



wwPDB EM Validation Summary Report ⓘ

Mar 27, 2026 – 04:00 PM UTC

PDB ID : 6S91 / pdb_00006s91
EMDB ID : EMD-10126
Title : Cryo-EM structure of the Type III-B Cmr-beta bound to cognate target RNA and AMPPnP, state 2
Authors : Sofos, N.; Montoya, G.; Stella, S.
Deposited on : 2019-07-11
Resolution : 2.68 Å(reported)

This is a wwPDB EM Validation Summary 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:

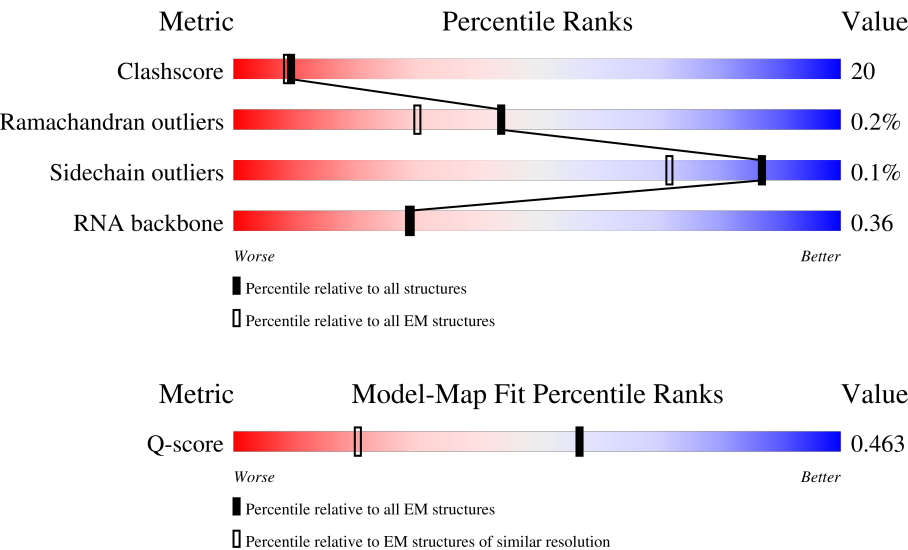
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data 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 2.68 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.



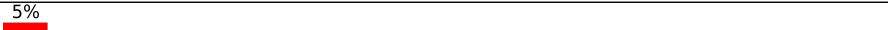
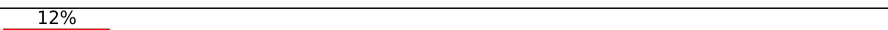
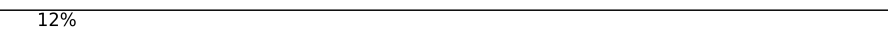
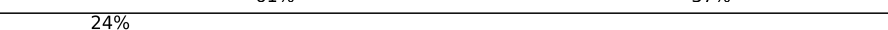
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
RNA backbone	8273	3508	-
Q-score	-	25397	9255 (2.18 - 3.18)

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	155	
1	B	155	
1	C	155	

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Mol	Chain	Length	Quality of chain
2	D	286	
2	E	286	
2	F	286	
2	G	286	
3	H	313	
4	I	296	
5	J	476	
6	K	1037	
7	U	46	
8	V	51	
9	L	174	
9	M	174	
9	N	174	
9	O	174	
9	P	174	
9	Q	174	
9	R	174	
9	S	174	
9	T	174	
9	W	174	
9	X	174	
9	l	174	
9	m	174	
9	n	174	
9	o	174	

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Mol	Chain	Length	Quality of chain
9	p	174	
9	q	174	
9	r	174	
9	s	174	
9	t	174	
9	w	174	
9	x	174	

2 Entry composition

There are 12 unique types of molecules in this entry. The entry contains 62045 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-associated protein, Cmr5 family.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	153	Total	C	N	O	S	0	0
			1253	817	205	230	1		
1	B	151	Total	C	N	O	S	0	0
			1235	809	200	225	1		
1	C	154	Total	C	N	O	S	0	0
			1261	823	206	231	1		

- Molecule 2 is a protein called CRISPR-associated RAMP protein, Cmr4 family.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	D	285	Total	C	N	O	S	0	0
			2274	1478	369	425	2		
2	E	285	Total	C	N	O	S	0	0
			2274	1478	369	425	2		
2	F	285	Total	C	N	O	S	0	0
			2274	1478	369	425	2		
2	G	283	Total	C	N	O	S	0	0
			2258	1467	367	422	2		

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	31	ALA	ASP	engineered mutation	UNP F0NDX6
E	31	ALA	ASP	engineered mutation	UNP F0NDX6
F	31	ALA	ASP	engineered mutation	UNP F0NDX6
G	31	ALA	ASP	engineered mutation	UNP F0NDX6

- Molecule 3 is a protein called CRISPR-associated protein, Cmr3 family.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	H	312	Total	C	N	O	S	0	0
			2528	1630	418	473	7		

- Molecule 4 is a protein called CRISPR-associated RAMP protein, Cmr6 family.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	I	284	Total	C	N	O	S	0	0
			2282	1470	381	427	4		

There are 13 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
I	284	ALA	-	expression tag	UNP F0NDX3
I	285	ALA	-	expression tag	UNP F0NDX3
I	286	ALA	-	expression tag	UNP F0NDX3
I	287	HIS	-	expression tag	UNP F0NDX3
I	288	HIS	-	expression tag	UNP F0NDX3
I	289	HIS	-	expression tag	UNP F0NDX3
I	290	HIS	-	expression tag	UNP F0NDX3
I	291	HIS	-	expression tag	UNP F0NDX3
I	292	HIS	-	expression tag	UNP F0NDX3
I	293	HIS	-	expression tag	UNP F0NDX3
I	294	HIS	-	expression tag	UNP F0NDX3
I	295	HIS	-	expression tag	UNP F0NDX3
I	296	HIS	-	expression tag	UNP F0NDX3

- Molecule 5 is a protein called Cmr1,CRISPR-associated RAMP protein, Cmr1 family.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	J	467	Total	C	N	O	S	0	0
			3833	2481	622	717	13		

- Molecule 6 is a protein called CRISPR-associated protein, Cmr2 family.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	K	1023	Total	C	N	O	S	0	0
			8408	5438	1392	1554	24		

- Molecule 7 is a RNA chain called Cognate target RNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	U	36	Total	C	N	O	P	0	0
			758	340	127	256	35		

- Molecule 8 is a RNA chain called crRNA (48-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
8	V	48	Total	C	N	O	P	0	0
			1025	461	192	325	47		

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
V	1	A	C	conflict	GB 323473489
V	3	U	G	conflict	GB 323473489

- Molecule 9 is a protein called CRISPR-associated protein CmrX.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	L	173	Total	C	N	O	S	0	0
			1378	880	227	269	2		
9	M	173	Total	C	N	O	S	0	0
			1378	880	227	269	2		
9	N	173	Total	C	N	O	S	0	0
			1378	880	227	269	2		
9	O	173	Total	C	N	O	S	0	0
			1378	880	227	269	2		
9	P	173	Total	C	N	O	S	0	0
			1378	880	227	269	2		
9	Q	173	Total	C	N	O	S	0	0
			1378	880	227	269	2		
9	S	173	Total	C	N	O	S	0	0
			1378	880	227	269	2		
9	T	173	Total	C	N	O	S	0	0
			1378	880	227	269	2		
9	W	173	Total	C	N	O	S	0	0
			1378	880	227	269	2		
9	X	173	Total	C	N	O	S	0	0
			1378	880	227	269	2		
9	l	173	Total	C	N	O	S	0	0
			1378	880	227	269	2		
9	m	173	Total	C	N	O	S	0	0
			1378	880	227	269	2		
9	n	173	Total	C	N	O	S	0	0
			1378	880	227	269	2		
9	o	173	Total	C	N	O	S	0	0
			1378	880	227	269	2		
9	p	173	Total	C	N	O	S	0	0
			1378	880	227	269	2		
9	q	173	Total	C	N	O	S	0	0
			1378	880	227	269	2		

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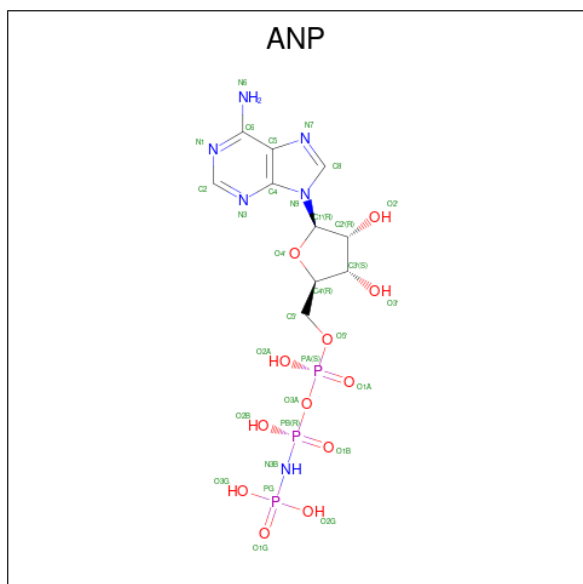
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Mol	Chain	Residues	Atoms					AltConf	Trace
9	r	173	Total	C	N	O	S	0	0
			1378	880	227	269	2		
9	s	173	Total	C	N	O	S	0	0
			1378	880	227	269	2		
9	t	173	Total	C	N	O	S	0	0
			1378	880	227	269	2		
9	w	173	Total	C	N	O	S	0	0
			1378	880	227	269	2		
9	x	173	Total	C	N	O	S	0	0
			1378	880	227	269	2		
9	R	173	Total	C	N	O	S	0	0
			1378	880	227	269	2		

- Molecule 10 is ZINC ION (CCD ID: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		AltConf
10	K	1	Total	Zn	0
			1	1	

- Molecule 11 is PHOSPHOAMINOPHOSPHONIC ACID-ADENYLATE ESTER (CCD ID: ANP) (formula: C₁₀H₁₇N₆O₁₂P₃) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					AltConf
11	K	1	Total	C	N	O	P	0
			31	10	6	12	3	

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Mol	Chain	Residues	Atoms					AltConf
11	K	1	Total	C	N	O	P	0
			31	10	6	12	3	

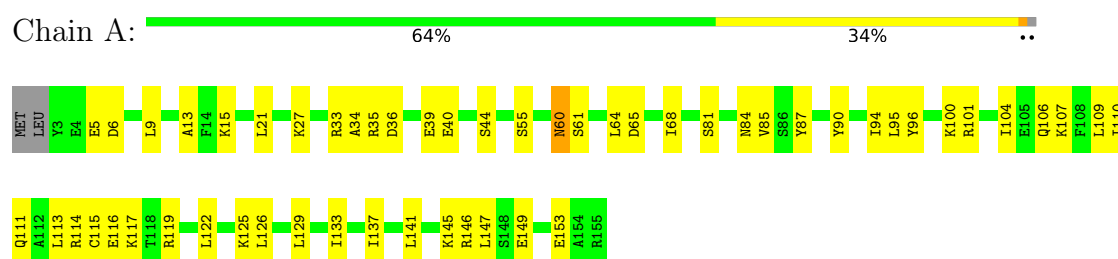
- Molecule 12 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		AltConf
12	K	3	Total	Mn	0
			3	3	

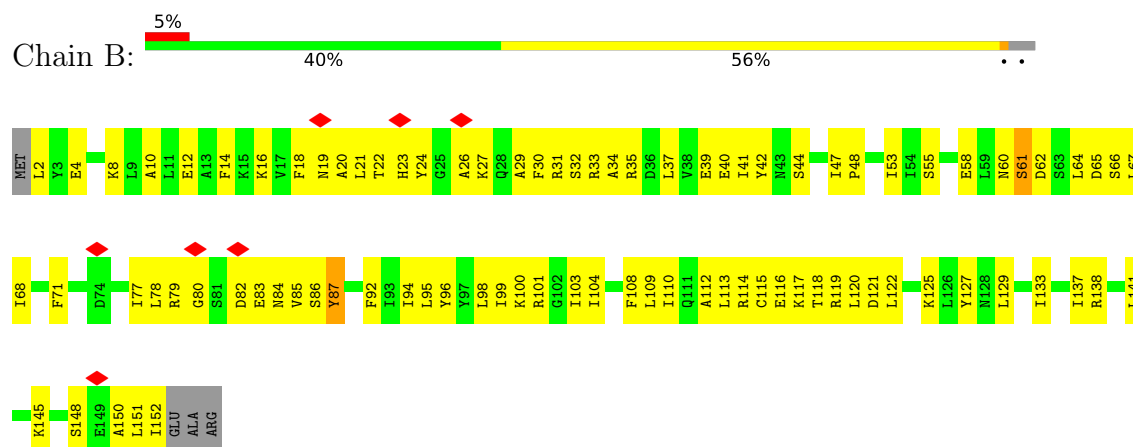
3 Residue-property plots

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.

