



Full wwPDB EM Validation Report ⓘ

Nov 7, 2022 – 05:46 AM EST

PDB ID : 6NUE
EMDB ID : EMD-0519
Title : Small conformation of apo CRISPR_Csm complex
Authors : Zhang, K.; Pintilie, G.; Li, S.; Zhu, Y.; Chiu, W.; Huang, Z.
Deposited on : 2019-01-31
Resolution : 3.30 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
Mogul : 1.8.5 (274361), CSD as541be (2020)
MolProbity : 4.02b-467
buster-report : 1.1.7 (2018)
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

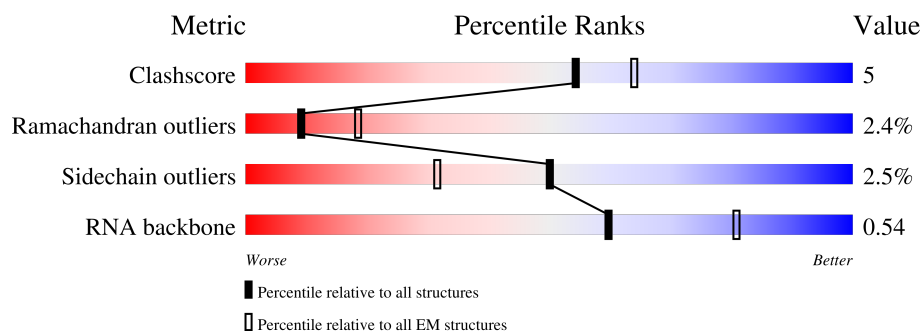
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.30 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
RNA backbone	4643	859

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	J	758	<div> <div>20%</div> <div>80%</div> <div>16%</div> <div>.</div> </div>
2	A	121	<div> <div>30%</div> <div>79%</div> <div>10%</div> <div>.</div> <div>11%</div> </div>
2	B	121	<div> <div>36%</div> <div>79%</div> <div>10%</div> <div>.</div> <div>11%</div> </div>
2	M	121	<div> <div>31%</div> <div>80%</div> <div>8%</div> <div>.</div> <div>11%</div> </div>
3	C	220	<div> <div>59%</div> <div>77%</div> <div>18%</div> <div>.</div> <div>.</div> </div>
3	E	220	<div> <div>15%</div> <div>73%</div> <div>21%</div> <div>.</div> <div>.</div> </div>
3	N	220	<div> <div>19%</div> <div>73%</div> <div>22%</div> <div>.</div> <div>.</div> </div>

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Mol	Chain	Length	Quality of chain
3	O	220	<div><div></div><div>17%</div><div>72%</div><div>23%</div><div></div><div></div></div>
3	P	220	<div><div></div><div>21%</div><div>73%</div><div>22%</div><div></div><div></div></div>
4	H	72	<div><div></div><div>6%</div><div>39%</div><div>12%</div><div></div><div>44%</div></div>
5	I	299	<div><div></div><div>15%</div><div>84%</div><div>14%</div><div></div><div></div></div>

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 20175 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 single-strand-specific deoxyribonuclease Cas10/Csm1 (subtype III-A).

Mol	Chain	Residues	Atoms					AltConf	Trace
1	J	733	Total	C	N	O	S	0	0
			5928	3796	990	1127	15		

- Molecule 2 is a protein called CRISPR system Cms protein Csm2.

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

- Molecule 3 is a protein called CRISPR type III-associated RAMP protein Csm3.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	212	Total	C	N	O	S	0	0
			1670	1058	285	325	2		
3	E	212	Total	C	N	O	S	0	0
			1670	1058	285	325	2		
3	N	212	Total	C	N	O	S	0	0
			1670	1058	285	325	2		
3	O	212	Total	C	N	O	S	0	0
			1670	1058	285	325	2		
3	P	212	Total	C	N	O	S	0	0
			1670	1058	285	325	2		

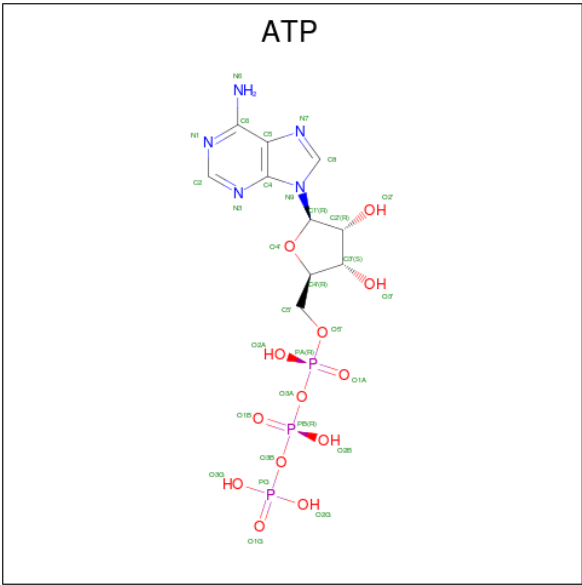
- Molecule 4 is a RNA chain called crRNA.

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

- Molecule 5 is a protein called CRISPR type III-associated RAMP protein Csm4.

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

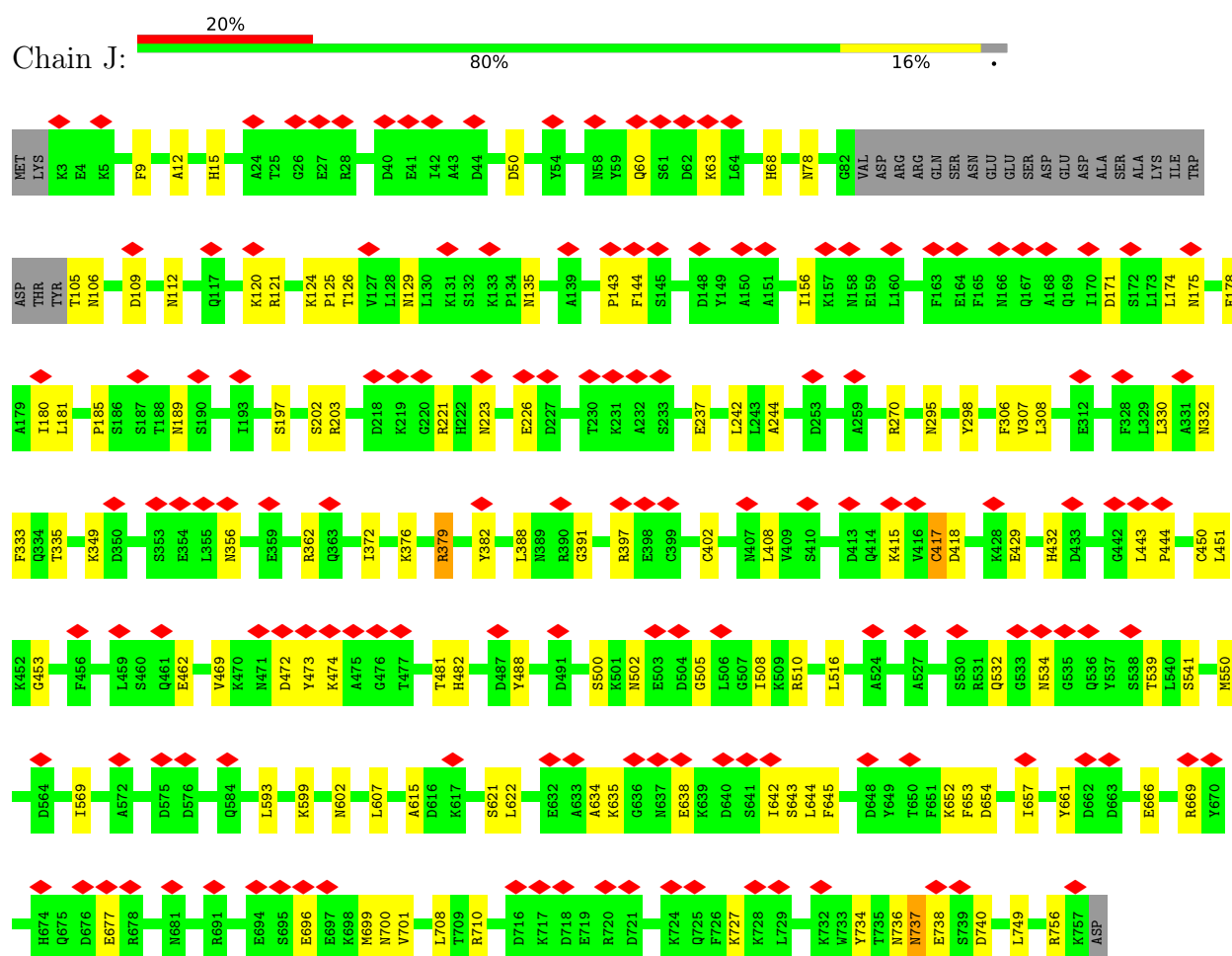
- Molecule 6 is ADENOSINE-5'-TRIPHOSPHATE (three-letter code: ATP) (formula: C₁₀H₁₆N₅O₁₃P₃).



