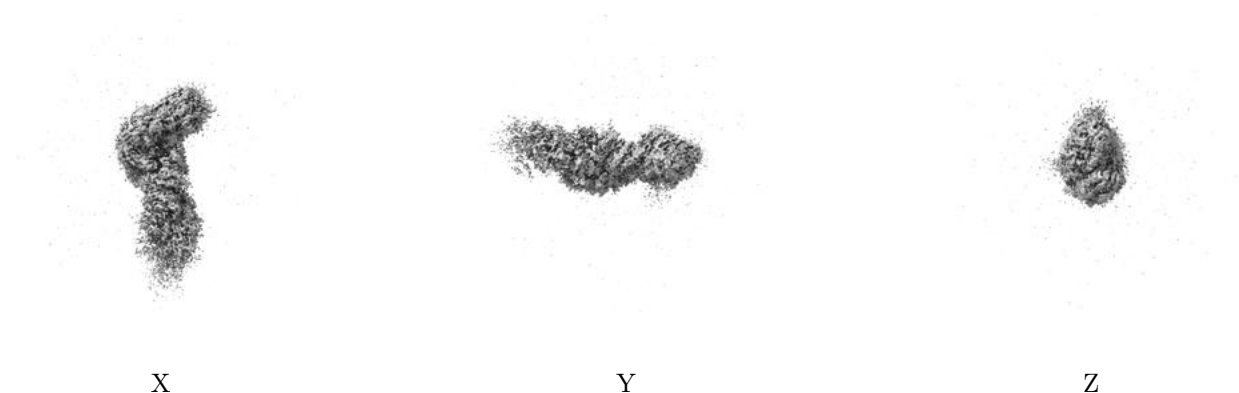


Z

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



The images above show the 3D surface view of the map at the recommended contour level 0.5. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

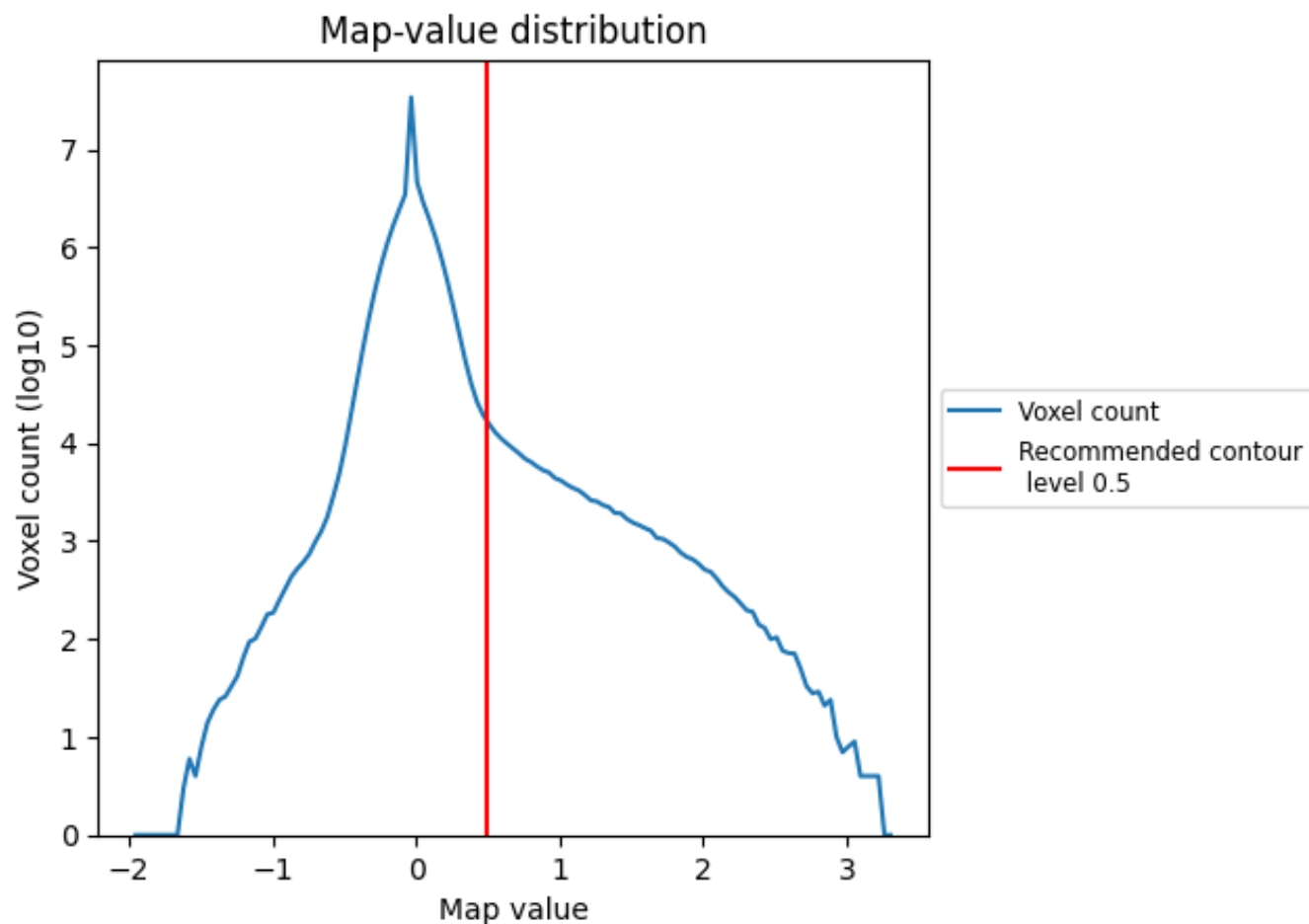
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