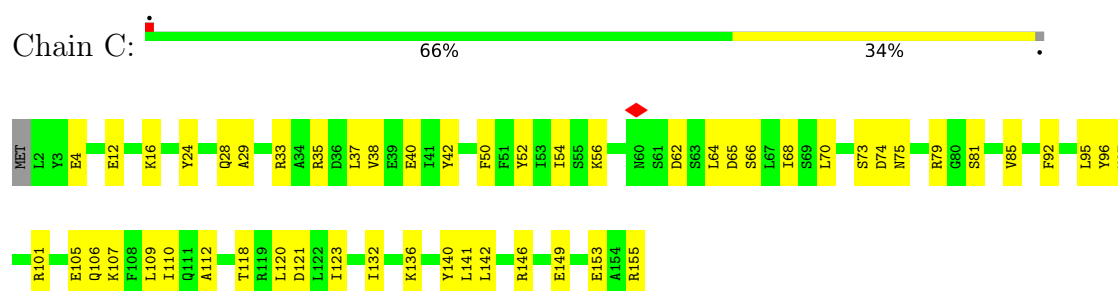
- Molecule 1: CRISPR-associated protein, Cmr5 family



- Molecule 1: CRISPR-associated protein, Cmr5 family

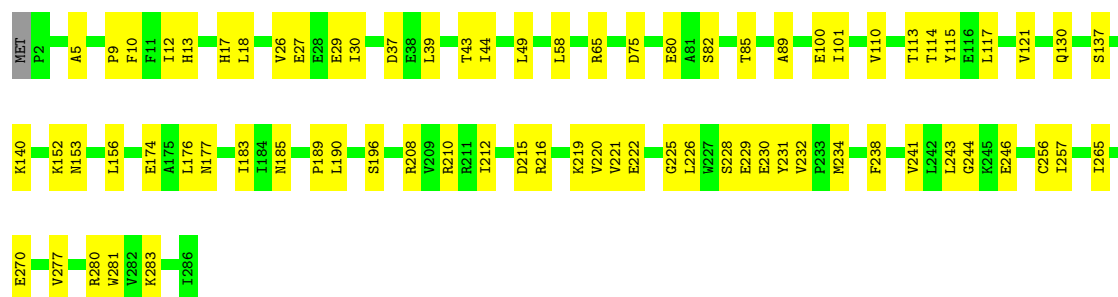


- Molecule 1: CRISPR-associated protein, Cmr5 family



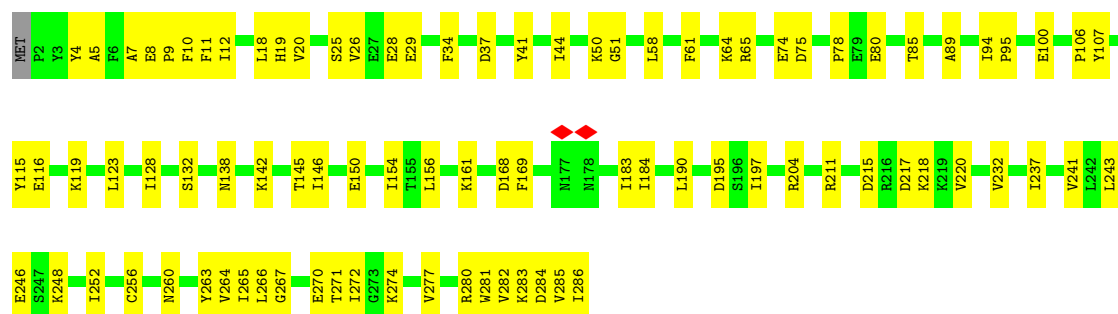
- Molecule 2: CRISPR-associated RAMP protein, Cmr4 family

Chain D:  73% 26%



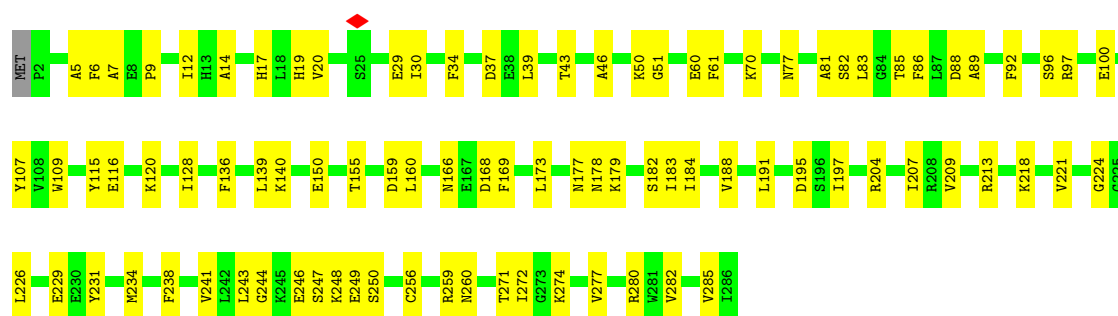
- Molecule 2: CRISPR-associated RAMP protein, Cmr4 family

Chain E:  69% 31%



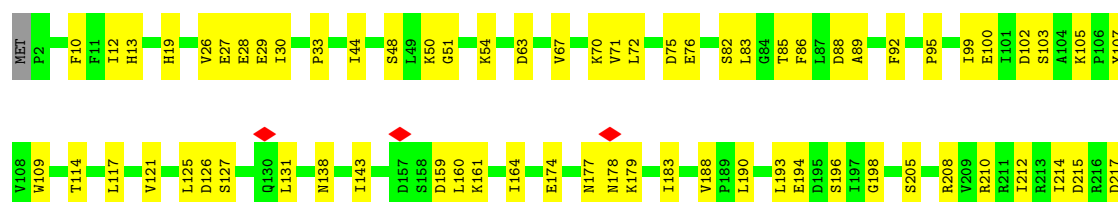
- Molecule 2: CRISPR-associated RAMP protein, Cmr4 family

Chain F:  68% 31%

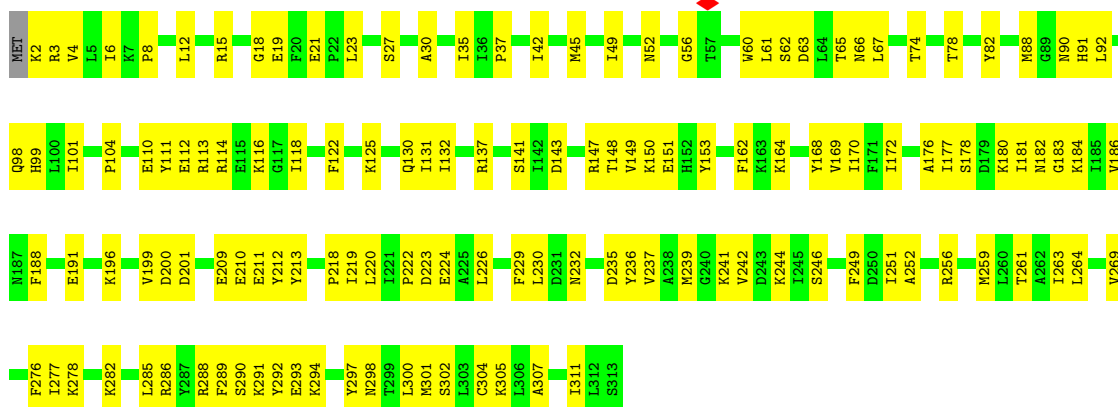


- Molecule 2: CRISPR-associated RAMP protein, Cmr4 family

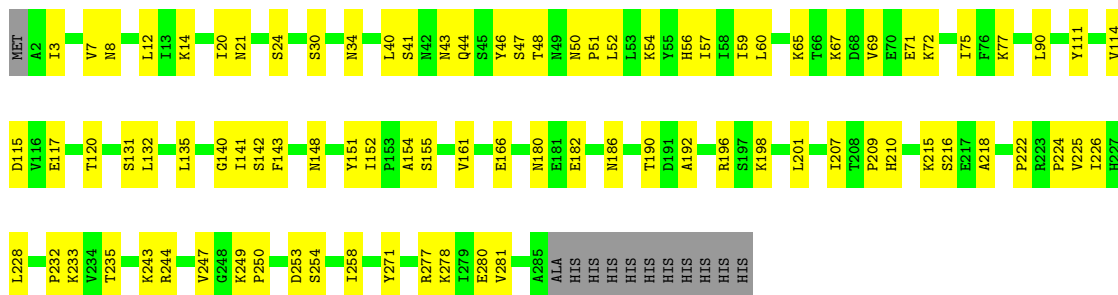
Chain G:  64% 35%



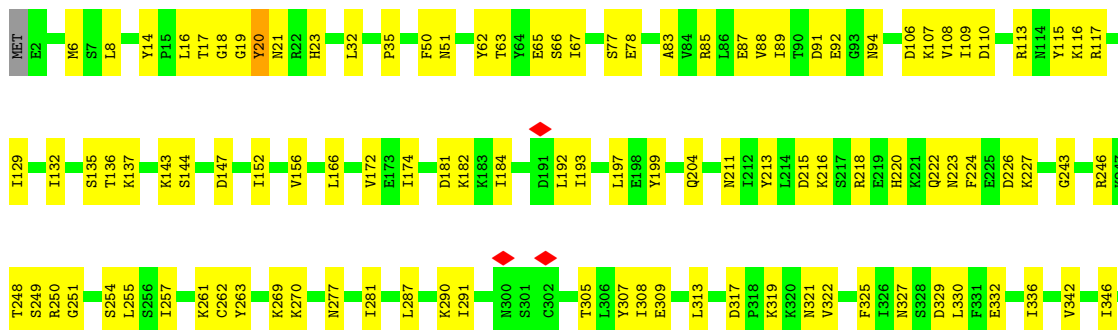
- Molecule 3: CRISPR-associated protein, Cmr3 family

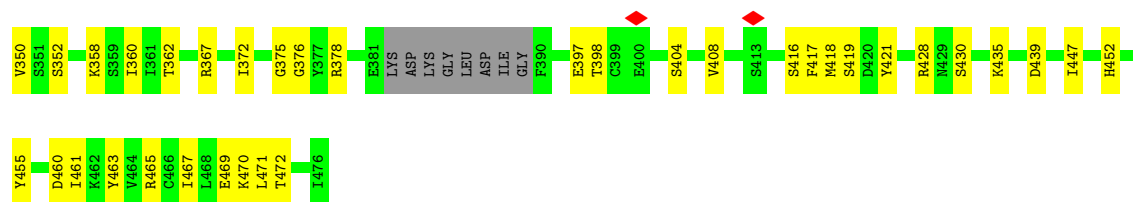


- Molecule 4: CRISPR-associated RAMP protein, Cmr6 family

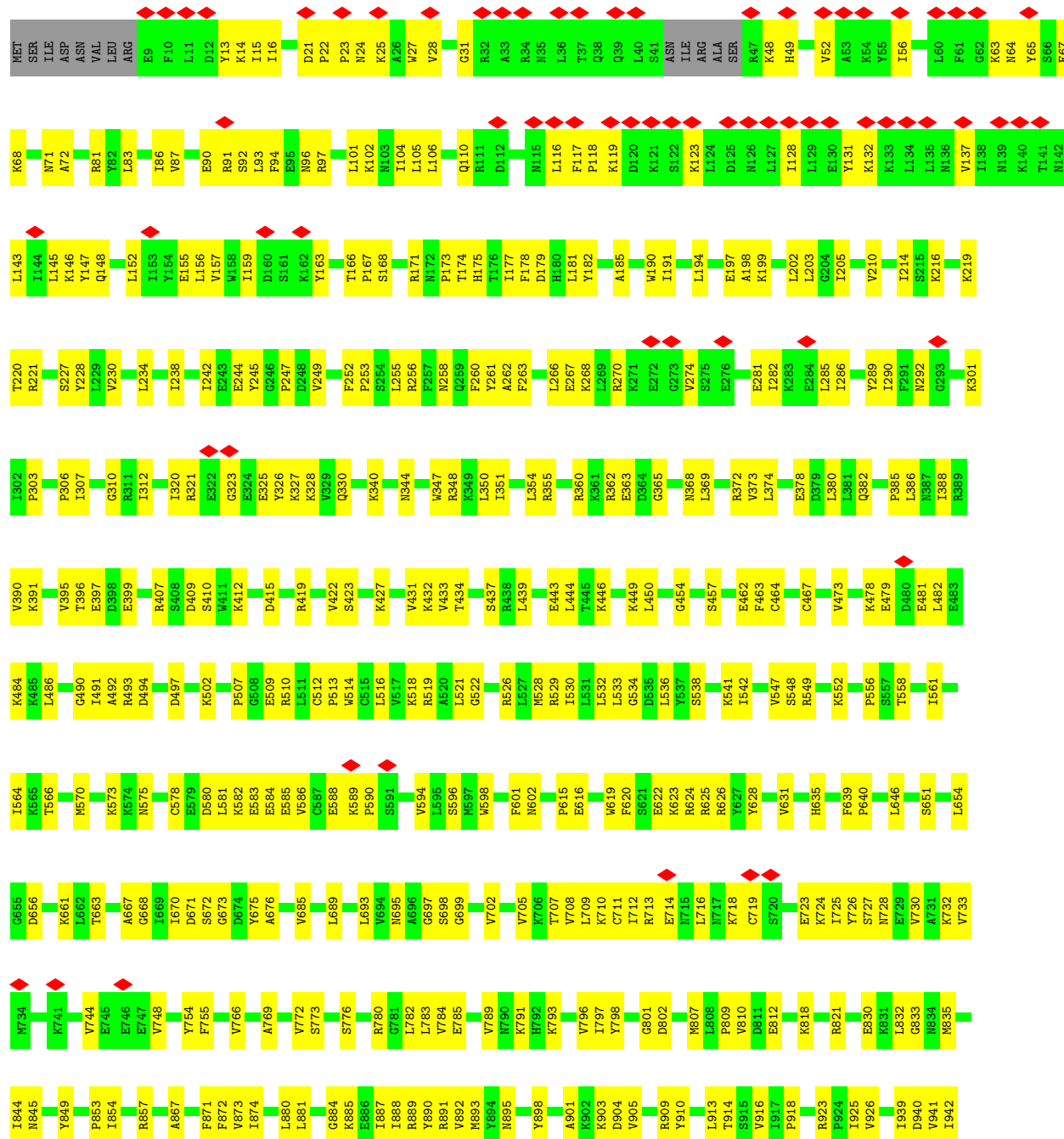


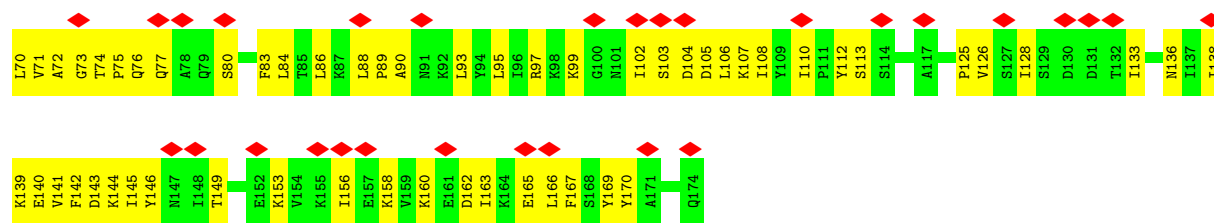
- Molecule 5: Cmr1, CRISPR-associated RAMP protein, Cmr1 family