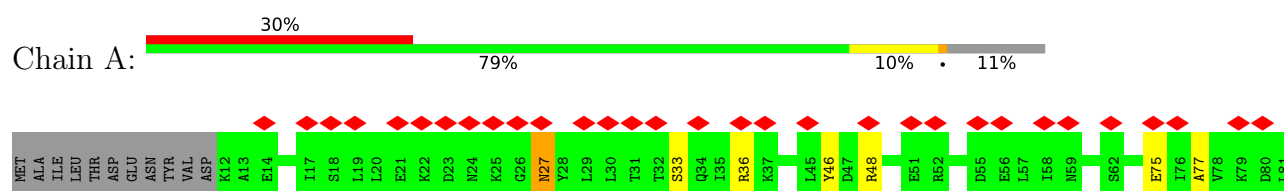
3 Residue-property plots [i](#)

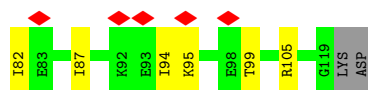
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion 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: CRISPR system single-strand-specific deoxyribonuclease Cas10/Csm1 (subtype III-A)

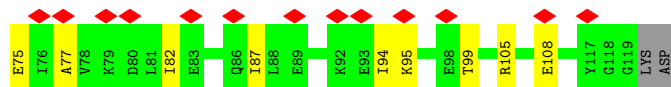
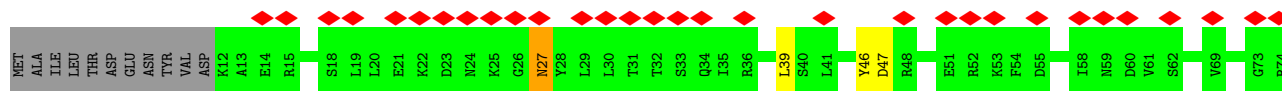
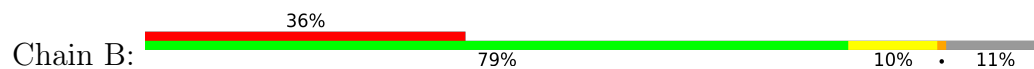


- Molecule 2: CRISPR system Cms protein Csm2

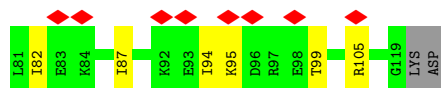
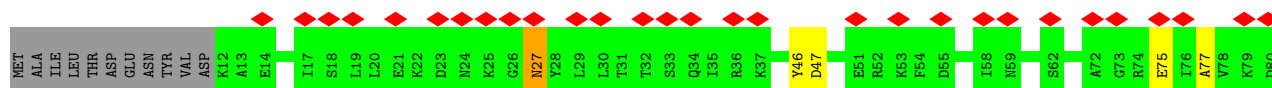
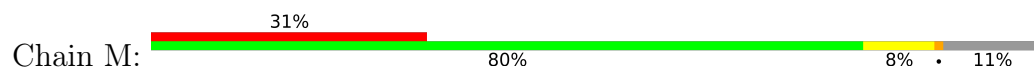




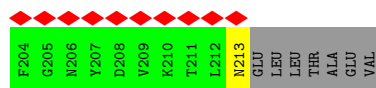
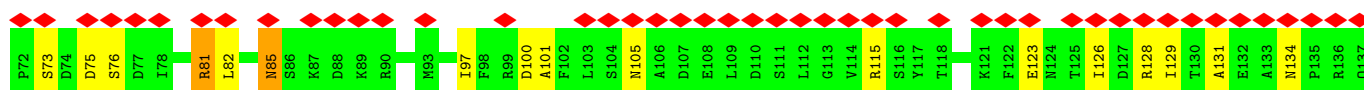
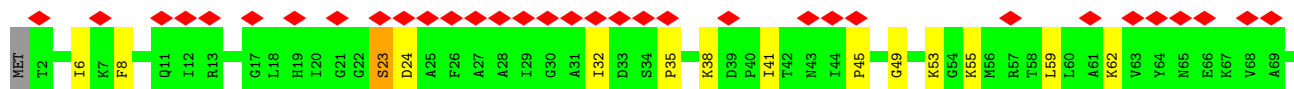
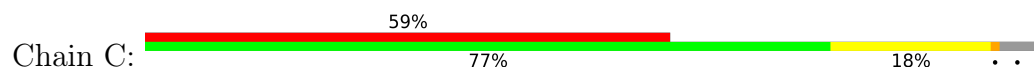
- Molecule 2: CRISPR system Cms protein Csm2



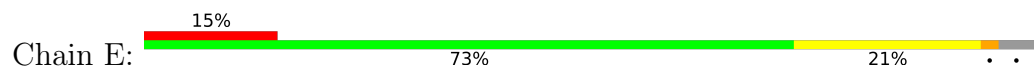
- Molecule 2: CRISPR system Cms protein Csm2

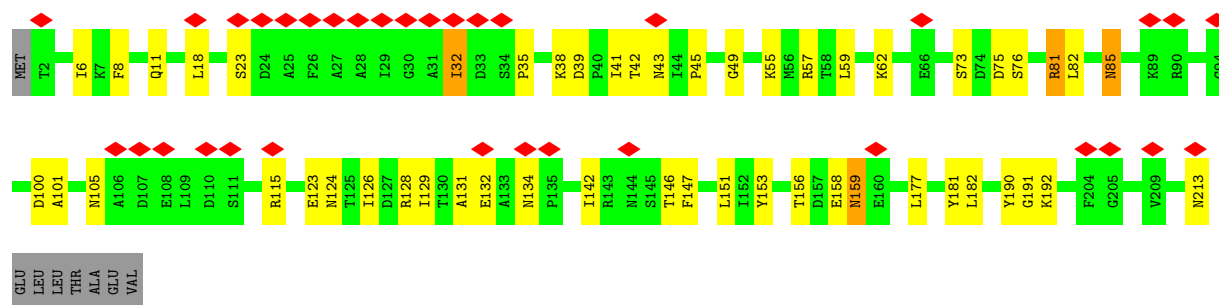


- Molecule 3: CRISPR type III-associated RAMP protein Csm3

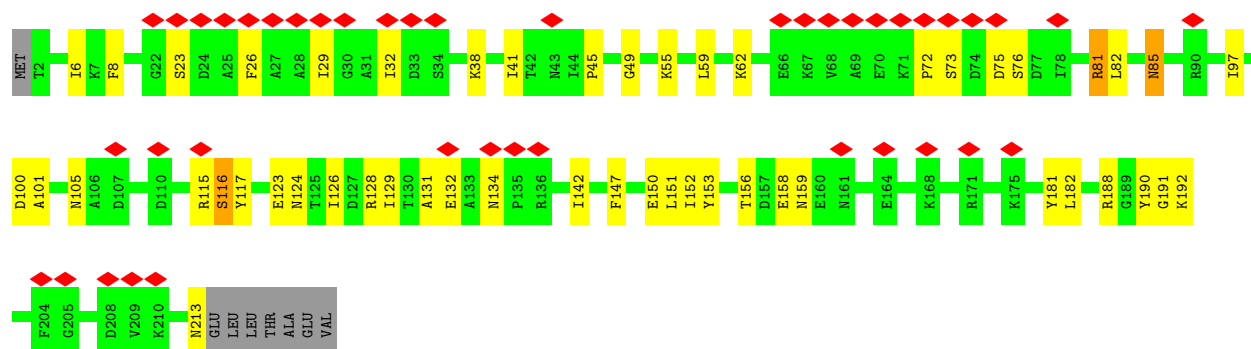


- Molecule 3: CRISPR type III-associated RAMP protein Csm3

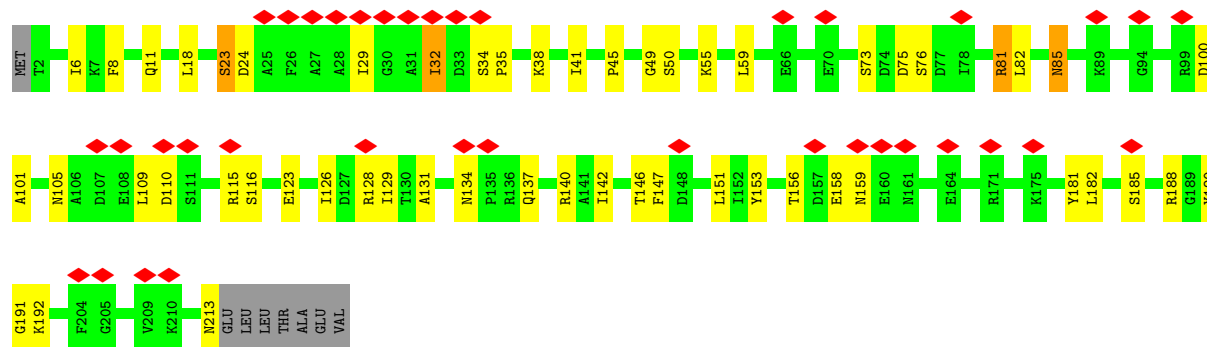




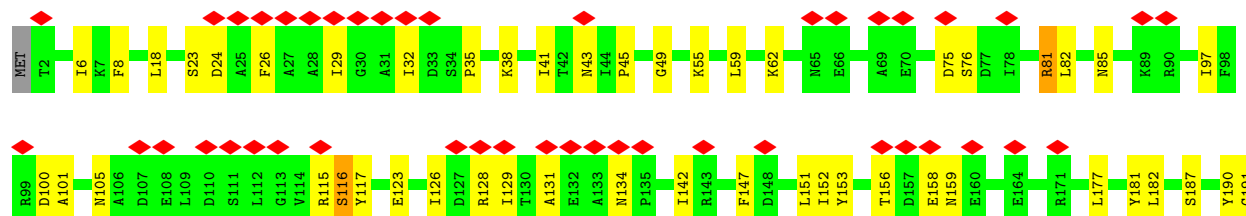
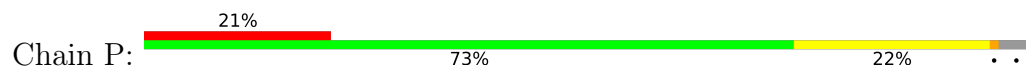
• Molecule 3: CRISPR type III-associated RAMP protein Csm3