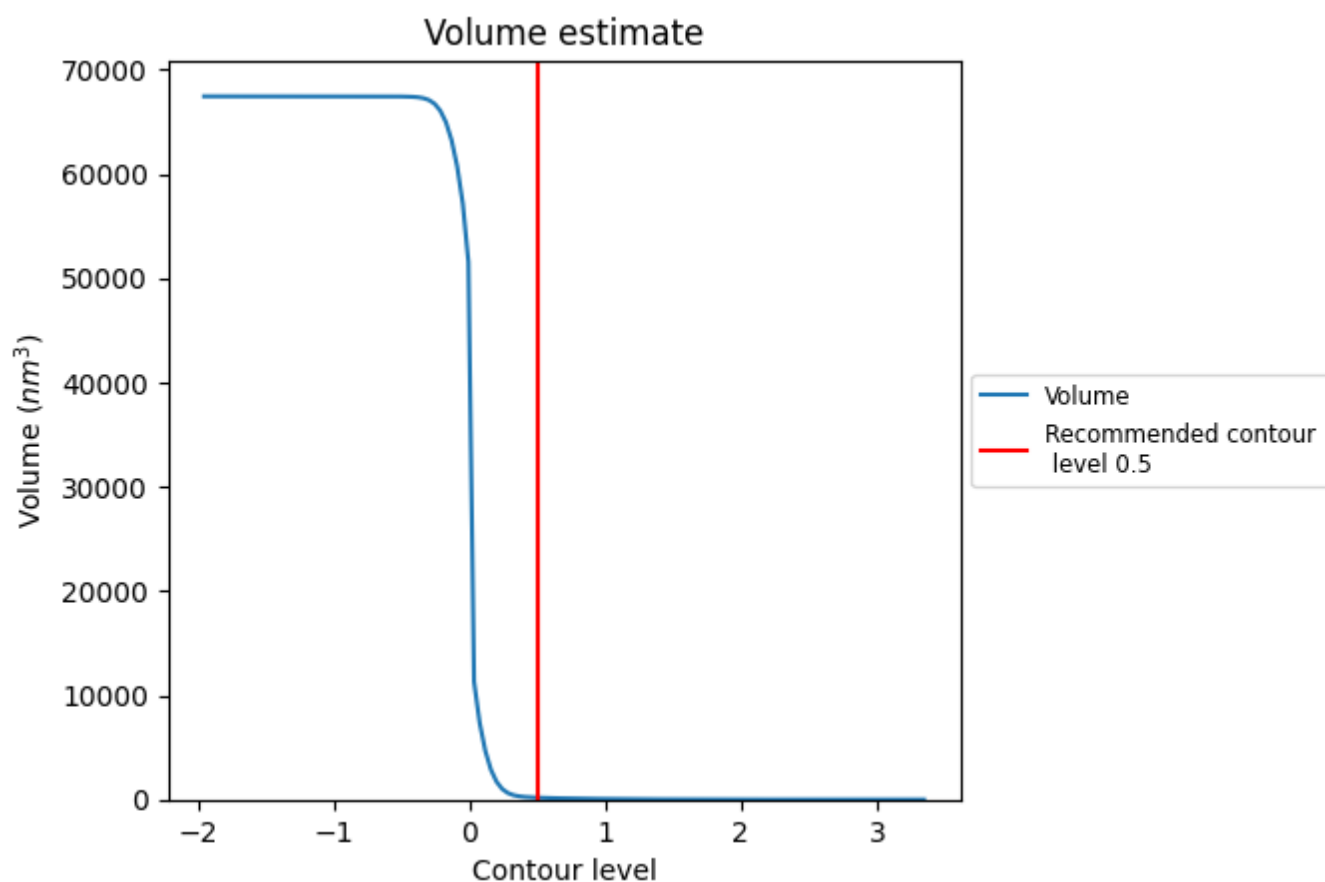
This section contains the results of statistical analysis of the map.