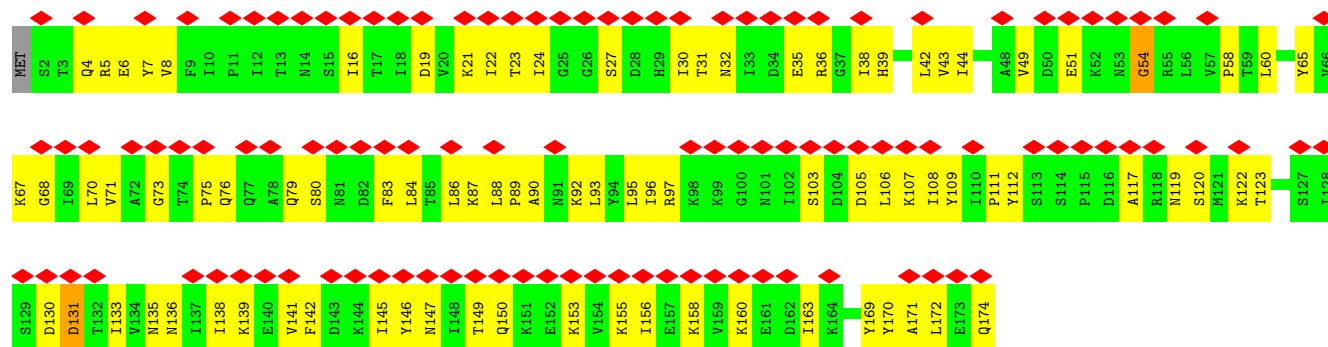


• Molecule 6: CRISPR-associated protein, Cmr2 family

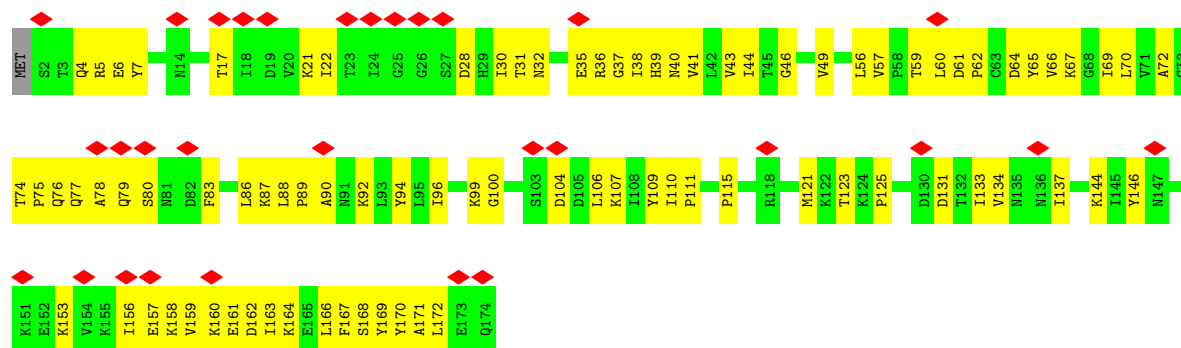




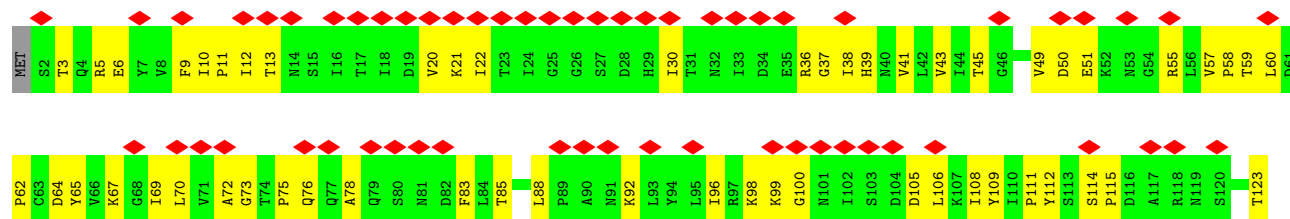
• Molecule 9: CRISPR-associated protein CmrX



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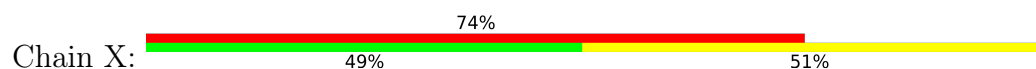


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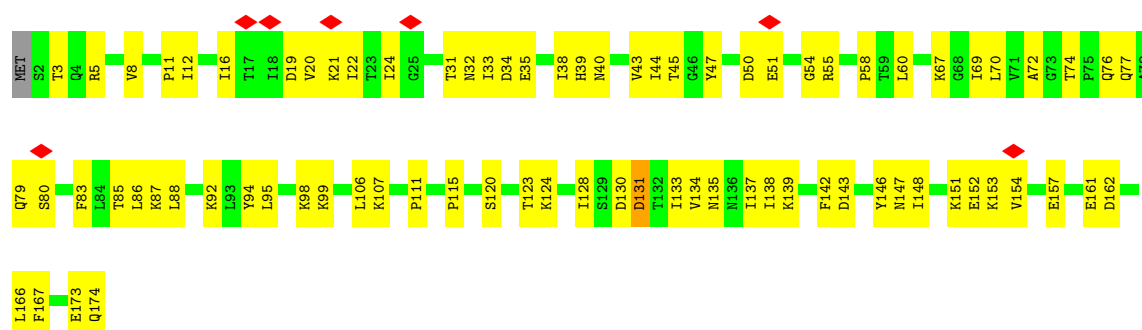




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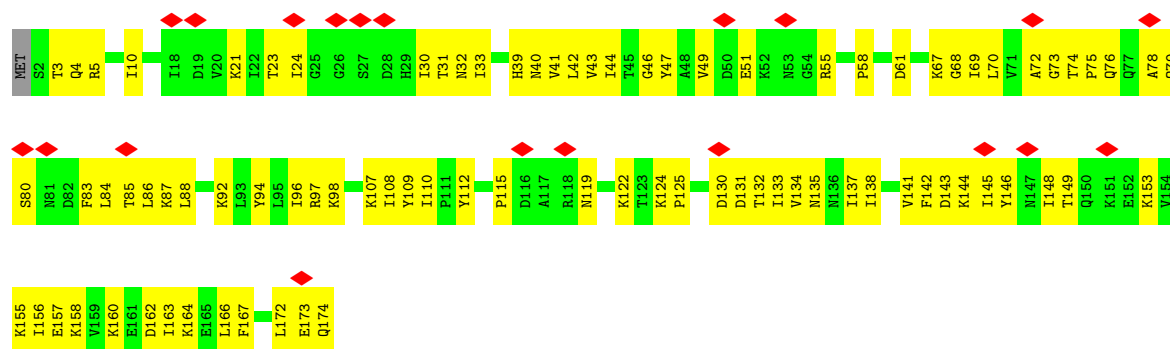


• Molecule 9: CRISPR-associated protein CmrX

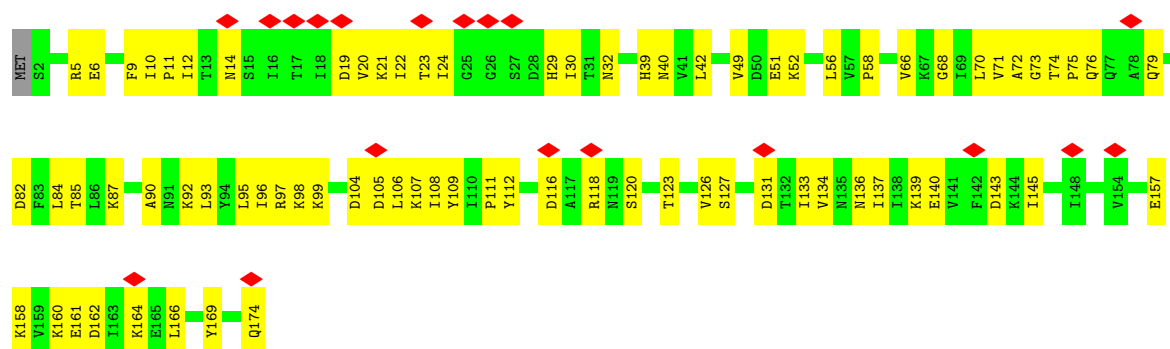


• Molecule 9: CRISPR-associated protein CmrX

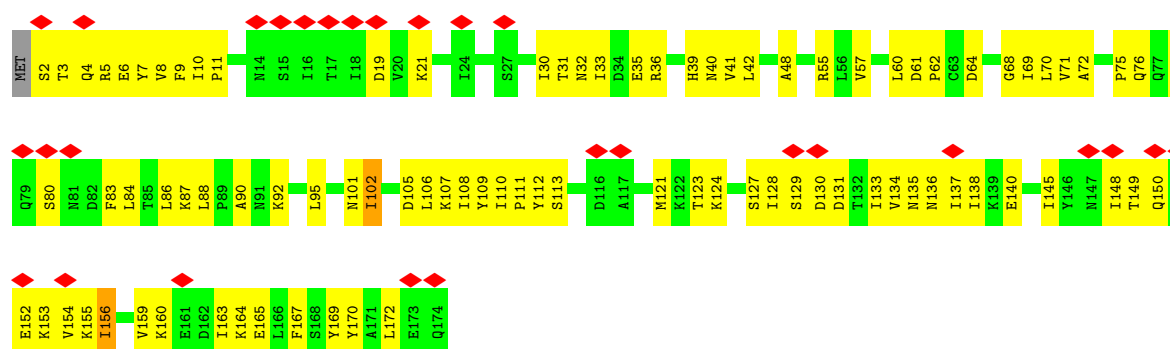




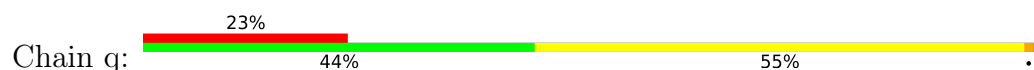
• Molecule 9: CRISPR-associated protein CmrX

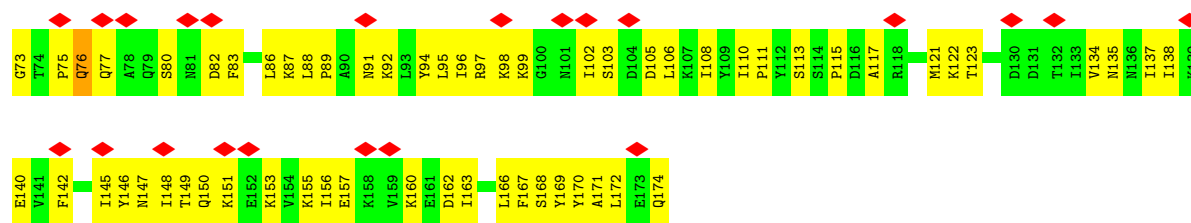


• Molecule 9: CRISPR-associated protein CmrX

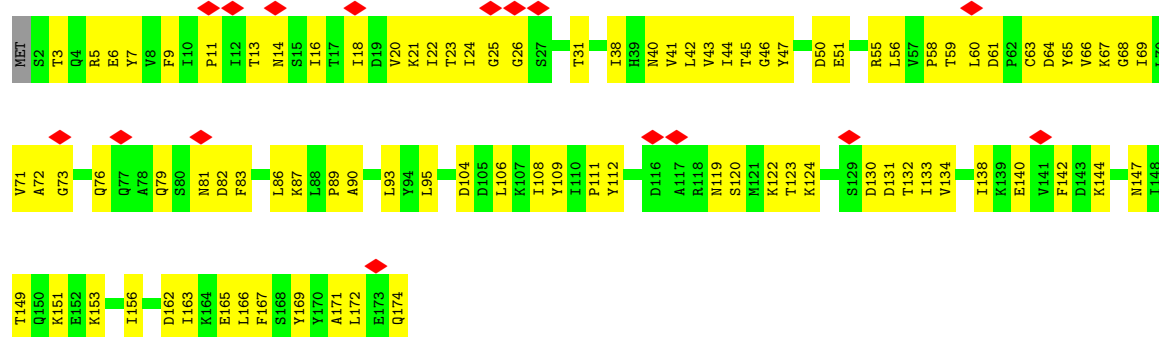


• Molecule 9: CRISPR-associated protein CmrX

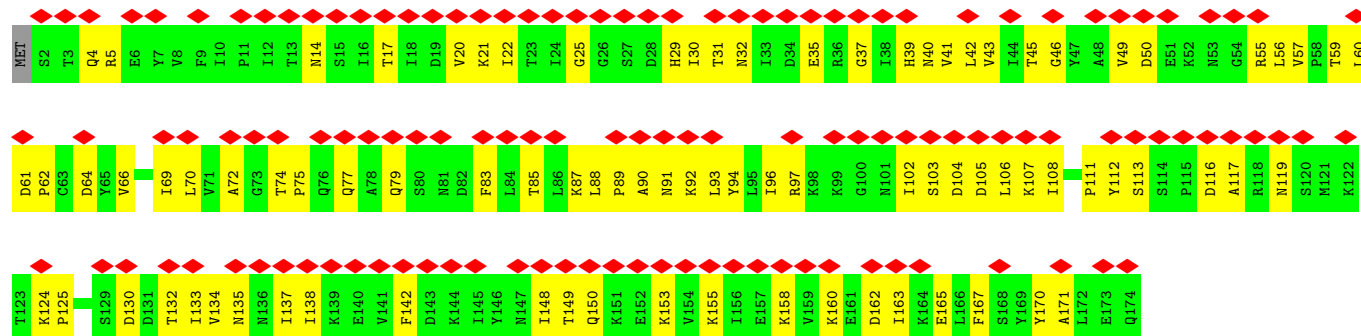




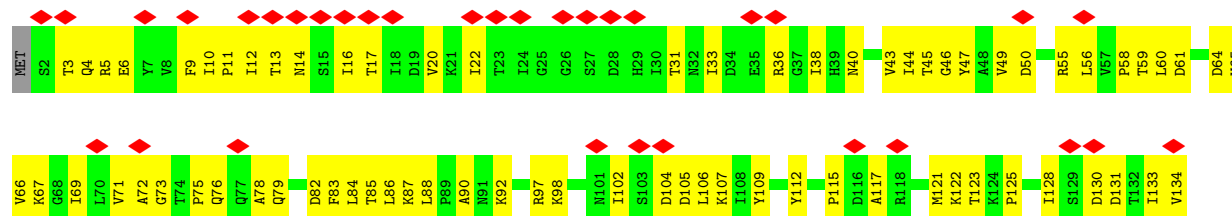
• Molecule 9: CRISPR-associated protein CmrX