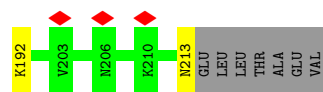


• Molecule 3: CRISPR type III-associated RAMP protein Csm3

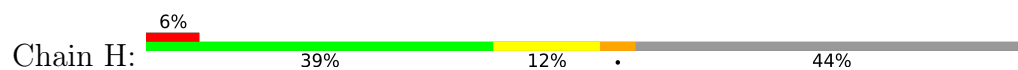


• Molecule 3: CRISPR type III-associated RAMP protein Csm3

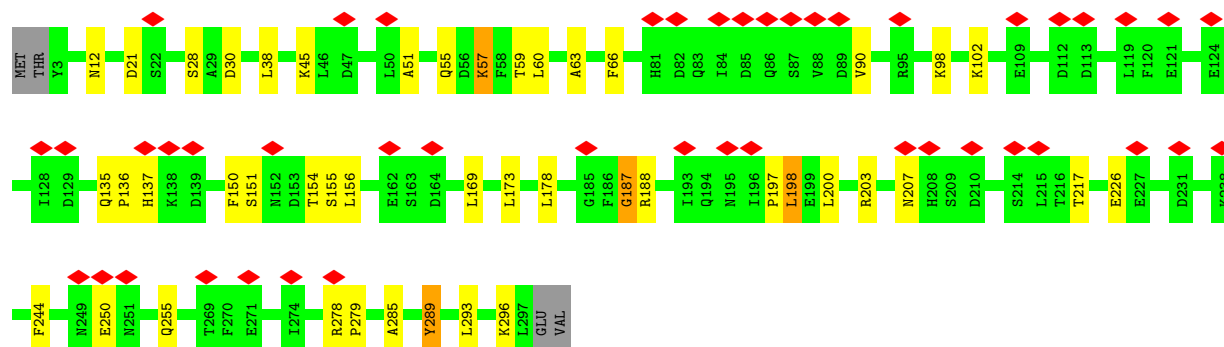
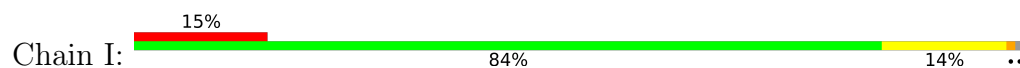




- Molecule 4: crRNA



- Molecule 5: CRISPR type III-associated RAMP protein Csm4



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	81540	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	5.120	Depositor
Minimum map value	-2.725	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.135	Depositor
Recommended contour level	0.8	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	J	0.24	0/6051	0.38	0/8149
2	A	0.23	0/905	0.34	0/1209
2	B	0.23	0/905	0.34	0/1209
2	M	0.23	0/905	0.34	0/1209
3	C	0.24	0/1695	0.40	0/2279
3	E	0.23	0/1695	0.40	0/2279
3	N	0.24	0/1695	0.41	0/2279
3	O	0.24	0/1695	0.40	0/2279
3	P	0.23	0/1695	0.40	0/2279
4	H	0.14	0/926	0.82	0/1437
5	I	0.24	0/2404	0.41	0/3247
All	All	0.23	0/20571	0.42	0/27855