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.

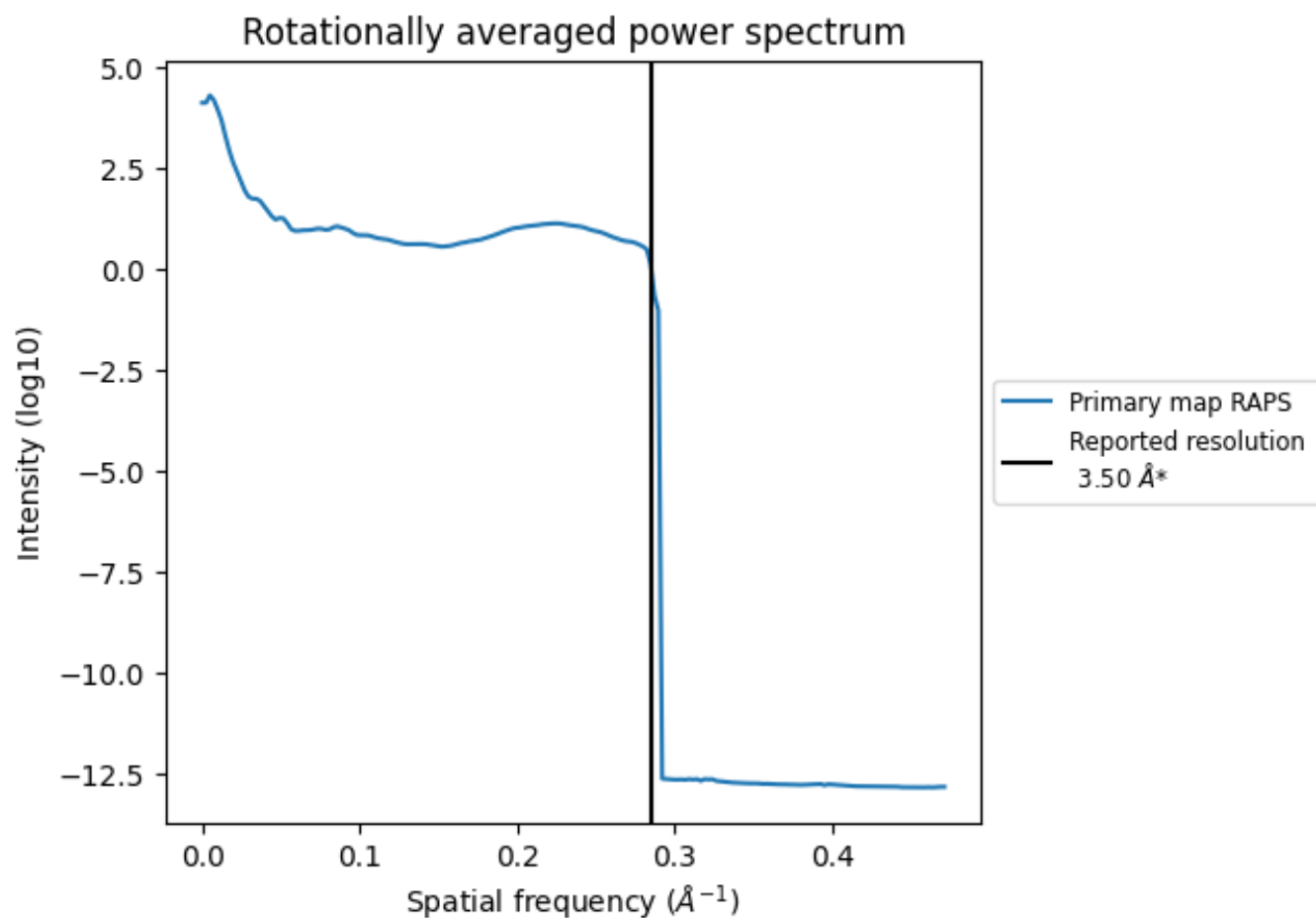
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 183 nm³; this corresponds to an approximate mass of 165 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 0.286 Å⁻¹