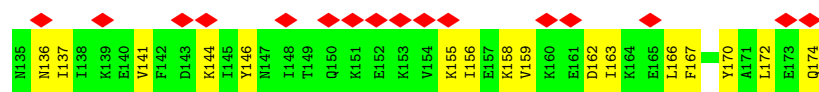


• Molecule 9: CRISPR-associated protein CmrX

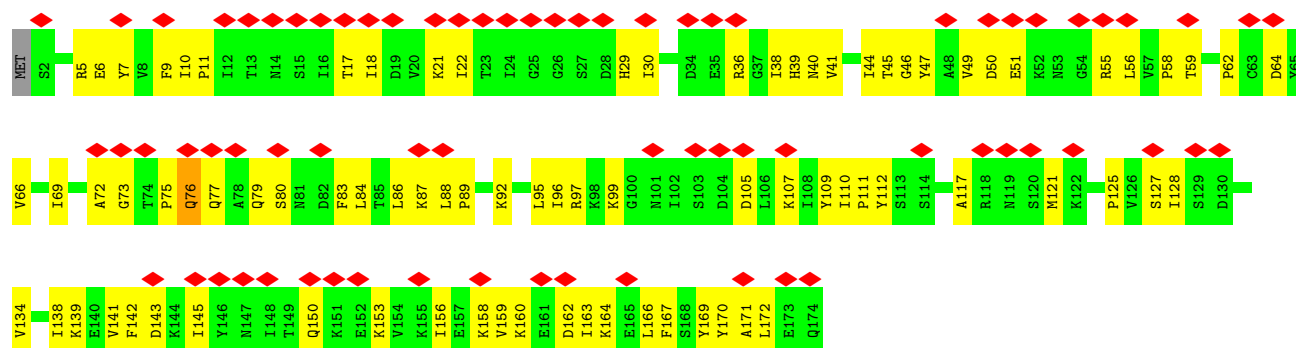
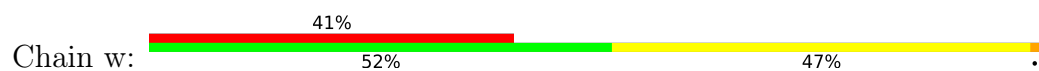


• Molecule 9: CRISPR-associated protein CmrX

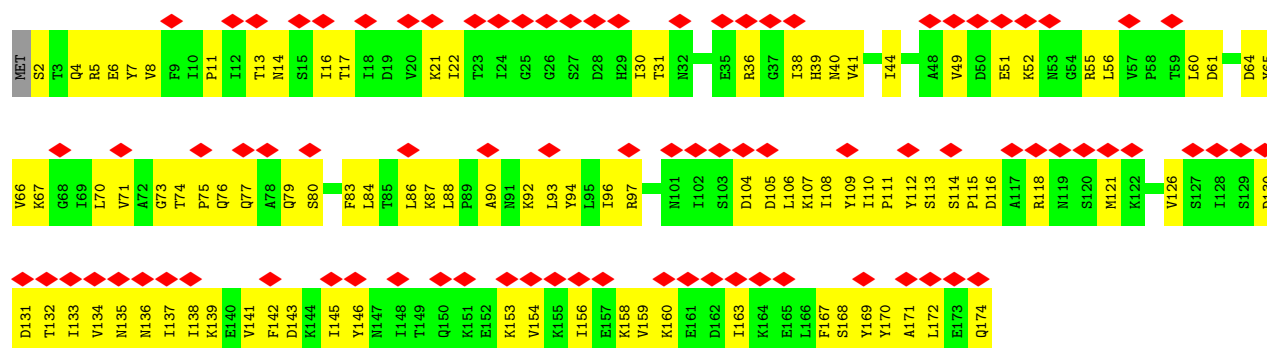




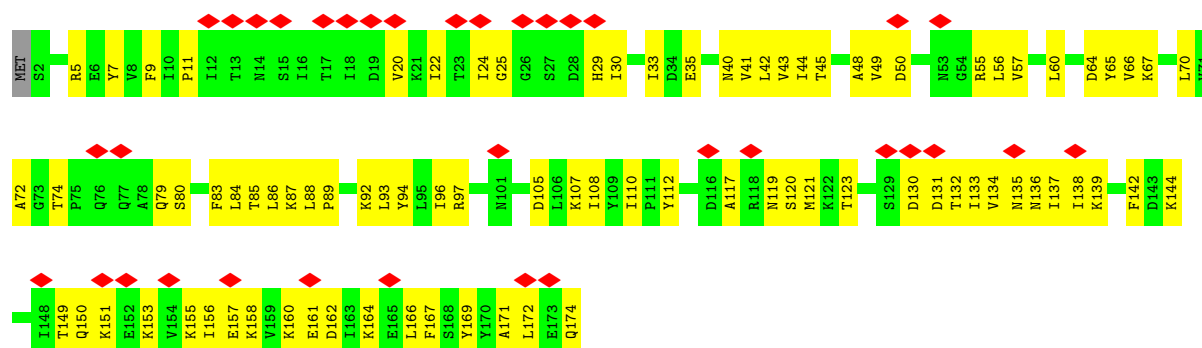
• Molecule 9: CRISPR-associated protein CmrX



• Molecule 9: CRISPR-associated protein CmrX



• Molecule 9: CRISPR-associated protein CmrX



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	51280	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 ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	1700	Depositor
Maximum defocus (nm)	2600	Depositor
Magnification	96000	Depositor
Image detector	FEI FALCON III (4k x 4k)	Depositor
Maximum map value	29.254	Depositor
Minimum map value	-13.609	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.688	Depositor
Recommended contour level	2.1	Depositor
Map size (Å)	416.0, 416.0, 416.0	wwPDB
Map dimensions	500, 500, 500	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.832, 0.832, 0.832	Depositor

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: MN, ANP, ZN

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.31	0/1275	0.41	0/1716
1	B	0.29	0/1257	0.62	1/1694 (0.1%)
1	C	0.33	0/1283	0.45	0/1727
2	D	0.38	0/2318	0.45	0/3134
2	E	0.39	0/2318	0.45	0/3134
2	F	0.41	0/2318	0.44	0/3134
2	G	0.36	0/2302	0.46	0/3113
3	H	0.28	0/2574	0.42	0/3469
4	I	0.35	0/2325	0.44	0/3138
5	J	0.27	0/3899	0.42	0/5242
6	K	0.23	0/8580	0.45	1/11588 (0.0%)
7	U	0.31	0/844	0.37	0/1310
8	V	0.47	0/1149	0.44	0/1789
9	L	0.25	0/1400	0.45	0/1895
9	M	0.22	0/1400	0.42	0/1895
9	N	0.19	0/1400	0.46	0/1895
9	O	0.18	0/1400	0.45	0/1895
9	P	0.18	0/1400	0.46	0/1895
9	Q	0.19	0/1400	0.50	0/1895
9	R	0.17	0/1400	0.47	0/1895
9	S	0.19	0/1400	0.54	1/1895 (0.1%)
9	T	0.17	0/1400	0.48	0/1895
9	W	0.18	0/1400	0.51	0/1895
9	X	0.18	0/1400	0.53	0/1895
9	l	0.25	0/1400	0.52	1/1895 (0.1%)
9	m	0.27	0/1400	0.52	0/1895
9	n	0.18	0/1400	0.45	0/1895
9	o	0.21	0/1400	0.47	0/1895
9	p	0.18	0/1400	0.45	0/1895
9	q	0.18	0/1400	0.46	0/1895
9	r	0.19	0/1400	0.45	0/1895
9	s	0.17	0/1400	0.51	1/1895 (0.1%)

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
9	t	0.19	0/1400	0.52	0/1895
9	w	0.17	0/1400	0.48	0/1895
9	x	0.19	0/1400	0.55	1/1895 (0.1%)
All	All	0.27	0/63242	0.47	6/85878 (0.0%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
6	K	0	1
9	N	0	1
9	O	0	1
9	Q	0	1
9	m	0	2
9	p	0	1
All	All	0	7

There are no bond length outliers.

The worst 5 of 6 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	s	148	ILE	N-CA-C	-8.52	105.19	113.53
9	l	154	VAL	N-CA-C	-7.48	106.20	113.53
1	B	61	SER	N-CA-C	-6.89	106.07	114.75
6	K	990	SER	N-CA-C	6.46	124.57	110.80
9	S	54	GLY	N-CA-C	6.27	119.81	111.03

There are no chirality outliers.

5 of 7 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
6	K	989	ASN	Peptide
9	N	79	GLN	Peptide
9	O	78	ALA	Peptide
9	Q	77	GLN	Peptide
9	m	15	SER	Peptide

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1253	0	1287	45	0
1	B	1235	0	1274	90	0
1	C	1261	0	1298	38	0
2	D	2274	0	2343	54	0
2	E	2274	0	2342	65	0
2	F	2274	0	2342	72	0
2	G	2258	0	2322	79	0
3	H	2528	0	2581	113	0
4	I	2282	0	2331	66	0
5	J	3833	0	3917	115	0
6	K	8408	0	8554	350	0
7	U	758	0	386	34	0
8	V	1025	0	520	43	0
9	L	1378	0	1418	59	0
9	M	1378	0	1418	61	0
9	N	1378	0	1417	70	0
9	O	1378	0	1418	57	0
9	P	1378	0	1418	71	0
9	Q	1378	0	1417	75	0
9	R	1378	0	1417	74	0
9	S	1378	0	1417	67	0
9	T	1378	0	1418	75	0
9	W	1378	0	1418	58	0
9	X	1378	0	1418	82	0
9	l	1378	0	1417	68	0
9	m	1378	0	1417	55	0
9	n	1378	0	1418	68	0
9	o	1378	0	1418	63	0
9	p	1378	0	1417	73	0
9	q	1378	0	1418	90	0
9	r	1378	0	1418	78	0
9	s	1378	0	1418	72	0
9	t	1378	0	1417	88	0
9	w	1378	0	1417	79	0
9	x	1378	0	1417	84	0
10	K	1	0	0	0	0
11	K	62	0	26	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
12	K	3	0	0	0	0
All	All	62045	0	62709	2495	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 20.

The worst 5 of 2495 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)
9:W:140:GLU:HG3	9:W:141:VAL:HG23	1.50	0.94
1:B:47:ILE:HG13	1:B:48:PRO:HD3	1.52	0.90
9:Q:136:ASN:O	9:Q:140:GLU:HB2	1.74	0.87
9:r:11:PRO:HG2	9:r:16:ILE:HG12	1.58	0.86
9:t:10:ILE:HD11	9:t:69:ILE:HG23	1.58	0.85

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	151/155 (97%)	139 (92%)	11 (7%)	1 (1%)	18	37
1	B	149/155 (96%)	122 (82%)	27 (18%)	0	100	100
1	C	152/155 (98%)	144 (95%)	8 (5%)	0	100	100
2	D	283/286 (99%)	265 (94%)	18 (6%)	0	100	100
2	E	283/286 (99%)	262 (93%)	21 (7%)	0	100	100
2	F	283/286 (99%)	265 (94%)	18 (6%)	0	100	100
2	G	281/286 (98%)	247 (88%)	34 (12%)	0	100	100
3	H	310/313 (99%)	283 (91%)	27 (9%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
4	I	282/296 (95%)	264 (94%)	18 (6%)	0	100	100
5	J	463/476 (97%)	415 (90%)	47 (10%)	1 (0%)	43	65
6	K	1019/1037 (98%)	916 (90%)	102 (10%)	1 (0%)	48	71
9	L	171/174 (98%)	143 (84%)	28 (16%)	0	100	100
9	M	171/174 (98%)	143 (84%)	28 (16%)	0	100	100
9	N	171/174 (98%)	139 (81%)	32 (19%)	0	100	100
9	O	171/174 (98%)	137 (80%)	34 (20%)	0	100	100
9	P	171/174 (98%)	141 (82%)	29 (17%)	1 (1%)	21	41
9	Q	171/174 (98%)	134 (78%)	37 (22%)	0	100	100
9	R	171/174 (98%)	145 (85%)	26 (15%)	0	100	100
9	S	171/174 (98%)	134 (78%)	36 (21%)	1 (1%)	21	41
9	T	171/174 (98%)	138 (81%)	33 (19%)	0	100	100
9	W	171/174 (98%)	137 (80%)	34 (20%)	0	100	100
9	X	171/174 (98%)	135 (79%)	36 (21%)	0	100	100
9	l	171/174 (98%)	144 (84%)	26 (15%)	1 (1%)	21	41
9	m	171/174 (98%)	138 (81%)	33 (19%)	0	100	100
9	n	171/174 (98%)	143 (84%)	26 (15%)	2 (1%)	10	24
9	o	171/174 (98%)	134 (78%)	37 (22%)	0	100	100
9	p	171/174 (98%)	145 (85%)	25 (15%)	1 (1%)	21	41
9	q	171/174 (98%)	141 (82%)	29 (17%)	1 (1%)	21	41
9	r	171/174 (98%)	146 (85%)	25 (15%)	0	100	100
9	s	171/174 (98%)	141 (82%)	29 (17%)	1 (1%)	21	41
9	t	171/174 (98%)	141 (82%)	30 (18%)	0	100	100
9	w	171/174 (98%)	138 (81%)	32 (19%)	1 (1%)	21	41
9	x	171/174 (98%)	142 (83%)	29 (17%)	0	100	100
All	All	7418/7559 (98%)	6401 (86%)	1005 (14%)	12 (0%)	44	65