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	J	5928	0	5801	65	0
2	A	894	0	923	7	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	894	0	923	7	0
2	M	894	0	923	5	0
3	C	1670	0	1684	22	0
3	E	1670	0	1684	28	0
3	N	1670	0	1684	29	0
3	O	1670	0	1684	27	0
3	P	1670	0	1684	26	0
4	H	832	0	425	5	0
5	I	2352	0	2298	28	0
6	J	31	0	12	2	0
All	All	20175	0	19725	216	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 (216) 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)
3:N:182:LEU:H	3:N:191:GLY:HA3	1.61	0.64
3:N:101:ALA:HB1	3:N:147:PHE:HB3	1.80	0.64
3:O:101:ALA:HB1	3:O:147:PHE:HB3	1.80	0.62
1:J:635:LYS:NZ	6:J:801:ATP:O1G	2.32	0.62
3:E:6:ILE:HB	3:E:153:TYR:HB3	1.82	0.61
3:E:101:ALA:HB1	3:E:147:PHE:HB3	1.82	0.61
3:C:6:ILE:HB	3:C:153:TYR:HB3	1.81	0.61
3:P:6:ILE:HB	3:P:153:TYR:HB3	1.82	0.61
3:P:101:ALA:HB1	3:P:147:PHE:HB3	1.81	0.61
3:O:6:ILE:HB	3:O:153:TYR:HB3	1.84	0.60
1:J:242:LEU:HB3	1:J:308:LEU:HB2	1.83	0.60
1:J:295:ASN:HD22	1:J:307:VAL:HG13	1.67	0.60
3:P:182:LEU:H	3:P:191:GLY:HA3	1.68	0.59
2:B:82:ILE:HG23	2:B:87:ILE:HB	1.84	0.59
3:E:126:ILE:HG22	3:E:128:ARG:H	1.68	0.59
1:J:109:ASP:HB3	1:J:112:ASN:HB2	1.83	0.59
1:J:666:GLU:HG3	1:J:669:ARG:HH21	1.68	0.58
5:I:59:THR:HG21	5:I:293:LEU:HD21	1.85	0.58
2:B:39:LEU:HB3	3:P:29:ILE:HG13	1.84	0.58
3:O:18:LEU:HA	3:O:190:TYR:HB3	1.86	0.58
3:N:8:PHE:HB2	3:N:151:LEU:HB3	1.85	0.58
5:I:12:ASN:ND2	3:N:100:ASP:OD2	2.38	0.57
1:J:756:ARG:NH2	3:N:32:ILE:O	2.37	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:8:PHE:HB2	3:C:151:LEU:HB3	1.87	0.57
1:J:657:ILE:HA	1:J:661:TYR:HB3	1.86	0.56
3:C:101:ALA:HB1	3:C:147:PHE:HB3	1.87	0.56
3:O:126:ILE:HG22	3:O:128:ARG:H	1.71	0.56
3:P:126:ILE:HG22	3:P:128:ARG:H	1.71	0.56
2:M:82:ILE:HG23	2:M:87:ILE:HB	1.88	0.56
2:A:82:ILE:HG23	2:A:87:ILE:HB	1.88	0.56
3:C:59:LEU:HA	3:C:62:LYS:HE3	1.88	0.55
1:J:105:THR:N	1:J:143:PRO:O	2.38	0.55
3:E:18:LEU:HA	3:E:190:TYR:HB3	1.88	0.55
3:E:177:LEU:HD11	3:E:182:LEU:HD13	1.89	0.55
4:H:15:A:O2'	3:N:124:ASN:ND2	2.40	0.55
5:I:45:LYS:HD2	5:I:169:LEU:HD21	1.89	0.55
3:N:59:LEU:HA	3:N:62:LYS:HE3	1.89	0.55
1:J:736:ASN:ND2	1:J:740:ASP:OD2	2.39	0.55
3:P:38:LYS:HG2	3:P:45:PRO:HA	1.90	0.54
1:J:156:ILE:HD11	1:J:181:LEU:HD11	1.90	0.54
2:A:75:GLU:HG2	2:A:77:ALA:H	1.73	0.54
3:N:156:THR:O	3:N:158:GLU:N	2.40	0.54
3:N:6:ILE:HB	3:N:153:TYR:HB3	1.90	0.54
3:E:59:LEU:HA	3:E:62:LYS:HE3	1.89	0.54
5:I:28:SER:OG	5:I:30:ASP:OD1	2.21	0.53
3:N:81:ARG:HG2	3:N:82:LEU:HD12	1.91	0.53
2:A:36:ARG:NH1	3:O:137:GLN:OE1	2.39	0.53
5:I:59:THR:OG1	5:I:289:TYR:OH	2.26	0.53
3:O:182:LEU:H	3:O:191:GLY:HA3	1.73	0.53
2:M:75:GLU:HG2	2:M:77:ALA:H	1.74	0.53
3:P:181:TYR:HA	3:P:192:LYS:H	1.74	0.53
1:J:112:ASN:HD21	1:J:120:LYS:HD3	1.74	0.53
3:P:59:LEU:HA	3:P:62:LYS:HE3	1.91	0.53
3:P:49:GLY:N	3:P:100:ASP:OD1	2.40	0.52
3:O:11:GLN:NE2	3:O:146:THR:OG1	2.42	0.52
3:P:8:PHE:HB2	3:P:151:LEU:HB3	1.91	0.52
1:J:749:LEU:HD13	3:N:29:ILE:HD12	1.90	0.52
3:E:38:LYS:HG2	3:E:45:PRO:HA	1.91	0.52
2:B:46:TYR:OH	2:B:105:ARG:NH2	2.43	0.52
3:C:182:LEU:H	3:C:191:GLY:HA3	1.74	0.52
1:J:124:LYS:HD2	1:J:125:PRO:HD2	1.90	0.52
2:B:47:ASP:OD2	3:P:43:ASN:ND2	2.42	0.52
2:A:48:ARG:NH2	3:O:110:ASP:OD1	2.43	0.52
1:J:126:THR:OG1	1:J:135:ASN:ND2	2.43	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:N:188:ARG:HG2	3:O:50:SER:HA	1.90	0.52
3:E:73:SER:O	3:E:85:ASN:ND2	2.43	0.52
2:A:94:ILE:HG22	2:A:99:THR:HG22	1.91	0.51
2:M:46:TYR:OH	2:M:105:ARG:NH2	2.43	0.51
1:J:510:ARG:HD2	1:J:615:ALA:HA	1.91	0.51
2:B:94:ILE:HG22	2:B:99:THR:HG22	1.92	0.51
2:B:75:GLU:HG2	2:B:77:ALA:H	1.74	0.51
3:N:124:ASN:HD22	3:N:132:GLU:HG3	1.75	0.51
3:E:156:THR:O	3:E:158:GLU:N	2.41	0.51
3:E:181:TYR:HA	3:E:192:LYS:H	1.76	0.51
3:E:11:GLN:NE2	3:E:146:THR:OG1	2.44	0.51
3:O:156:THR:O	3:O:158:GLU:N	2.41	0.51
3:C:81:ARG:HG2	3:C:82:LEU:HD12	1.94	0.50
5:I:278:ARG:HH21	5:I:285:ALA:HB2	1.77	0.50
2:M:94:ILE:HG22	2:M:99:THR:HG22	1.92	0.50
3:N:73:SER:O	3:N:85:ASN:ND2	2.44	0.50
1:J:382:TYR:HB2	5:I:226:GLU:HB2	1.93	0.50
3:C:126:ILE:HG22	3:C:128:ARG:H	1.75	0.50
3:O:81:ARG:HG2	3:O:82:LEU:HD12	1.94	0.50
1:J:450:CYS:SG	1:J:451:LEU:N	2.83	0.50
1:J:516:LEU:HD13	1:J:593:LEU:HD23	1.94	0.50
5:I:98:LYS:HG2	5:I:102:LYS:HE3	1.93	0.50
3:P:97:ILE:HB	3:P:152:ILE:HB	1.94	0.50
3:E:41:ILE:HD12	3:O:109:LEU:HD11	1.93	0.50
3:E:49:GLY:N	3:E:100:ASP:OD1	2.39	0.50
5:I:188:ARG:HH12	3:N:150:GLU:HB2	1.77	0.50
1:J:443:LEU:HD22	1:J:444:PRO:HD2	1.94	0.50
3:E:8:PHE:HB2	3:E:151:LEU:HB3	1.93	0.50
3:O:181:TYR:HA	3:O:192:LYS:H	1.77	0.50
3:C:156:THR:O	3:C:158:GLU:N	2.41	0.49
3:E:55:LYS:HE3	3:E:59:LEU:HD11	1.94	0.49
1:J:121:ARG:NH1	1:J:599:LYS:O	2.45	0.49
1:J:244:ALA:HB3	1:J:306:PHE:HB2	1.93	0.49
1:J:408:LEU:HA	1:J:417:CYS:HA	1.95	0.49
3:N:126:ILE:HG22	3:N:128:ARG:H	1.77	0.49
3:O:8:PHE:HB2	3:O:151:LEU:HB3	1.93	0.49
2:A:46:TYR:OH	2:A:105:ARG:NH2	2.46	0.49
1:J:9:PHE:HZ	1:J:174:LEU:HD11	1.78	0.49
1:J:532:GLN:HG2	1:J:534:ASN:H	1.78	0.48
3:C:38:LYS:HG2	3:C:45:PRO:HA	1.95	0.48
3:N:116:SER:OG	3:N:117:TYR:N	2.45	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:100:ASP:OD2	3:O:140:ARG:NH2	2.44	0.48
2:A:33:SER:O	2:A:36:ARG:NE	2.42	0.48
1:J:516:LEU:HD21	1:J:607:LEU:HD11	1.95	0.48
4:H:15:A:N6	3:O:34:SER:OG	2.46	0.48
3:P:156:THR:O	3:P:158:GLU:N	2.41	0.48
3:E:182:LEU:H	3:E:191:GLY:HA3	1.78	0.47
1:J:737:ASN:HD22	1:J:738:GLU:H	1.61	0.47
1:J:408:LEU:HG	1:J:417:CYS:HA	1.95	0.47
3:E:81:ARG:HG2	3:E:82:LEU:HD12	1.95	0.47
1:J:621:SER:OG	5:I:21:ASP:OD2	2.33	0.47
3:C:55:LYS:HE3	3:C:59:LEU:HD11	1.96	0.47
1:J:185:PRO:HA	1:J:197:SER:HA	1.97	0.47
3:C:97:ILE:HB	3:C:152:ILE:HB	1.96	0.47
3:P:81:ARG:HG2	3:P:82:LEU:HD12	1.96	0.47
1:J:12:ALA:O	1:J:15:HIS:ND1	2.45	0.47
3:P:24:ASP:HB3	3:P:26:PHE:HD2	1.79	0.47
1:J:550:MET:SD	6:J:801:ATP:N6	2.88	0.47
3:E:43:ASN:ND2	2:M:47:ASP:OD2	2.48	0.47
5:I:178:LEU:H	5:I:187:GLY:HA3	1.80	0.46
3:C:49:GLY:N	3:C:100:ASP:OD1	2.43	0.46
3:E:142:ILE:HD12	3:P:41:ILE:HD11	1.97	0.46
3:E:123:GLU:OE2	3:E:190:TYR:OH	2.33	0.46
3:O:73:SER:O	3:O:85:ASN:ND2	2.49	0.46
1:J:645:PHE:HB3	3:N:26:PHE:HE1	1.79	0.46
5:I:150:PHE:HB3	5:I:154:THR:HB	1.98	0.46
5:I:173:LEU:HD11	5:I:178:LEU:HD23	1.96	0.46
3:E:41:ILE:HD11	3:O:142:ILE:HD12	1.96	0.46
3:O:38:LYS:HG2	3:O:45:PRO:HA	1.98	0.46
1:J:417:CYS:SG	1:J:418:ASP:N	2.89	0.46
3:N:142:ILE:HD12	3:O:41:ILE:HD11	1.98	0.46
5:I:203:ARG:HD3	5:I:296:LYS:HB2	1.98	0.46
1:J:488:TYR:HB3	1:J:569:ILE:HG22	1.97	0.45
3:N:123:GLU:OE2	3:N:190:TYR:OH	2.33	0.45
1:J:15:HIS:HE1	1:J:202:SER:HA	1.81	0.45
5:I:63:ALA:HB1	5:I:156:LEU:HD13	1.98	0.45
3:N:181:TYR:HA	3:N:192:LYS:H	1.80	0.45
1:J:642:ILE:HB	1:J:653:PHE:HB2	1.98	0.45
1:J:372:ILE:HG22	1:J:376:LYS:HE2	1.99	0.45
1:J:696:GLU:HG3	1:J:699:MET:HB3	1.98	0.45
3:O:55:LYS:HE3	3:O:59:LEU:HD11	1.97	0.45
1:J:178:GLU:OE1	1:J:203:ARG:NH1	2.49	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:124:ASN:HD22	3:E:132:GLU:HG3	1.82	0.45
1:J:362:ARG:HH11	5:I:90:VAL:HB	1.82	0.44
5:I:135:GLN:HE21	5:I:137:HIS:CE1	2.35	0.44
3:N:49:GLY:N	3:N:100:ASP:OD1	2.51	0.44
3:O:49:GLY:N	3:O:100:ASP:OD1	2.43	0.44
3:C:73:SER:O	3:C:85:ASN:ND2	2.51	0.44
1:J:175:ASN:HD22	1:J:482:HIS:HE1	1.65	0.44
1:J:500:SER:OG	1:J:508:ILE:O	2.30	0.44
3:E:57:ARG:NH1	4:H:19:G:OP1	2.48	0.44
1:J:429:GLU:HB3	1:J:453:GLY:HA3	1.98	0.44
1:J:710:ARG:HH12	3:O:32:ILE:HD12	1.83	0.44
3:O:123:GLU:OE2	3:O:190:TYR:OH	2.35	0.44
3:P:177:LEU:HD11	3:P:182:LEU:HD13	1.99	0.44
5:I:60:LEU:O	5:I:289:TYR:OH	2.29	0.43
3:P:18:LEU:HA	3:P:190:TYR:HB3	2.01	0.43
1:J:622:LEU:HD23	5:I:21:ASP:HB2	2.00	0.43
1:J:397:ARG:HG2	1:J:415:LYS:HD3	2.00	0.43
1:J:652:LYS:O	1:J:654:ASP:N	2.43	0.43
3:C:23:SER:OG	3:C:24:ASP:N	2.51	0.43
3:E:153:TYR:OH	3:E:159:ASN:OD1	2.35	0.43
3:P:55:LYS:HE3	3:P:59:LEU:HD11	2.00	0.43
3:C:123:GLU:OE2	3:C:190:TYR:OH	2.35	0.43
5:I:66:PHE:N	5:I:155:SER:O	2.51	0.43
3:N:97:ILE:HB	3:N:152:ILE:HB	2.01	0.43
1:J:332:ASN:OD1	1:J:333:PHE:N	2.47	0.42
1:J:539:THR:HG22	1:J:541:SER:H	1.84	0.42
4:H:12:C:O2'	4:H:13:G:OP1	2.30	0.42
1:J:708:LEU:HB3	1:J:727:LYS:HG3	2.00	0.42
5:I:137:HIS:HE1	3:N:72:PRO:HG3	1.85	0.42
5:I:38:LEU:HD11	5:I:244:PHE:HD2	1.84	0.42
1:J:330:LEU:HD11	1:J:388:LEU:HD11	2.02	0.42
1:J:736:ASN:OD1	1:J:737:ASN:N	2.52	0.42
1:J:429:GLU:HA	1:J:432:HIS:HD2	1.84	0.42
3:C:41:ILE:HD11	3:P:142:ILE:HD12	2.01	0.42
3:N:38:LYS:HG2	3:N:45:PRO:HA	2.02	0.42
3:N:55:LYS:HE3	3:N:59:LEU:HD11	2.02	0.42
3:O:185:SER:HB2	3:O:188:ARG:HH12	1.85	0.42
5:I:51:ALA:O	5:I:55:GLN:NE2	2.53	0.41
5:I:198:LEU:HD13	5:I:200:LEU:H	1.85	0.41
1:J:391:GLY:HA2	5:I:255:GLN:HE21	1.85	0.41
3:N:75:ASP:OD1	3:N:75:ASP:N	2.52	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:P:123:GLU:OE2	3:P:190:TYR:OH	2.39	0.41
1:J:237:GLU:HA	1:J:349:LYS:HD3	2.03	0.41
1:J:634:ALA:HB2	1:J:643:SER:HB2	2.03	0.41
4:H:10:U:H1'	5:I:136:PRO:HB3	2.02	0.41
3:P:116:SER:OG	3:P:117:TYR:N	2.52	0.41
2:B:108:GLU:HG2	3:P:29:ILE:HD13	2.02	0.41
3:C:75:ASP:OD2	3:P:128:ARG:NH2	2.54	0.41
1:J:50:ASP:HB3	1:J:68:HIS:HE1	1.85	0.41
1:J:171:ASP:HB2	1:J:481:THR:HG22	2.03	0.41
3:C:32:ILE:HD12	3:C:32:ILE:HA	1.96	0.41
3:C:38:LYS:HE3	3:C:45:PRO:HG3	2.02	0.41
3:C:75:ASP:OD1	3:C:75:ASP:N	2.54	0.41
5:I:55:GLN:O	5:I:57:LYS:N	2.47	0.41
1:J:472:ASP:O	1:J:474:LYS:N	2.53	0.41
1:J:696:GLU:HG2	1:J:700:ASN:HB2	2.02	0.41
1:J:701:VAL:HG22	1:J:734:TYR:HE2	1.85	0.41
3:E:75:ASP:OD1	3:E:75:ASP:N	2.54	0.41
1:J:335:THR:O	1:J:379:ARG:NH1	2.54	0.41
3:C:53:LYS:NZ	3:P:187:SER:OG	2.54	0.41
3:O:23:SER:OG	3:O:24:ASP:N	2.54	0.41
1:J:180:ILE:HG23	1:J:181:LEU:HD12	2.03	0.40
1:J:242:LEU:HD12	1:J:308:LEU:HD12	2.04	0.40
1:J:270:ARG:NH1	1:J:402:CYS:SG	2.95	0.40
3:E:39:ASP:OD2	3:E:42:THR:N	2.42	0.40
5:I:151:SER:HB3	3:N:41:ILE:HG13	2.04	0.40
3:P:75:ASP:OD1	3:P:75:ASP:N	2.54	0.40
3:C:177:LEU:HD11	3:C:182:LEU:HD13	2.03	0.40
3:E:32:ILE:HD12	3:E:32:ILE:HA	1.97	0.40
3:N:128:ARG:NH2	3:O:75:ASP:OD2	2.54	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	J	729/758 (96%)	595 (82%)	120 (16%)	14 (2%)	8	34
2	A	106/121 (88%)	95 (90%)	9 (8%)	2 (2%)	8	34
2	B	106/121 (88%)	96 (91%)	8 (8%)	2 (2%)	8	34
2	M	106/121 (88%)	96 (91%)	8 (8%)	2 (2%)	8	34
3	C	210/220 (96%)	169 (80%)	36 (17%)	5 (2%)	6	28
3	E	210/220 (96%)	169 (80%)	35 (17%)	6 (3%)	4	24
3	N	210/220 (96%)	168 (80%)	37 (18%)	5 (2%)	6	28
3	O	210/220 (96%)	169 (80%)	33 (16%)	8 (4%)	3	19
3	P	210/220 (96%)	170 (81%)	33 (16%)	7 (3%)	4	22
5	I	293/299 (98%)	240 (82%)	47 (16%)	6 (2%)	7	32
All	All	2390/2520 (95%)	1967 (82%)	366 (15%)	57 (2%)	9	28