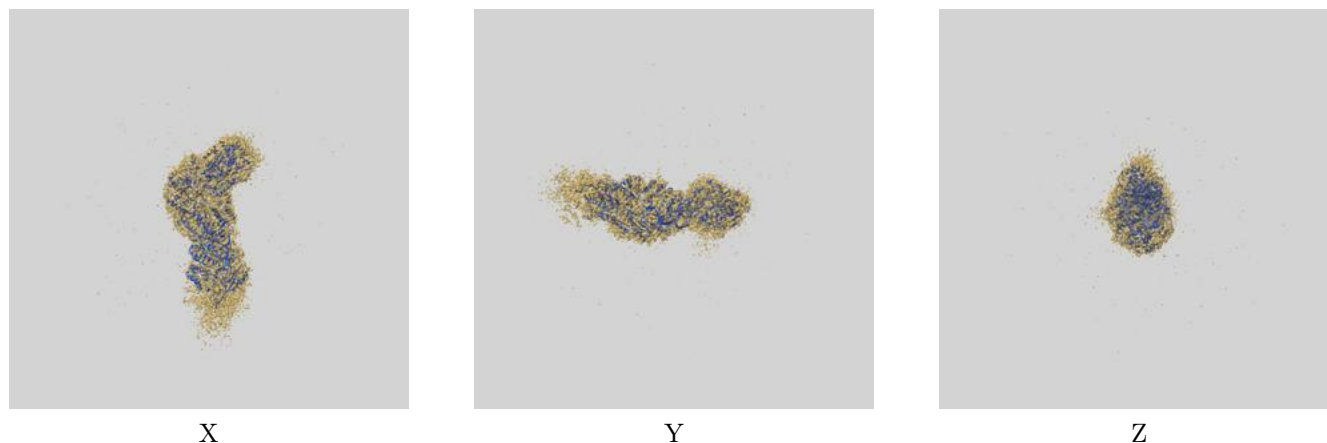
8 Fourier-Shell correlation ⓘ

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

9 Map-model fit [i](#)

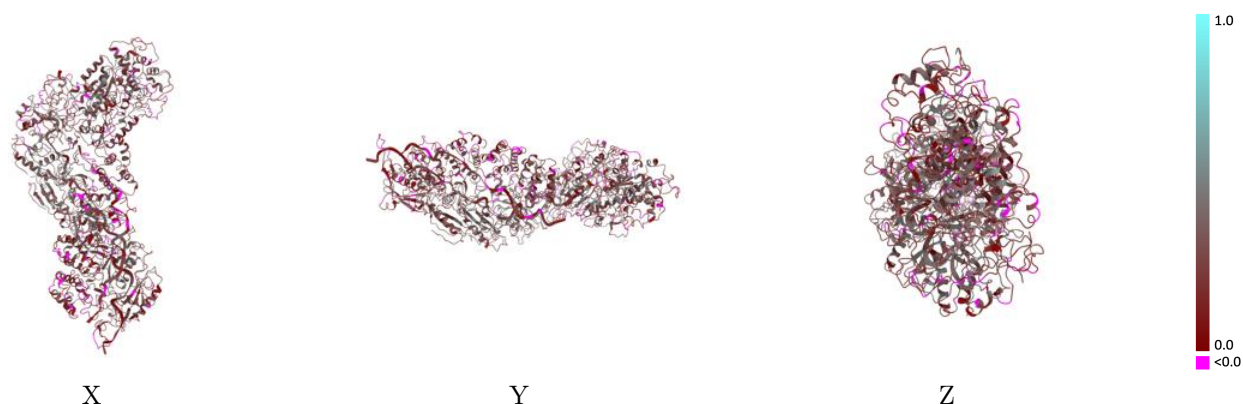
This section contains information regarding the fit between EMDB map EMD-0516 and PDB model 6NUD. Per-residue