5 of 12 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
6	K	991	LEU
9	P	131	ASP
9	n	131	ASP

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Mol	Chain	Res	Type
9	q	76	GLN
9	s	35	GLU

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	134/136 (98%)	134 (100%)	0	100	100
1	B	133/136 (98%)	132 (99%)	1 (1%)	73	87
1	C	135/136 (99%)	135 (100%)	0	100	100
2	D	252/253 (100%)	252 (100%)	0	100	100
2	E	252/253 (100%)	252 (100%)	0	100	100
2	F	252/253 (100%)	252 (100%)	0	100	100
2	G	250/253 (99%)	250 (100%)	0	100	100
3	H	279/280 (100%)	279 (100%)	0	100	100
4	I	255/266 (96%)	254 (100%)	1 (0%)	84	93
5	J	438/445 (98%)	438 (100%)	0	100	100
6	K	936/949 (99%)	936 (100%)	0	100	100
9	L	159/160 (99%)	159 (100%)	0	100	100
9	M	159/160 (99%)	159 (100%)	0	100	100
9	N	159/160 (99%)	159 (100%)	0	100	100
9	O	159/160 (99%)	158 (99%)	1 (1%)	78	90
9	P	159/160 (99%)	158 (99%)	1 (1%)	78	90
9	Q	159/160 (99%)	159 (100%)	0	100	100
9	R	159/160 (99%)	159 (100%)	0	100	100
9	S	159/160 (99%)	158 (99%)	1 (1%)	78	90
9	T	159/160 (99%)	159 (100%)	0	100	100
9	W	159/160 (99%)	159 (100%)	0	100	100
9	X	159/160 (99%)	159 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
9	l	159/160 (99%)	159 (100%)	0	100	100
9	m	159/160 (99%)	159 (100%)	0	100	100
9	n	159/160 (99%)	159 (100%)	0	100	100
9	o	159/160 (99%)	159 (100%)	0	100	100
9	p	159/160 (99%)	158 (99%)	1 (1%)	78	90
9	q	159/160 (99%)	159 (100%)	0	100	100
9	r	159/160 (99%)	159 (100%)	0	100	100
9	s	159/160 (99%)	159 (100%)	0	100	100
9	t	159/160 (99%)	159 (100%)	0	100	100
9	w	159/160 (99%)	159 (100%)	0	100	100
9	x	159/160 (99%)	159 (100%)	0	100	100
All	All	6814/6880 (99%)	6808 (100%)	6 (0%)	87	96

5 of 6 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
9	P	163	ILE
9	S	163	ILE
9	p	156	ILE
4	I	271	TYR
1	B	87	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 104 such sidechains are listed below:

Mol	Chain	Res	Type
9	Q	101	ASN
9	m	14	ASN
9	x	77	GLN
9	S	76	GLN
9	T	79	GLN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
7	U	34/46 (73%)	11 (32%)	1 (2%)
8	V	47/51 (92%)	12 (25%)	0

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Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
All	All	81/97 (83%)	23 (28%)	1 (1%)

5 of 23 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
7	U	6	A
7	U	7	G
7	U	12	G
7	U	14	U
7	U	18	C

All (1) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
7	U	5	A

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](#)

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 for Mogul analysis.

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	$\# Z > 2$
11	ANP	K	1102	12	33,33,33	1.06	4 (12%)	45,52,52	0.61	1 (2%)
11	ANP	K	1103	12	33,33,33	0.93	4 (12%)	45,52,52	0.61	0

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
11	ANP	K	1102	12	-	5/18/38/38	0/3/3/3
11	ANP	K	1103	12	-	3/18/38/38	0/3/3/3

The worst 5 of 8 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
11	K	1102	ANP	PB-O3A	-3.46	1.54	1.59
11	K	1103	ANP	PB-O3A	-2.97	1.55	1.59
11	K	1102	ANP	PG-N3B	2.47	1.69	1.63
11	K	1102	ANP	PA-O3A	-2.22	1.57	1.59
11	K	1103	ANP	PG-N3B	2.20	1.69	1.63

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
11	K	1102	ANP	O2G-PG-O1G	-2.06	108.28	113.45

There are no chirality outliers.

5 of 8 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
11	K	1102	ANP	C5'-O5'-PA-O1A
11	K	1103	ANP	PG-N3B-PB-O1B
11	K	1102	ANP	PB-O3A-PA-O1A
11	K	1102	ANP	C4'-C5'-O5'-PA
11	K	1102	ANP	C2'-C1'-N9-C8

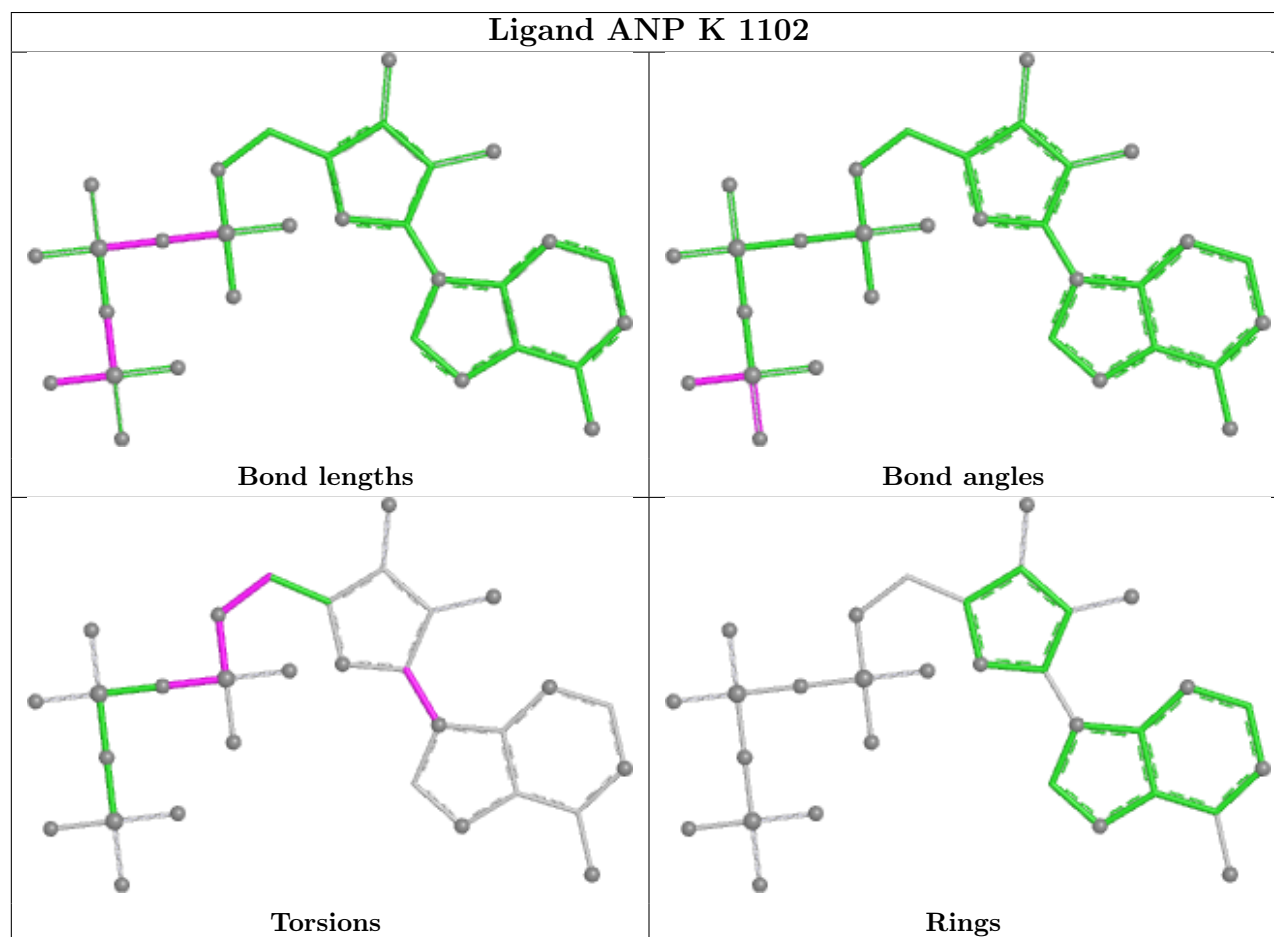
There are no ring outliers.

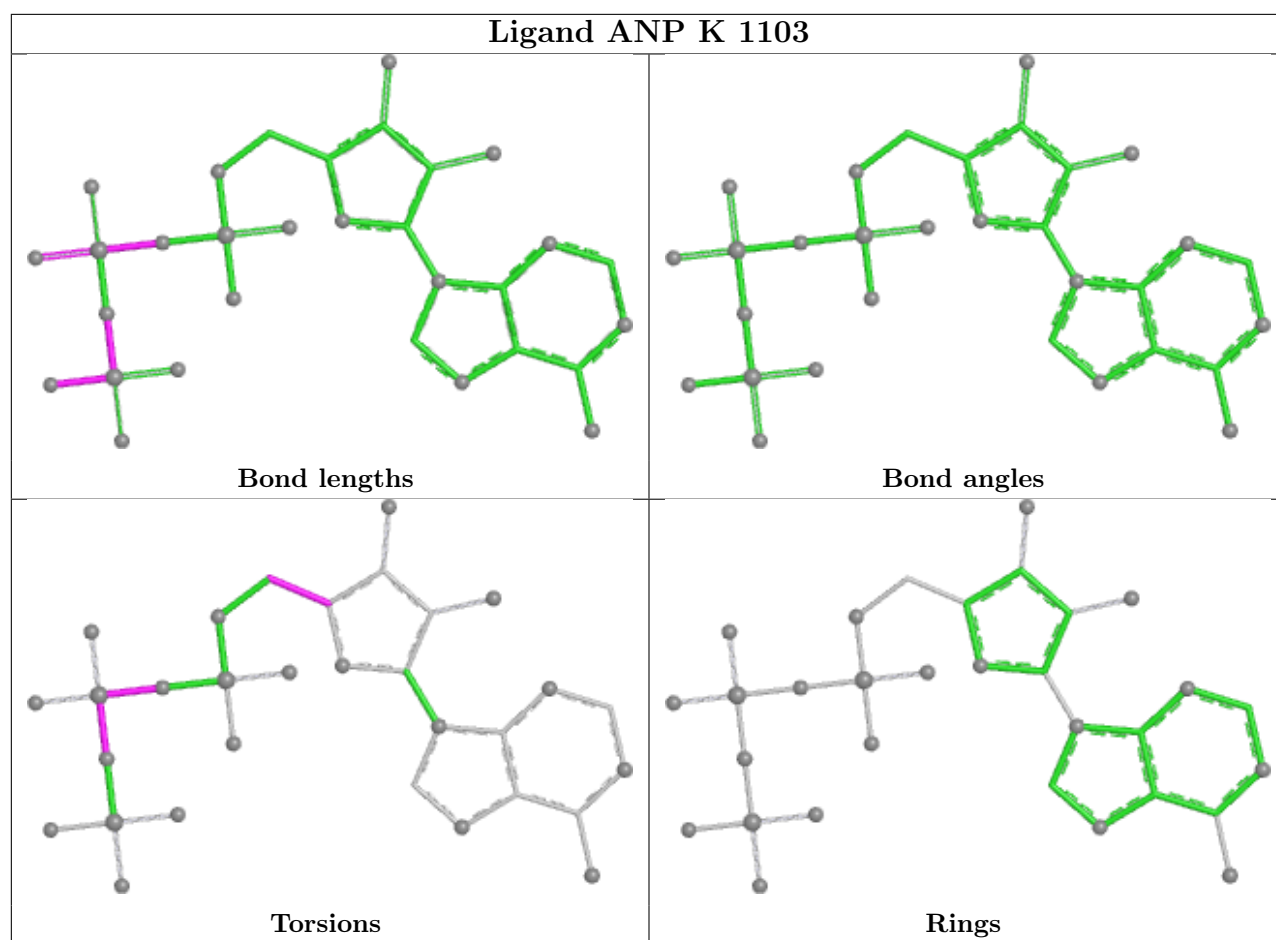
2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
11	K	1102	ANP	1	0
11	K	1103	ANP	4	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.

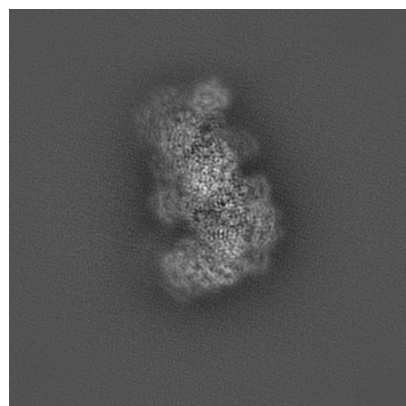
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-10126. These allow visual inspection of the internal detail of the map and identification of artifacts.