All (57) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	J	63	LYS
1	J	417	CYS
1	J	473	TYR
1	J	505	GLY
5	I	197	PRO
3	P	35	PRO
1	J	60	GLN
1	J	644	LEU
1	J	677	GLU
3	C	35	PRO
3	C	131	ALA
3	E	35	PRO
5	I	279	PRO
5	I	289	TYR
3	N	116	SER
3	N	131	ALA
3	O	23	SER
3	O	35	PRO
3	P	23	SER
3	P	131	ALA
1	J	226	GLU
1	J	298	TYR
1	J	379	ARG
1	J	462	GLU

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Mol	Chain	Res	Type
2	A	27	ASN
2	B	27	ASN
3	C	23	SER
3	E	23	SER
3	E	131	ALA
5	I	217	THR
2	M	27	ASN
3	N	23	SER
3	O	116	SER
3	O	129	ILE
3	O	131	ALA
3	P	116	SER
1	J	469	VAL
1	J	638	GLU
2	A	95	LYS
2	B	95	LYS
3	C	76	SER
3	C	129	ILE
3	E	76	SER
3	E	129	ILE
5	I	187	GLY
2	M	95	LYS
3	N	76	SER
3	N	129	ILE
3	O	76	SER
3	P	76	SER
3	P	129	ILE
1	J	144	PHE
5	I	250	GLU
3	E	32	ILE
3	O	32	ILE
3	P	32	ILE
3	O	29	ILE

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	J	626/649 (96%)	616 (98%)	10 (2%)	62	79
2	A	96/108 (89%)	95 (99%)	1 (1%)	76	86
2	B	96/108 (89%)	95 (99%)	1 (1%)	76	86
2	M	96/108 (89%)	95 (99%)	1 (1%)	76	86
3	C	181/188 (96%)	174 (96%)	7 (4%)	32	62
3	E	181/188 (96%)	174 (96%)	7 (4%)	32	62
3	N	181/188 (96%)	174 (96%)	7 (4%)	32	62
3	O	181/188 (96%)	174 (96%)	7 (4%)	32	62
3	P	181/188 (96%)	174 (96%)	7 (4%)	32	62
5	I	259/263 (98%)	256 (99%)	3 (1%)	71	83
All	All	2078/2176 (96%)	2027 (98%)	51 (2%)	50	72

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	J	78	ASN
1	J	106	ASN
1	J	129	ASN
1	J	189	ASN
1	J	221	ARG
1	J	223	ASN
1	J	356	ASN
1	J	502	ASN
1	J	602	ASN
1	J	737	ASN
2	A	27	ASN
2	B	27	ASN
3	C	81	ARG
3	C	85	ASN
3	C	105	ASN
3	C	115	ARG
3	C	134	ASN
3	C	159	ASN
3	C	213	ASN
3	E	81	ARG
3	E	85	ASN
3	E	105	ASN
3	E	115	ARG
3	E	134	ASN

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Mol	Chain	Res	Type
3	E	159	ASN
3	E	213	ASN
5	I	57	LYS
5	I	198	LEU
5	I	207	ASN
2	M	27	ASN
3	N	81	ARG
3	N	85	ASN
3	N	105	ASN
3	N	115	ARG
3	N	134	ASN
3	N	159	ASN
3	N	213	ASN
3	O	81	ARG
3	O	85	ASN
3	O	105	ASN
3	O	115	ARG
3	O	134	ASN
3	O	159	ASN
3	O	213	ASN
3	P	81	ARG
3	P	85	ASN
3	P	105	ASN
3	P	115	ARG
3	P	134	ASN
3	P	159	ASN
3	P	213	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (49) such sidechains are listed below:

Mol	Chain	Res	Type
1	J	68	HIS
1	J	106	ASN
1	J	112	ASN
1	J	129	ASN
1	J	135	ASN
1	J	169	GLN
1	J	189	ASN
1	J	223	ASN
1	J	389	ASN
1	J	403	HIS
1	J	432	HIS

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Mol	Chain	Res	Type
1	J	434	HIS
1	J	482	HIS
1	J	489	GLN
1	J	502	ASN
1	J	602	ASN
1	J	737	ASN
2	A	27	ASN
2	B	27	ASN
3	C	11	GLN
3	C	85	ASN
3	C	105	ASN
3	C	134	ASN
3	C	213	ASN
3	E	11	GLN
3	E	85	ASN
3	E	134	ASN
3	E	213	ASN
5	I	83	GLN
5	I	137	HIS
5	I	207	ASN
5	I	233	HIS
5	I	246	HIS
5	I	255	GLN
2	M	27	ASN
3	N	11	GLN
3	N	85	ASN
3	N	124	ASN
3	N	134	ASN
3	N	213	ASN
3	O	11	GLN
3	O	85	ASN
3	O	134	ASN
3	O	213	ASN
3	P	11	GLN
3	P	85	ASN
3	P	105	ASN
3	P	134	ASN
3	P	213	ASN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
4	H	39/72 (54%)	9 (23%)	1 (2%)

All (9) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
4	H	2	C
4	H	3	G
4	H	13	G
4	H	15	A
4	H	21	U
4	H	27	C
4	H	29	G
4	H	33	A
4	H	39	U

All (1) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
4	H	12	C

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](#)

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
6	ATP	J	801	-	26,33,33	0.92	1 (3%)	31,52,52	1.59	5 (16%)

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
6	ATP	J	801	-	-	1/18/38/38	0/3/3/3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	J	801	ATP	C5-C4	2.49	1.47	1.40

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	J	801	ATP	PB-O3B-PG	-3.85	119.60	132.83
6	J	801	ATP	PA-O3A-PB	-3.68	120.19	132.83
6	J	801	ATP	C3'-C2'-C1'	3.20	105.79	100.98
6	J	801	ATP	N3-C2-N1	-3.19	123.69	128.68
6	J	801	ATP	C4-C5-N7	-2.68	106.60	109.40