inclusion information can be found in [section 3](#) on [page 6](#).

9.1 Map-model overlay [i](#)



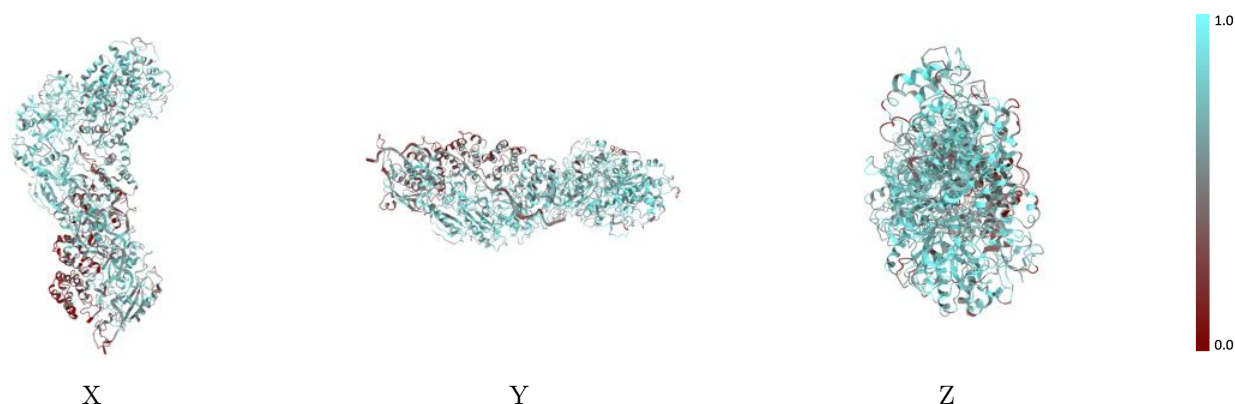
The images above show the 3D surface view of the map at the recommended contour level 0.5 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