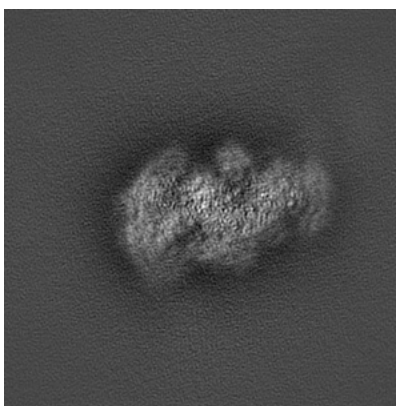
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

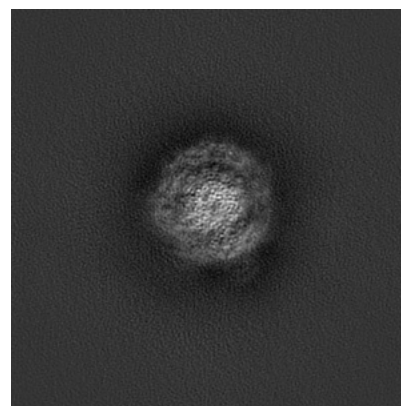
6.1.1 Primary map



X

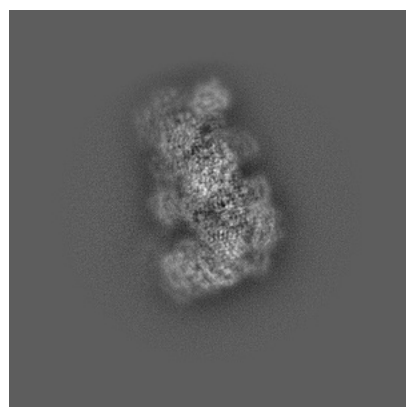


Y

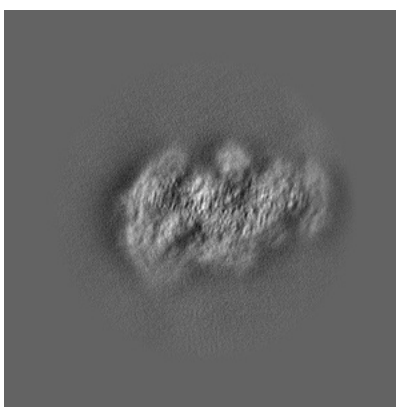


Z

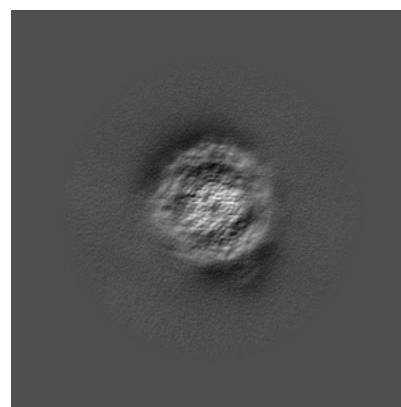
6.1.2 Raw map



X



Y

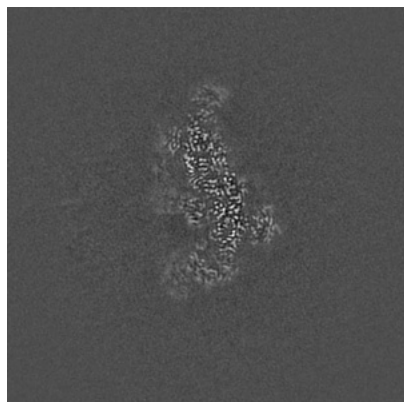


Z

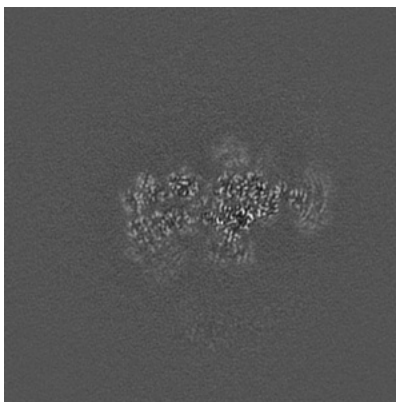
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

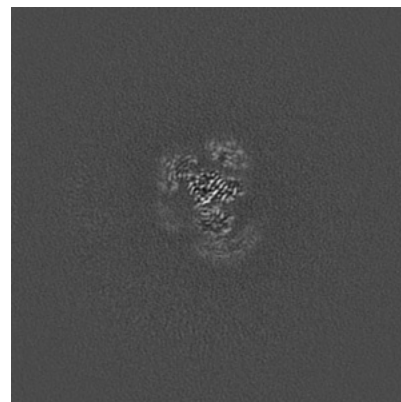
6.2.1 Primary map



X Index: 250

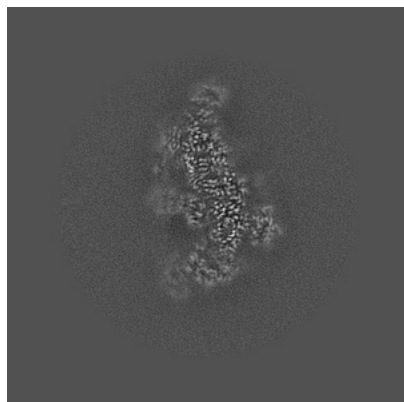


Y Index: 250

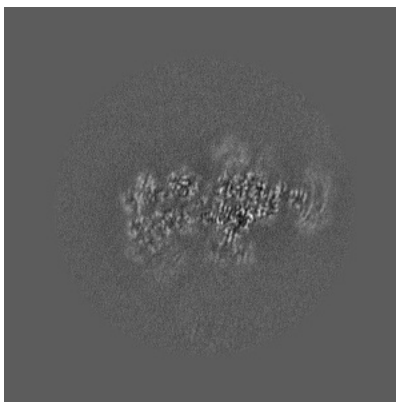


Z Index: 250

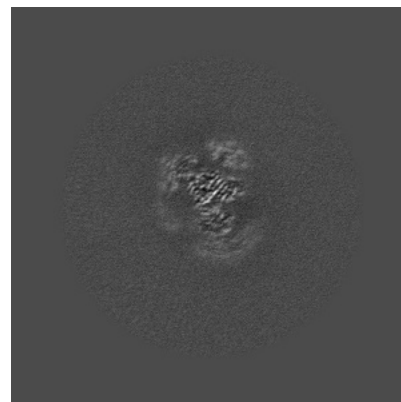
6.2.2 Raw map



X Index: 250



Y Index: 250

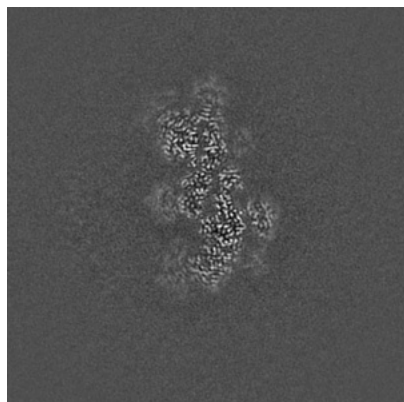


Z Index: 250

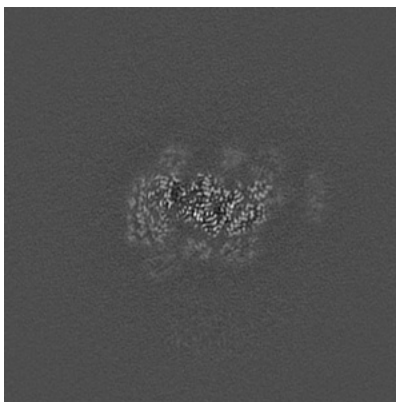
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

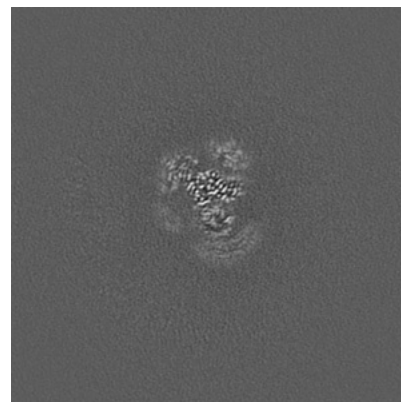
6.3.1 Primary map



X Index: 267

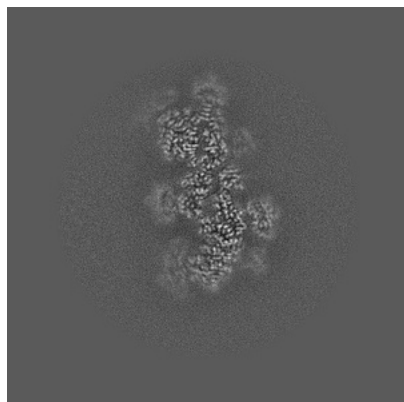


Y Index: 272

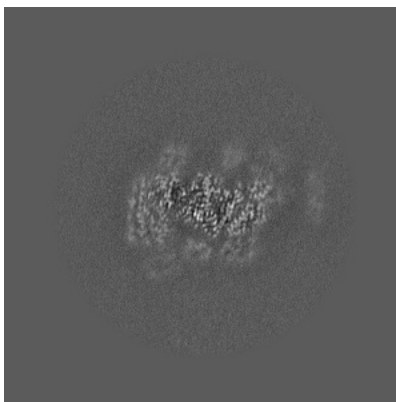


Z Index: 251

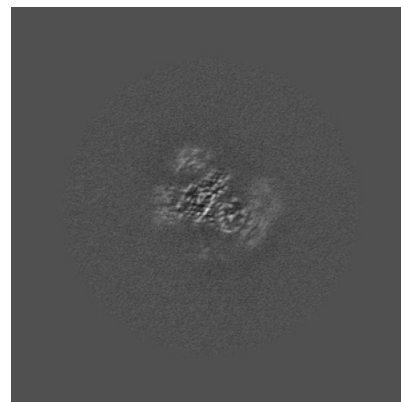
6.3.2 Raw map



X Index: 267



Y Index: 272

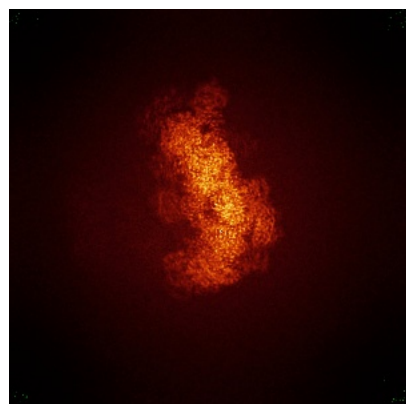


Z Index: 281

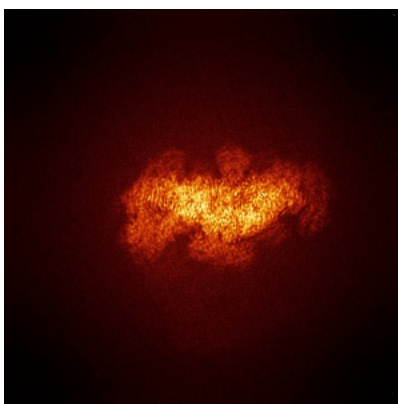
The images above show the largest variance slices of the map in three orthogonal directions.

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

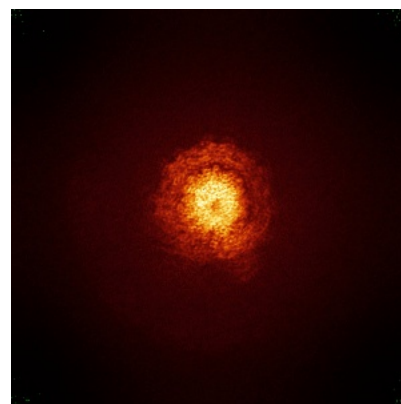
6.4.1 Primary map



X

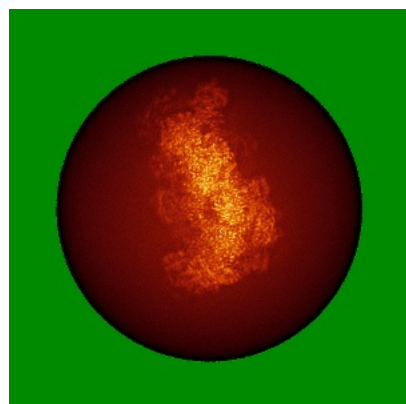


Y

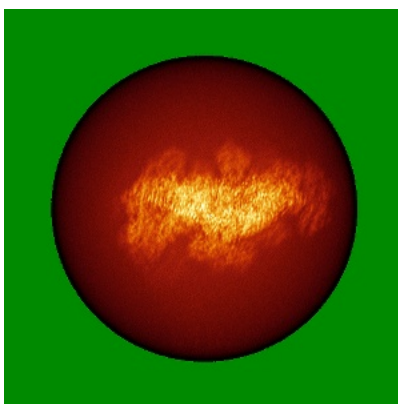


Z

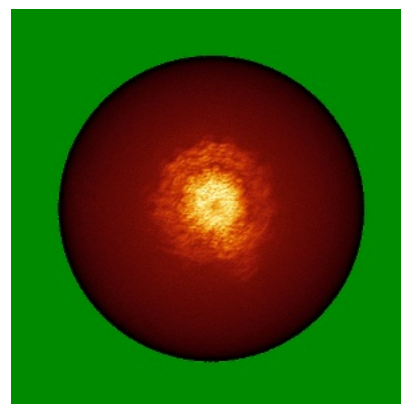
6.4.2 Raw map



X



Y

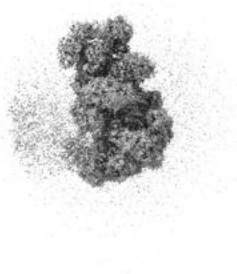


Z

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](#)

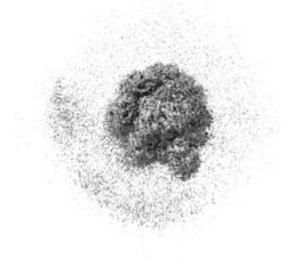
6.5.1 Primary map



X



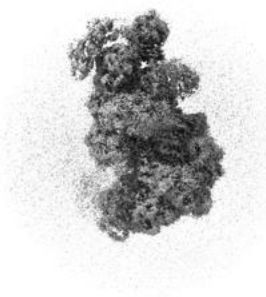
Y



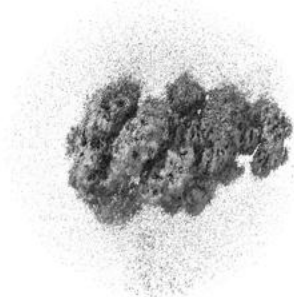
Z

The images above show the 3D surface view of the map at the recommended contour level 2.1. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



X



Y



Z

These images show the 3D surface of the raw map. The raw map's contour level was selected so that its surface encloses the same volume as the primary map does at its recommended contour level.