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	J	801	ATP	C5'-O5'-PA-O1A

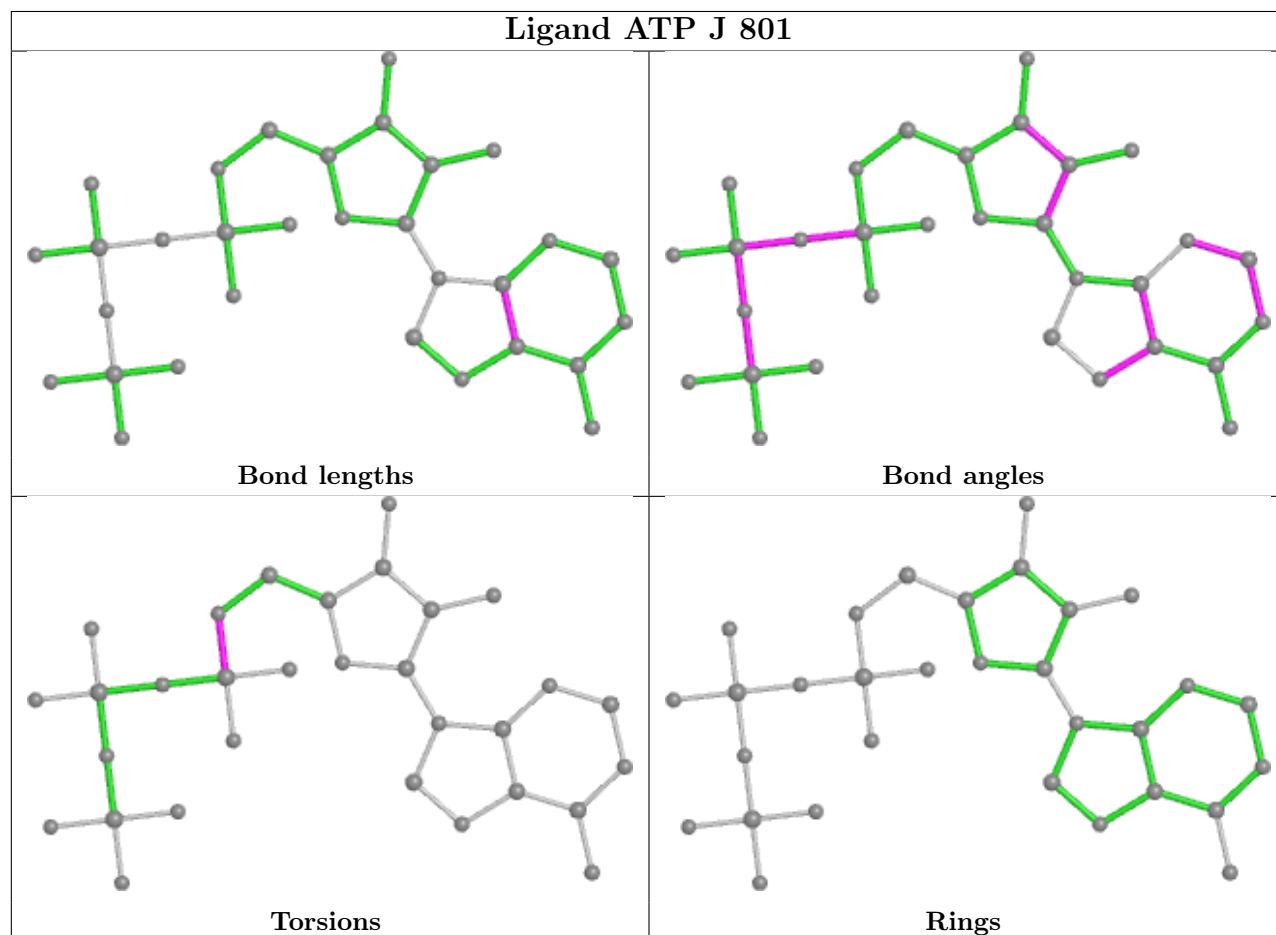
There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	J	801	ATP	2	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.

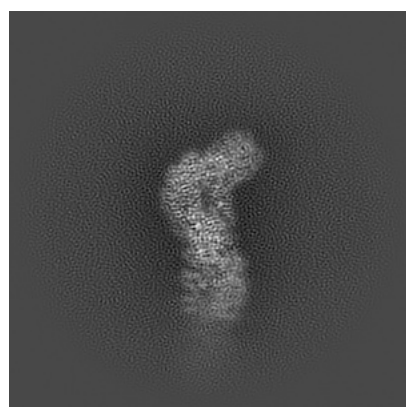
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-0519. These allow visual inspection of the internal detail of the map and identification of artifacts.

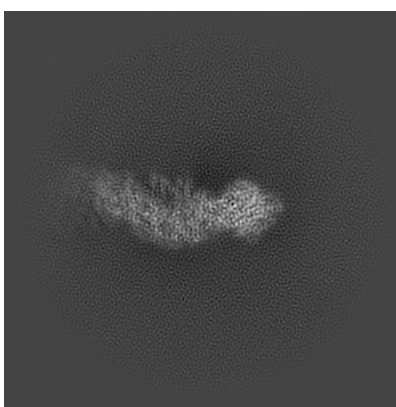
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

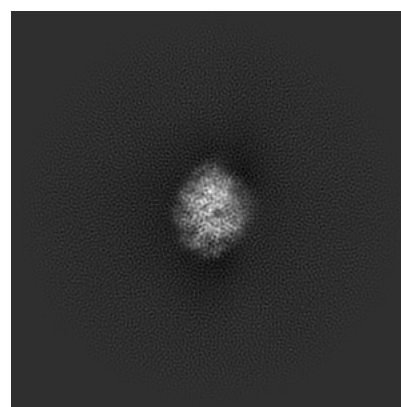
6.1.1 Primary map



X



Y

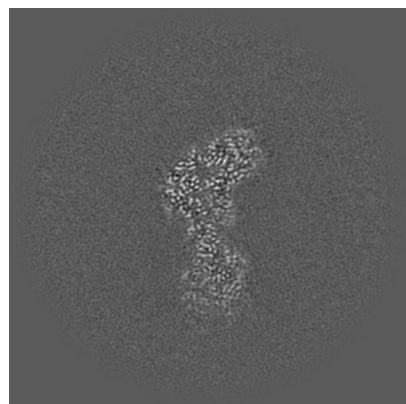


Z

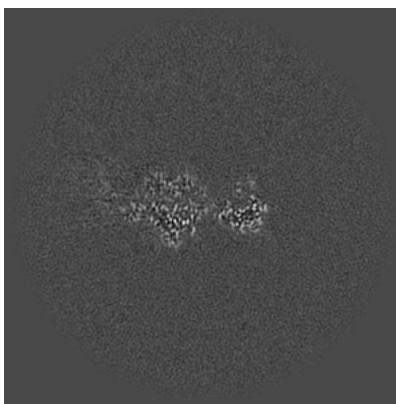
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

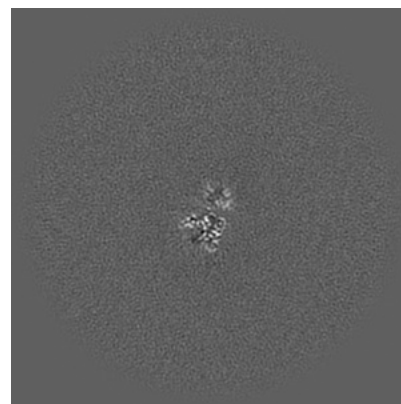
6.2.1 Primary map



X Index: 192



Y Index: 192

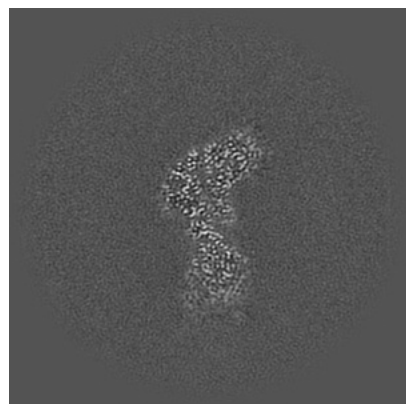


Z Index: 192

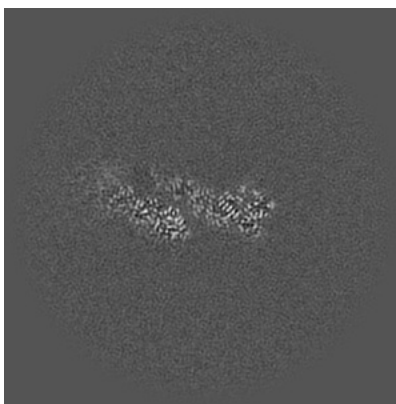
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