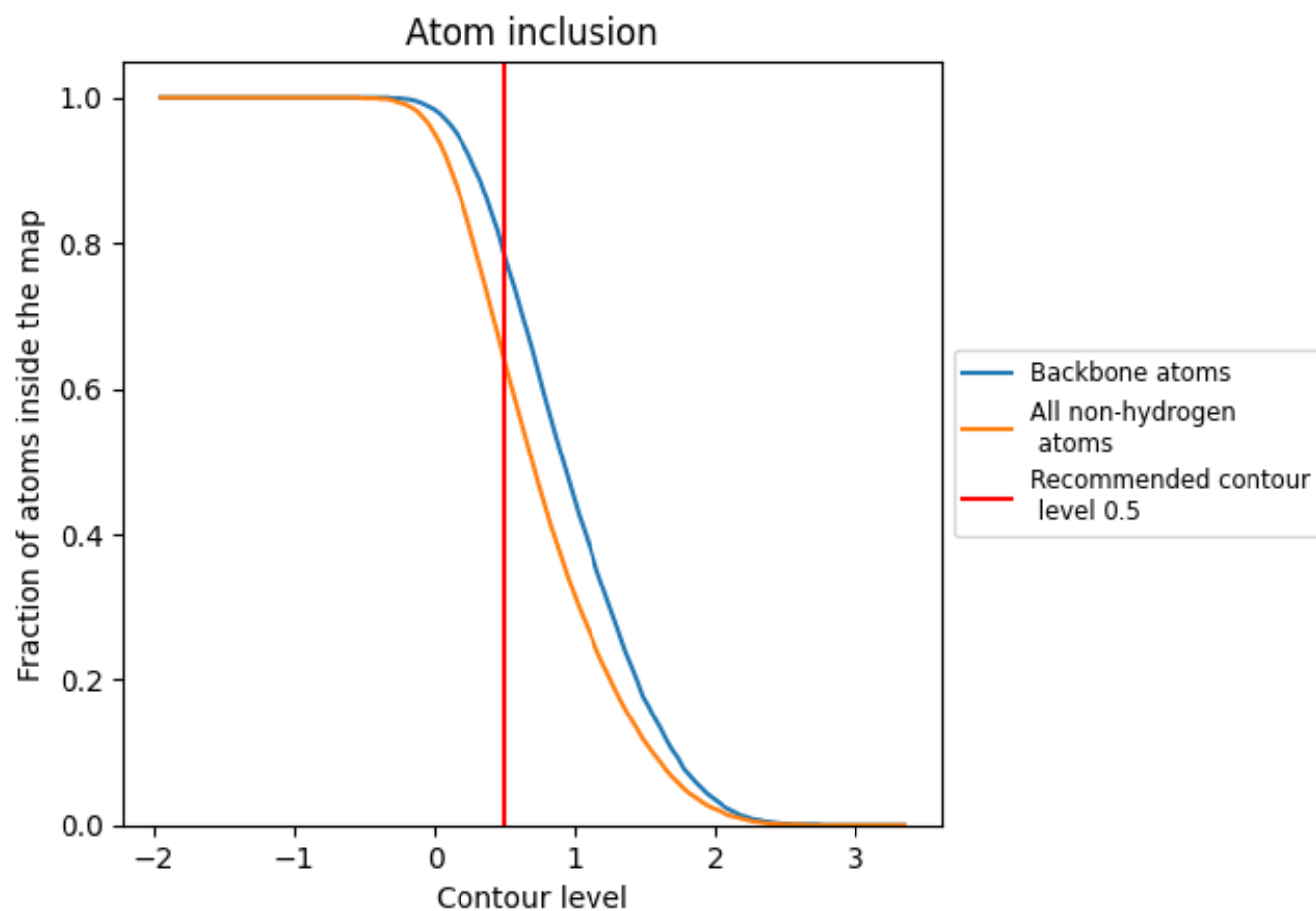
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.5).

9.4 Atom inclusion [i](#)



At the recommended contour level, 78% of all backbone atoms, 64% 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 (0.5) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.6360	 0.2540
A	 0.4220	 0.1660
B	 0.2200	 0.1300
C	 0.5310	 0.2180
E	 0.7170	 0.3210
H	 0.7600	 0.2750
I	 0.7010	 0.2500
J	 0.6890	 0.2530
M	 0.3220	 0.1590
N	 0.7160	 0.3190
O	 0.7430	 0.3240
P	 0.6880	 0.2940
U	 0.4620	 0.1250