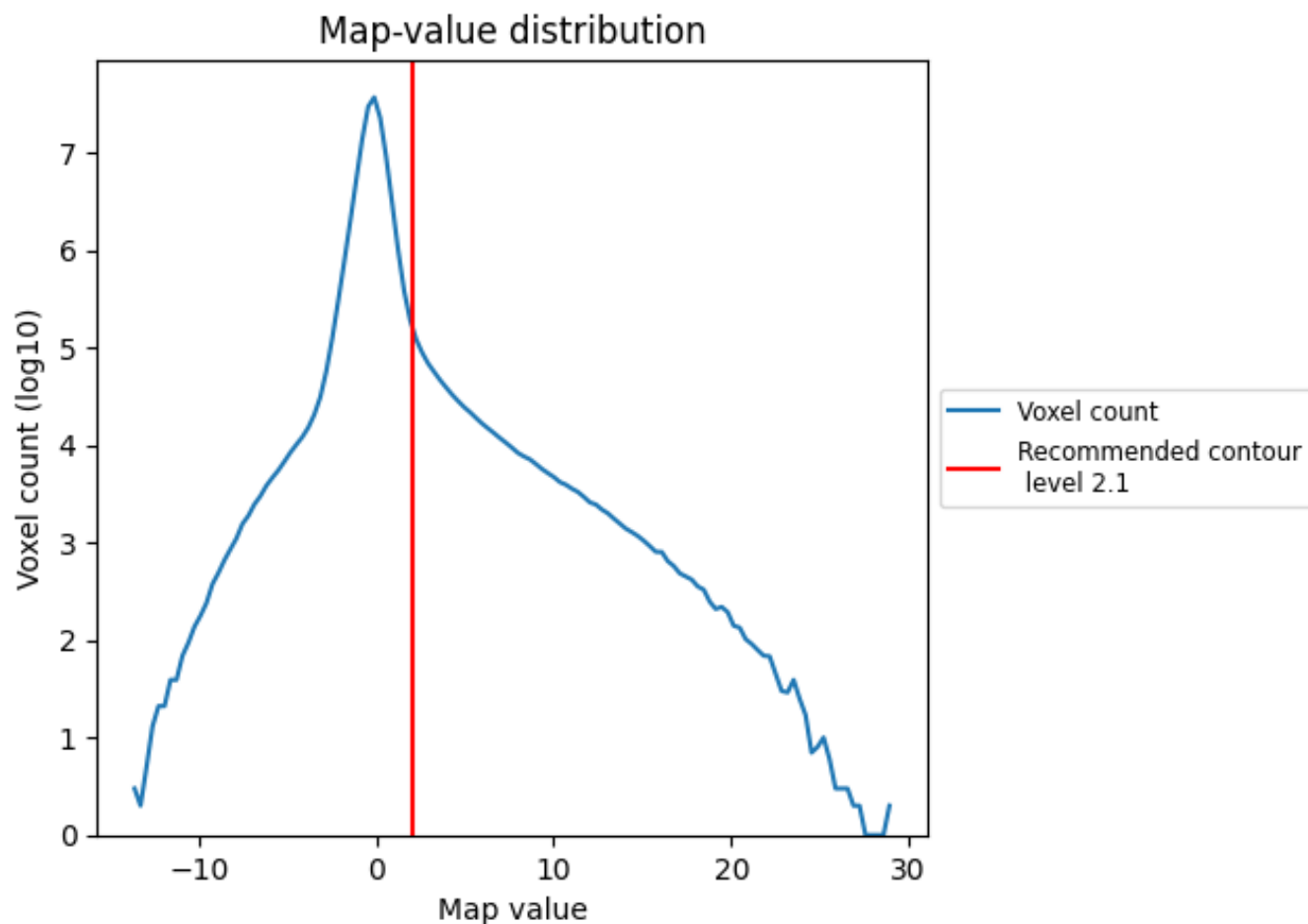
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

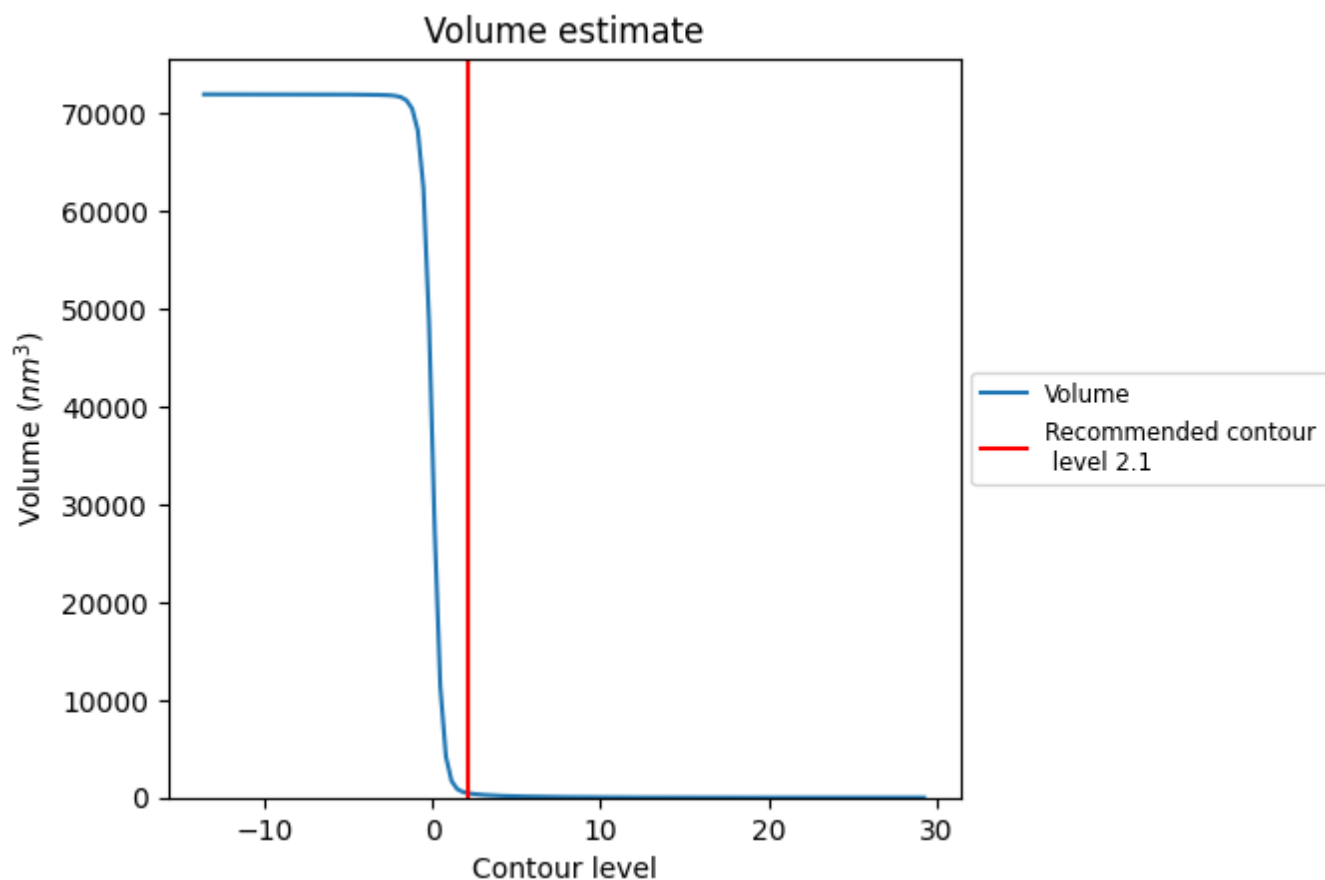
This section contains the results of statistical analysis of the map.

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.

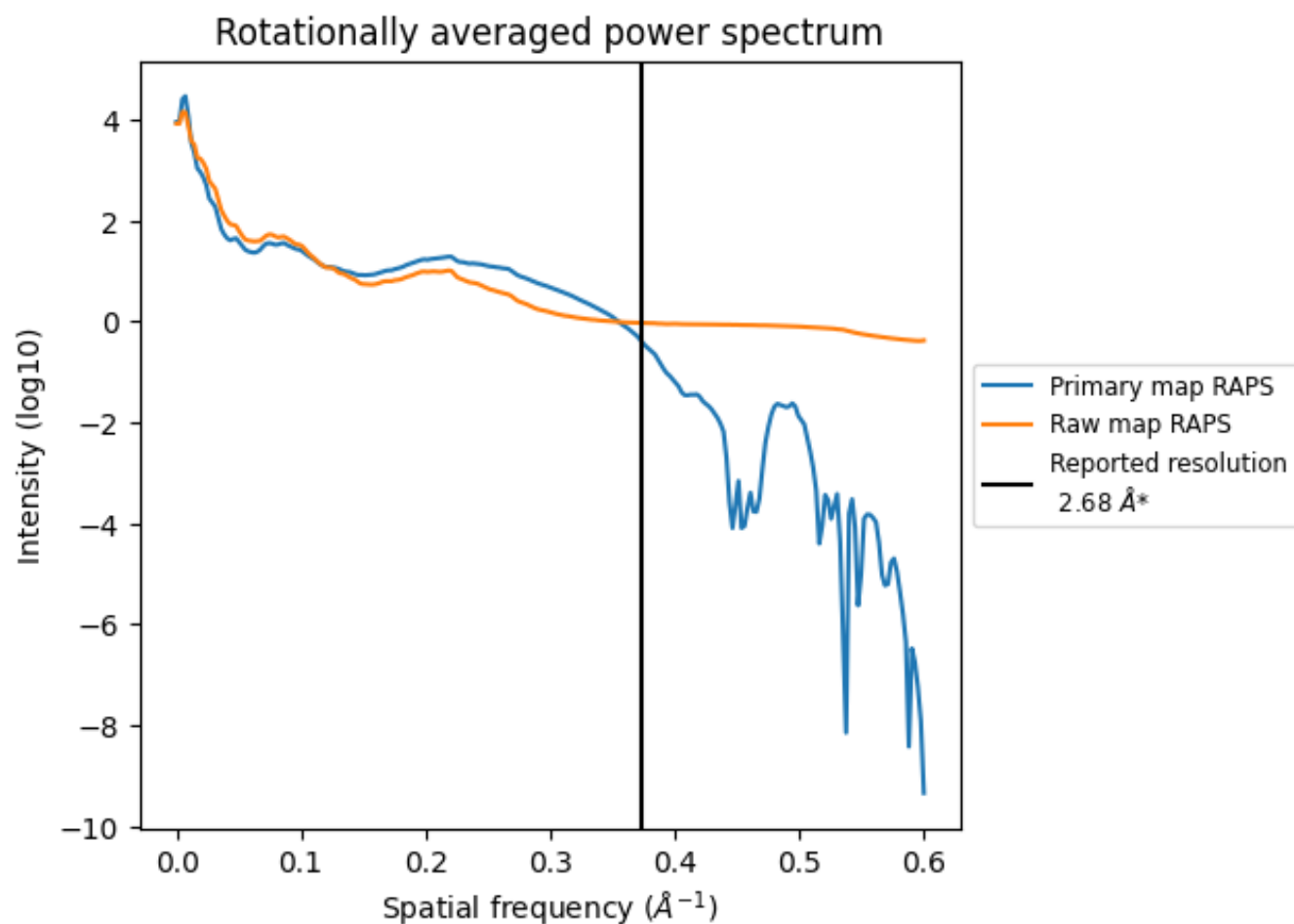
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 454 nm^3 ; this corresponds to an approximate mass of 410 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 ⓘ