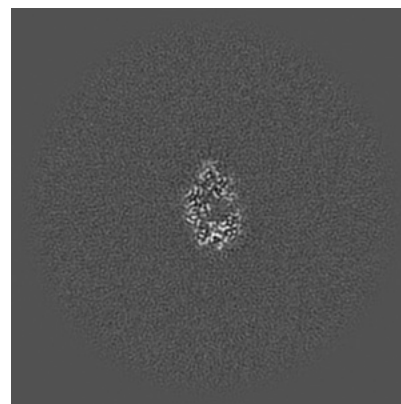
6.3.1 Primary map



X Index: 190



Y Index: 200



Z Index: 230

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

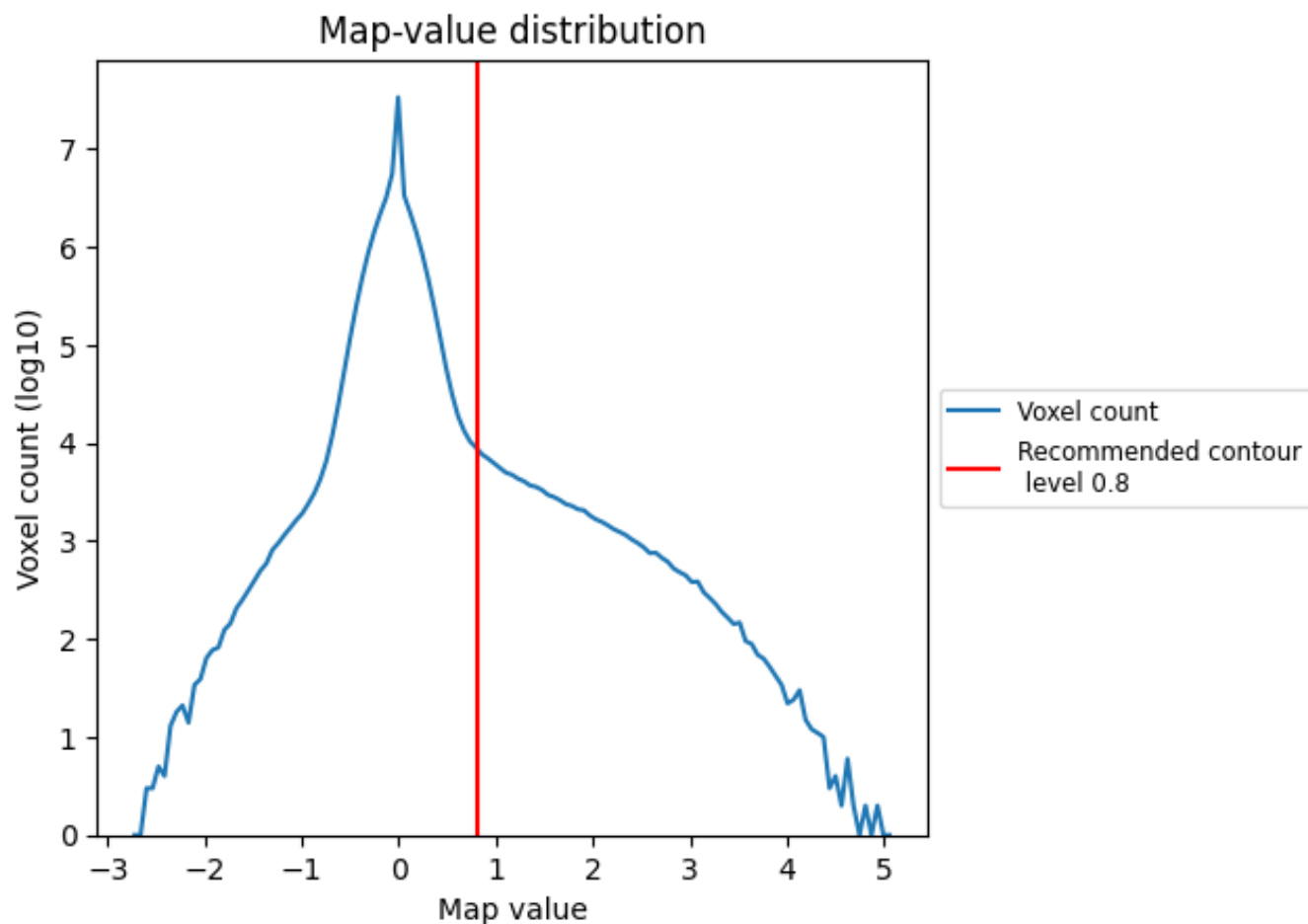
6.5 Mask visualisation

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

7 Map analysis [i](#)

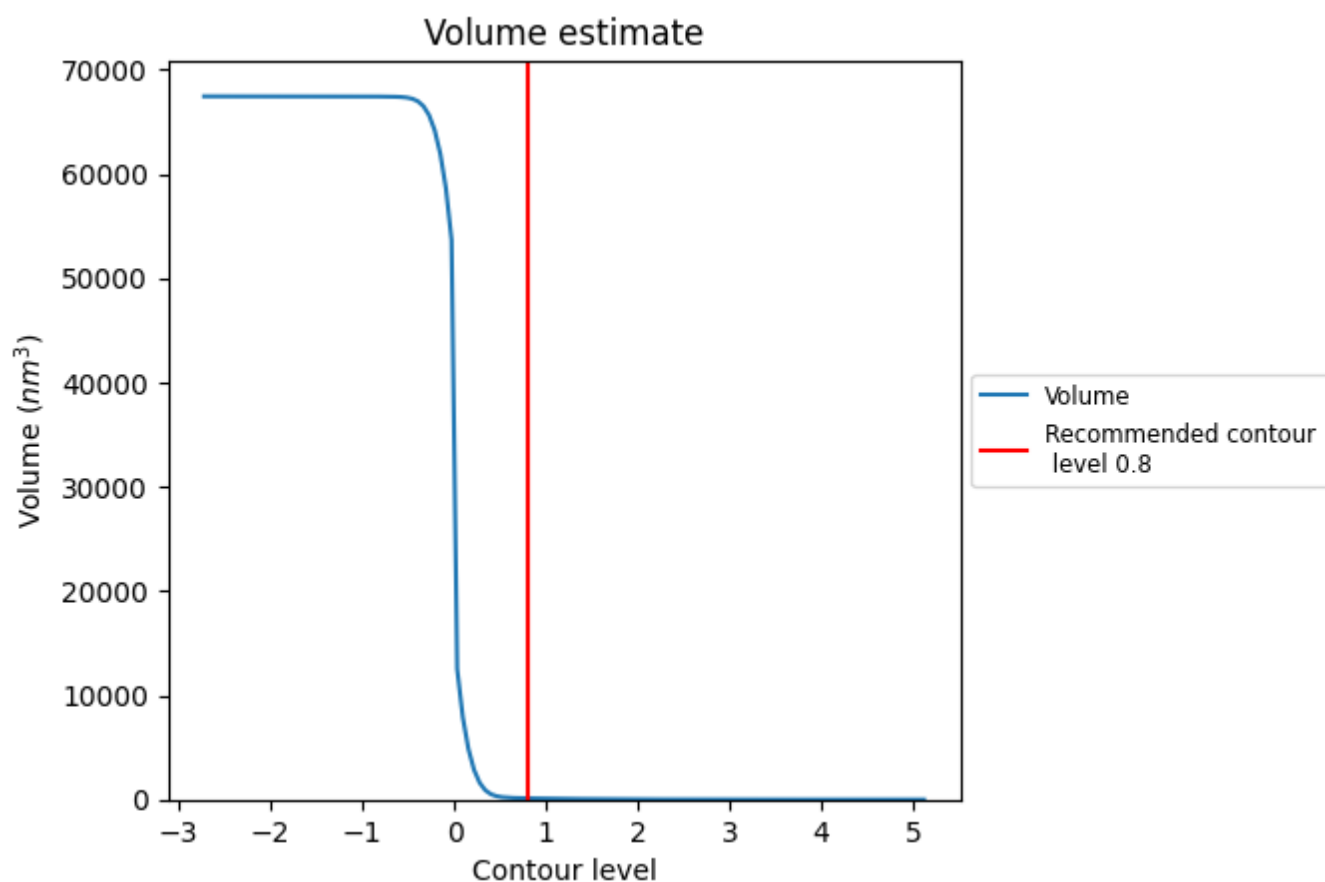
This section contains the results of statistical analysis of the map.

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.

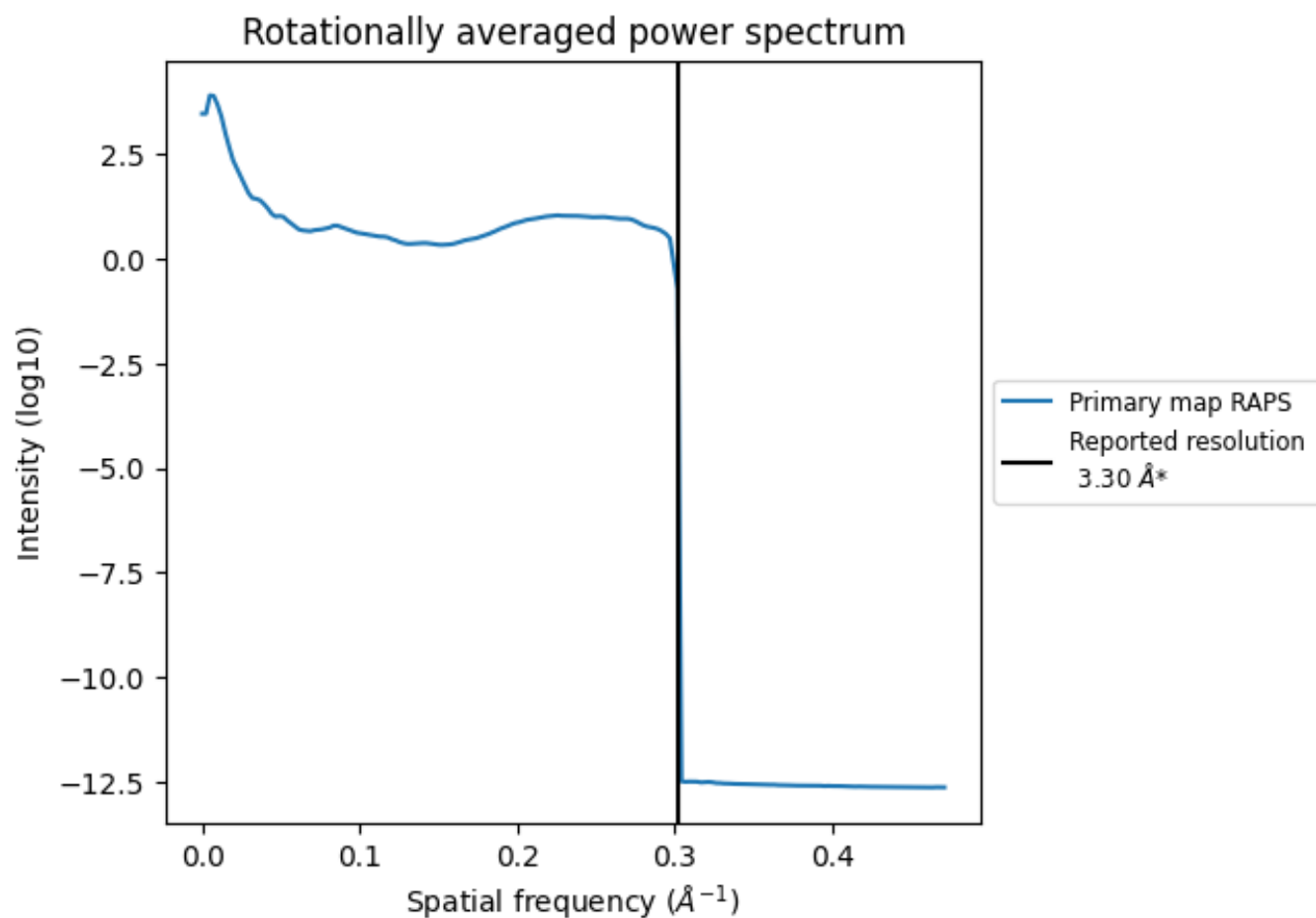
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 120 nm³; this corresponds to an approximate mass of 108 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 0.303 Å⁻¹