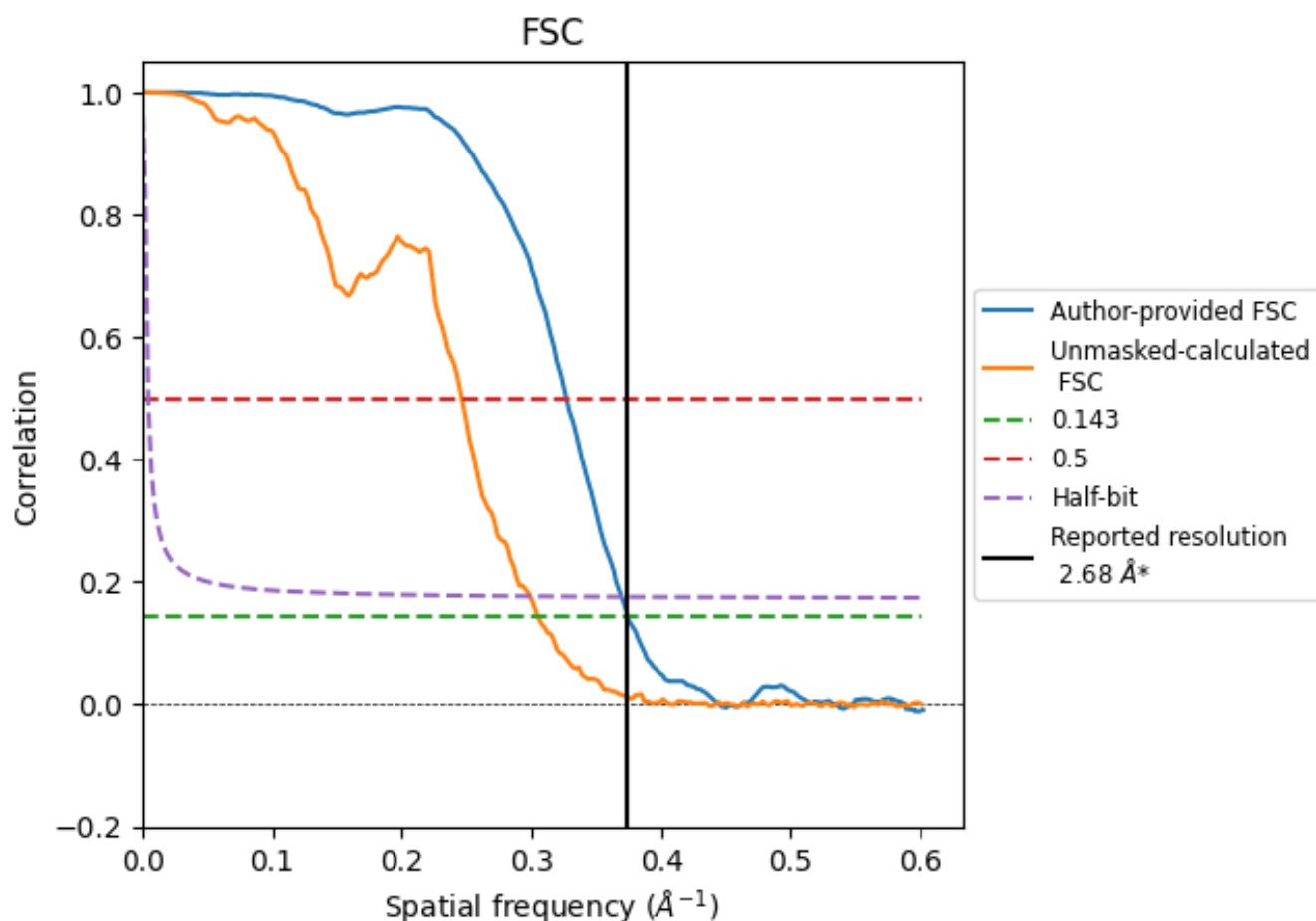


*Reported resolution corresponds to spatial frequency of 0.373 Å⁻¹

8 Fourier-Shell correlation [i](#)

Fourier-Shell Correlation (FSC) is the most commonly used method to estimate the resolution of single-particle and subtomogram-averaged maps. The shape of the curve depends on the imposed symmetry, mask and whether or not the two 3D reconstructions used were processed from a common reference. The reported resolution is shown as a black line. A curve is displayed for the half-bit criterion in addition to lines showing the 0.143 gold standard cut-off and 0.5 cut-off.

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of $0.373 \text{ \AA}^{-1}$

8.2 Resolution estimates [i](#)

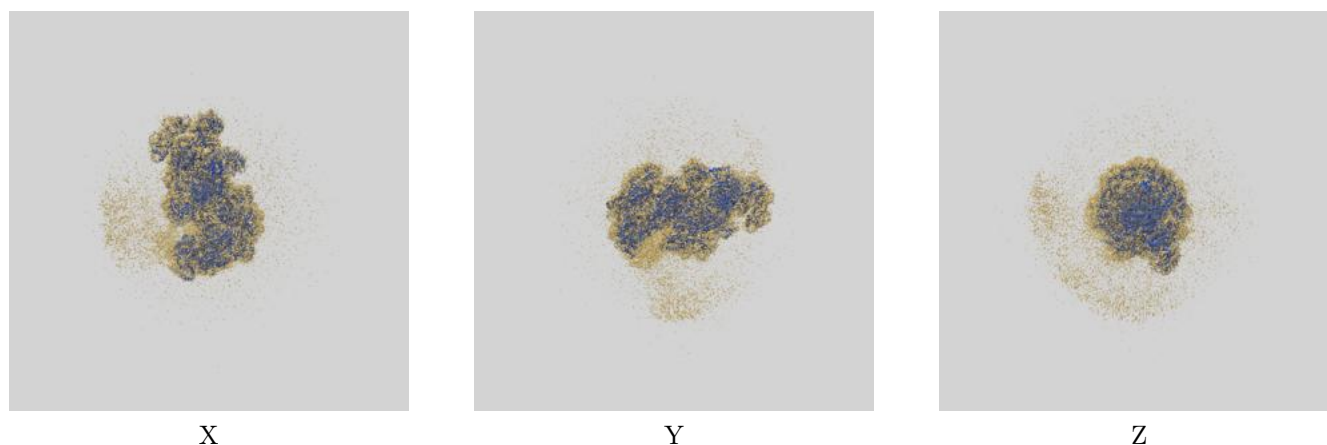
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.68	-	-
Author-provided FSC curve	2.67	3.06	2.71
Unmasked-calculated*	3.28	4.06	3.34

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 3.28 differs from the reported value 2.68 by more than 10 %

9 Map-model fit [i](#)

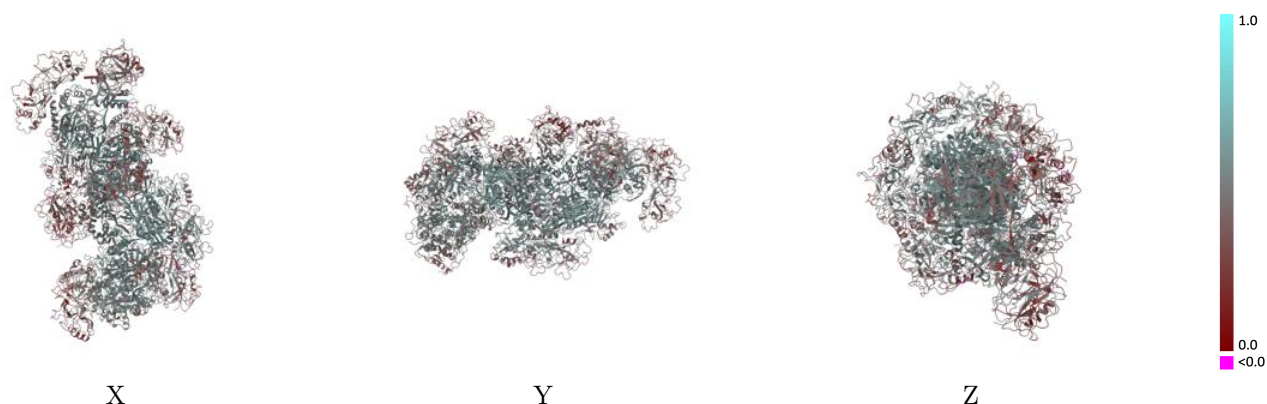
This section contains information regarding the fit between EMDB map EMD-10126 and PDB model 6S91. Per-residue inclusion information can be found in section [3](#) on page [10](#).

9.1 Map-model overlay [i](#)



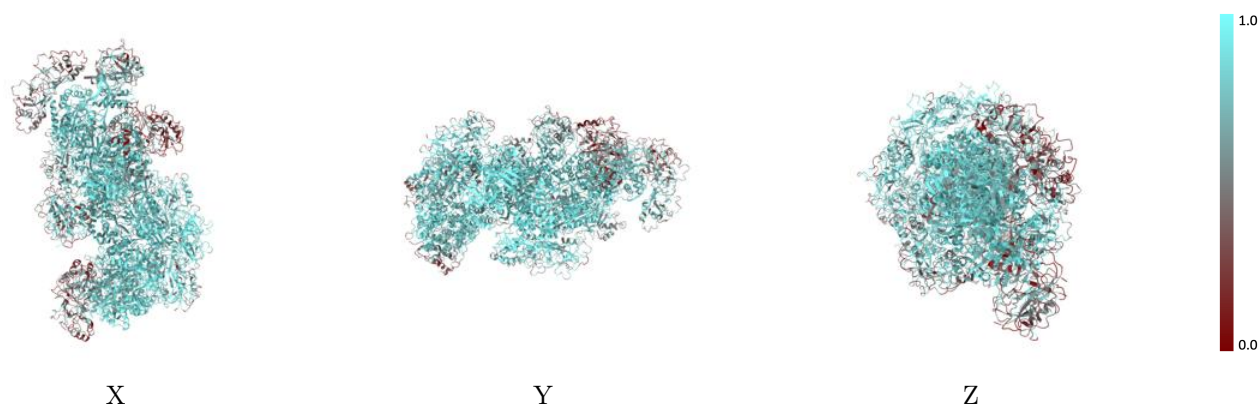
The images above show the 3D surface view of the map at the recommended contour level 2.1 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



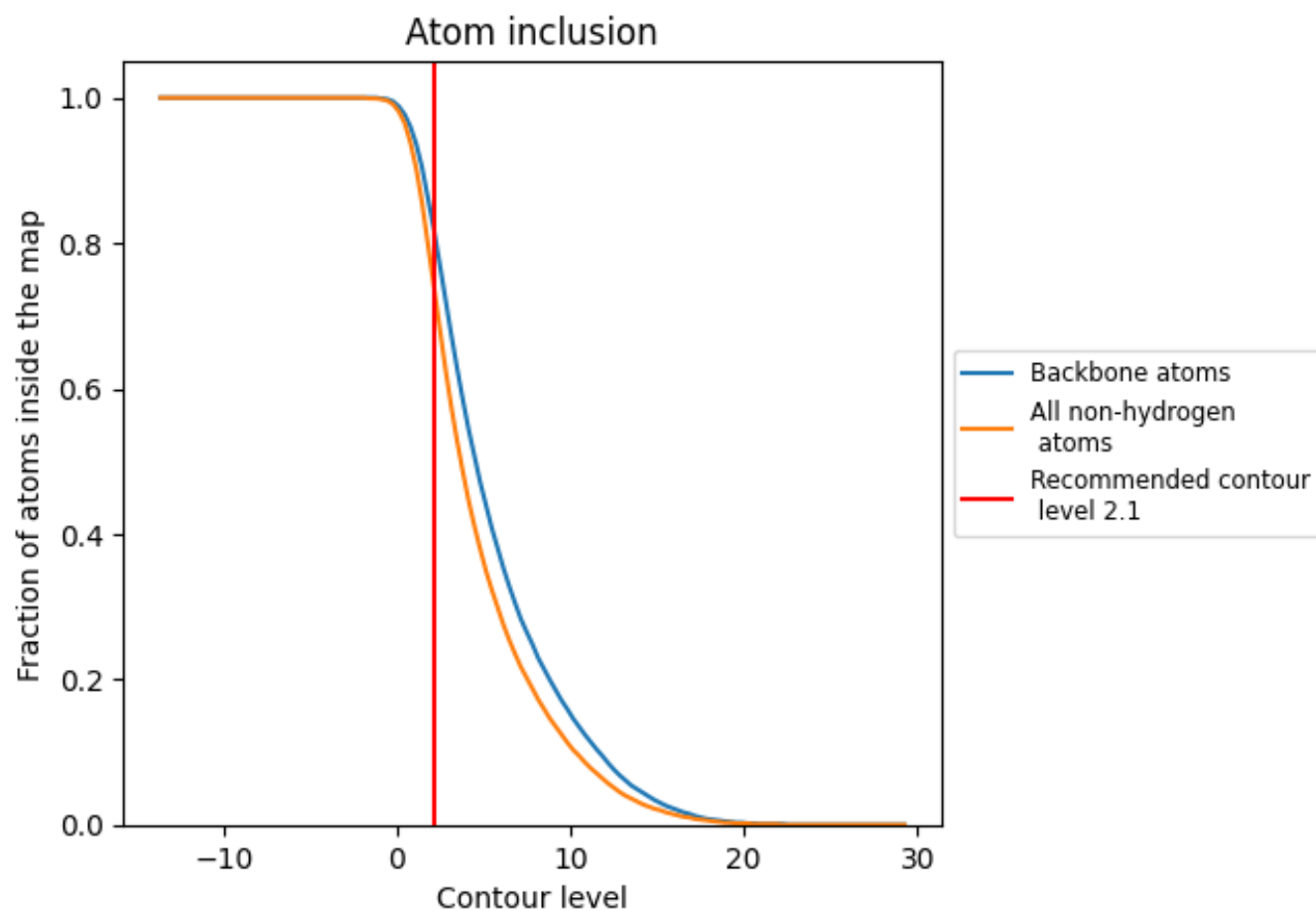
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (2.1).

9.4 Atom inclusion [i](#)



At the recommended contour level, 82% of all backbone atoms, 75% 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 (2.1) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.7480	 0.4630
A	 0.9020	 0.5410
B	 0.8340	 0.4550
C	 0.9030	 0.5560
D	 0.9320	 0.5720
E	 0.9210	 0.5690
F	 0.9220	 0.5700
G	 0.9170	 0.5500
H	 0.8990	 0.5190
I	 0.9280	 0.5630
J	 0.8620	 0.5130
K	 0.7990	 0.4790
L	 0.8260	 0.4670
M	 0.8710	 0.4660
N	 0.7110	 0.4150
O	 0.7320	 0.4310
P	 0.6280	 0.4370
Q	 0.5840	 0.3450
R	 0.6230	 0.3360
S	 0.3250	 0.3670
T	 0.6470	 0.3870
U	 0.8810	 0.5280
V	 0.9700	 0.5810
W	 0.4180	 0.3830
X	 0.2620	 0.3190
l	 0.8430	 0.4610
m	 0.8680	 0.4680
n	 0.6920	 0.3770
o	 0.7510	 0.4380
p	 0.6600	 0.4370
q	 0.6130	 0.3450
r	 0.7480	 0.3740
s	 0.2940	 0.3720
t	 0.5400	 0.3810
w	 0.4580	 0.3750
x	 0.4030	 0.3440