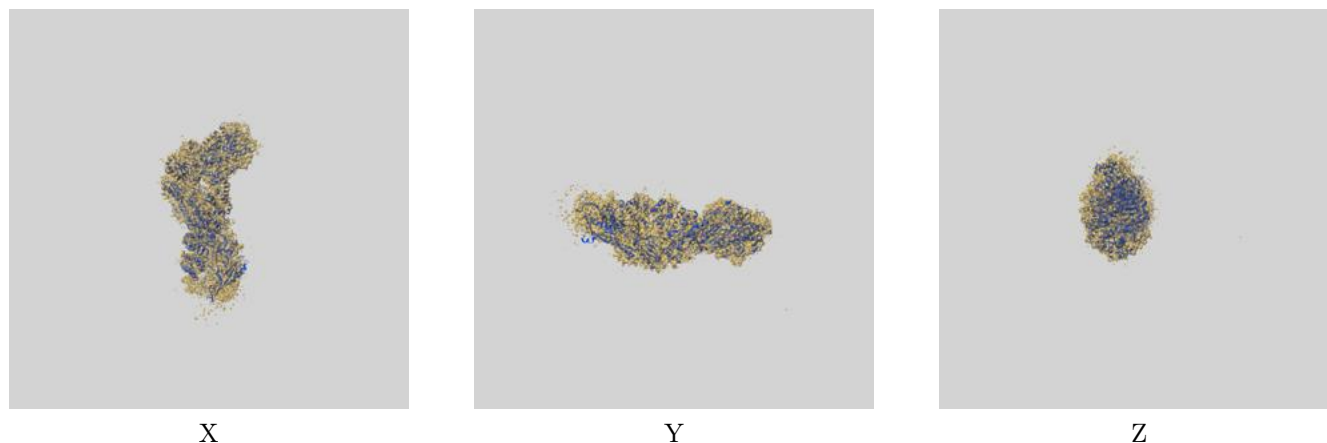
8 Fourier-Shell correlation

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

9 Map-model fit [i](#)

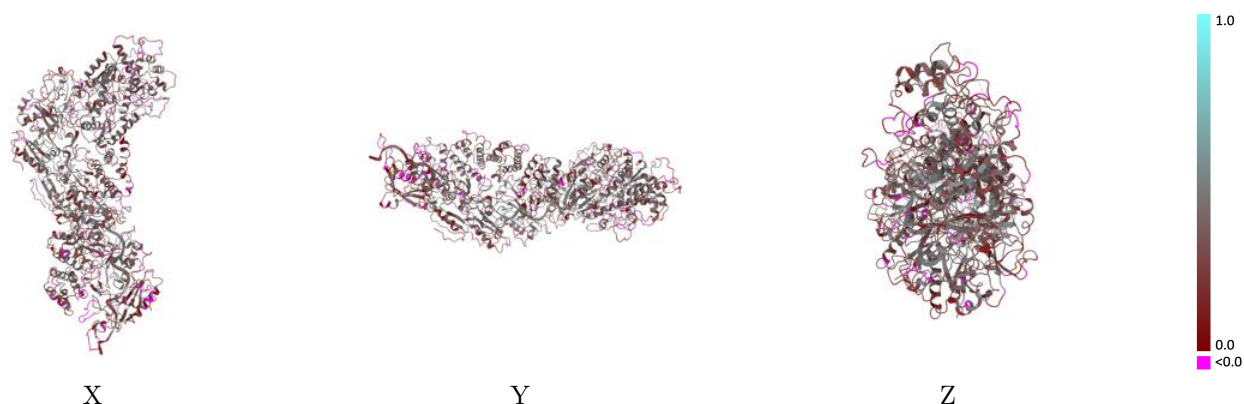
This section contains information regarding the fit between EMDB map EMD-0519 and PDB model 6NUE. Per-residue inclusion information can be found in section [3](#) on page [6](#).

9.1 Map-model overlay [i](#)



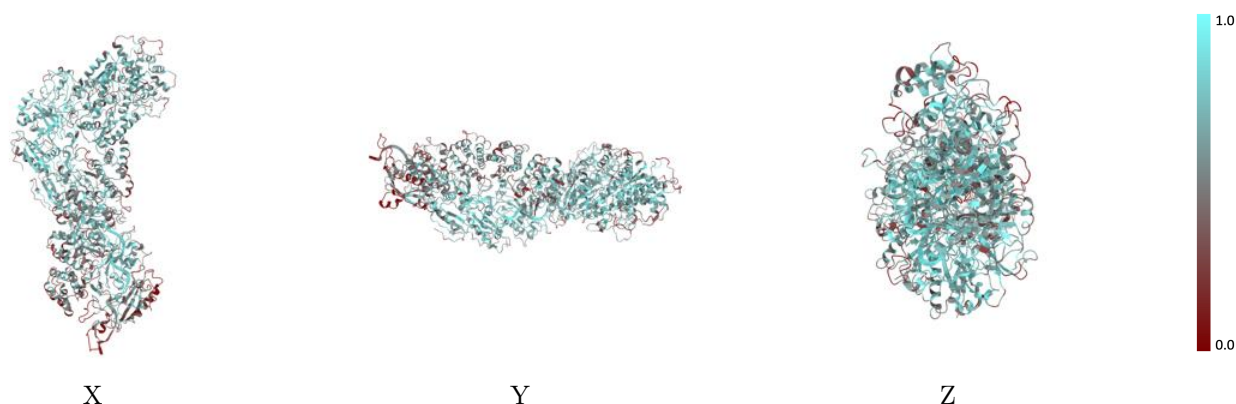
The images above show the 3D surface view of the map at the recommended contour level 0.8 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



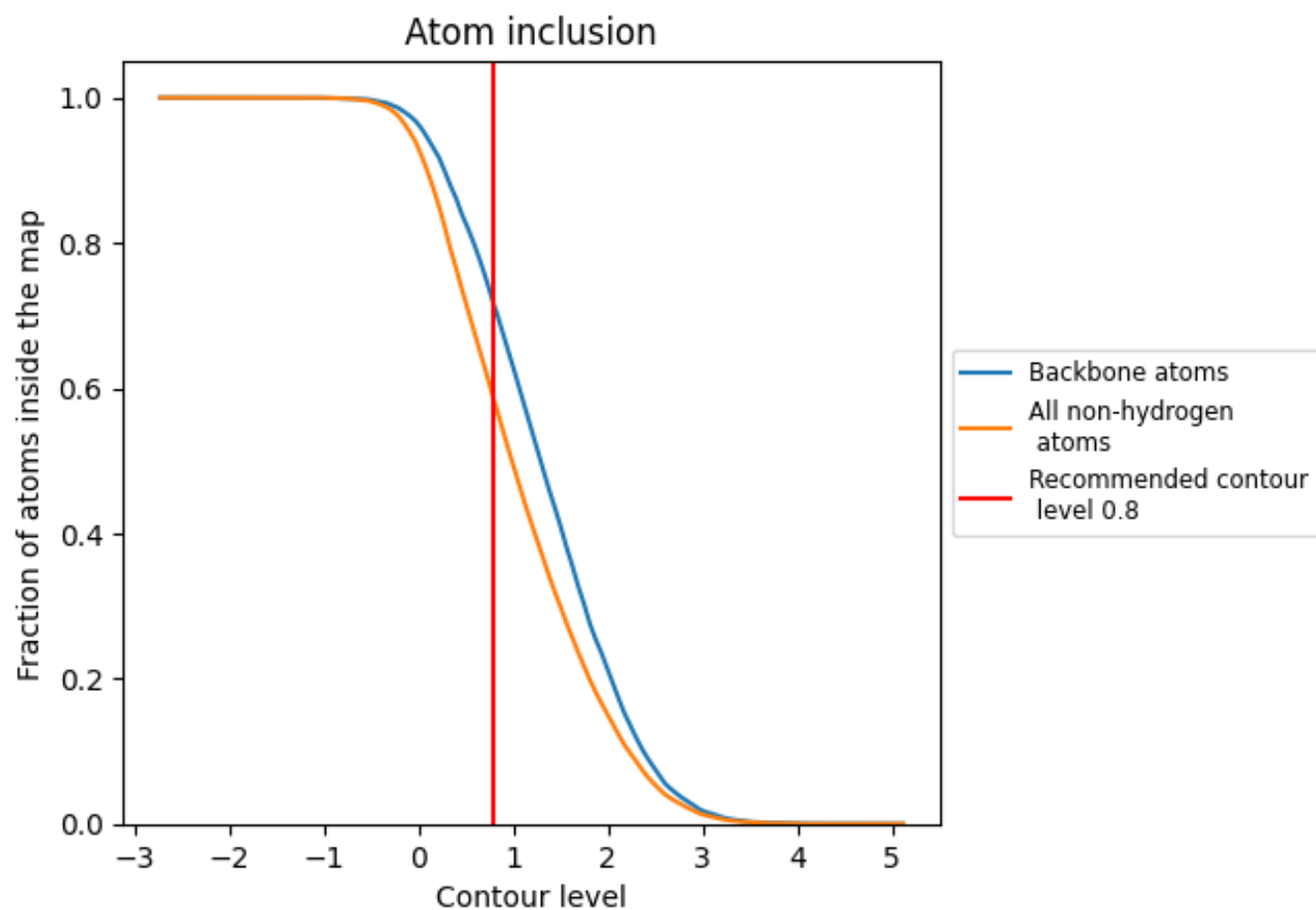
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.8).

9.4 Atom inclusion [i](#)



At the recommended contour level, 71% of all backbone atoms, 58% 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 (0.8) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div></div> 0.5824	<div></div> 0.3050
A	<div></div> 0.5063	<div></div> 0.3130
B	<div></div> 0.4419	<div></div> 0.2820
C	<div></div> 0.3223	<div></div> 0.1880
E	<div></div> 0.6349	<div></div> 0.3120
H	<div></div> 0.7452	<div></div> 0.3920
I	<div></div> 0.6444	<div></div> 0.3220
J	<div></div> 0.6066	<div></div> 0.3090
M	<div></div> 0.5293	<div></div> 0.3150
N	<div></div> 0.6178	<div></div> 0.3160
O	<div></div> 0.6343	<div></div> 0.3200
P	<div></div> 0.5885	<div></div> 0.3060

1.0

0.0

<0.0