



Full wwPDB EM Validation Report ⓘ

Nov 19, 2022 – 08:19 pm GMT

PDB ID : 6EMW
EMDB ID : EMD-3897
Title : Structure of S.aureus ClpC in complex with MecA
Authors : Carroni, M.; Mogk, A.; Bukau, B.; Franke, K.
Deposited on : 2017-10-03
Resolution : 11.00 Å (reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

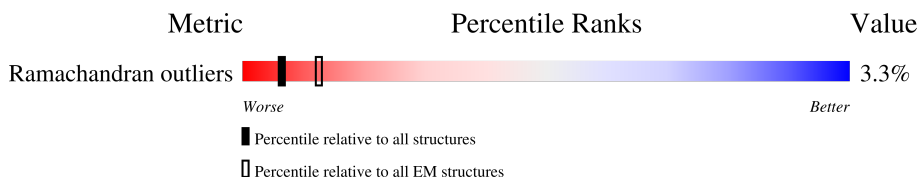
EMDB validation analysis : 0.0.1.dev43
MolProbity : 4.02b-467
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ : 1.9.9
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.31.2

1 Overall quality at a glance

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

The reported resolution of this entry is 11.00 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)
Ramachandran outliers	154571	4023

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	82	96% .
1	G	82	96% .
1	M	82	96% .
1	S	82	96% .
1	Y	82	96% .
1	k	82	96% .
2	B	225	89% .. 7%
2	H	225	89% .. 7%
2	N	225	89% .. 7%
2	T	225	89% .. 7%
2	Z	225	89% .. 7%

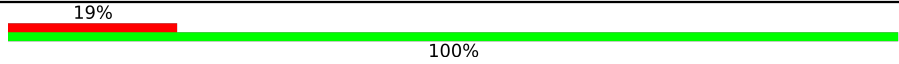
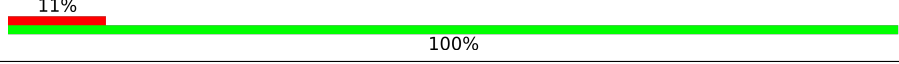
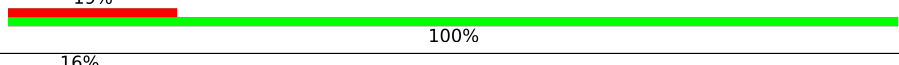
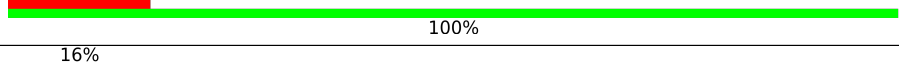
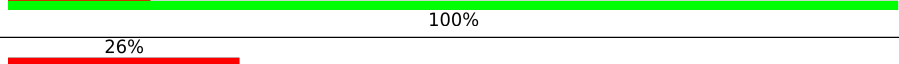
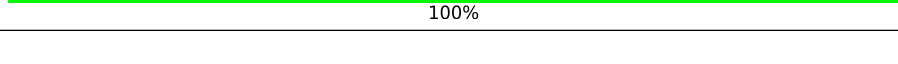
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Mol	Chain	Length	Quality of chain
2	l	225	8% 89% 7%
3	C	55	96%
3	I	55	96%
3	O	55	40% 96%
3	U	55	18% 96%
3	a	55	96%
3	m	55	13% 96%
4	D	145	6% 61% 38%
4	J	145	6% 61% 38%
4	P	145	6% 60% 38%
4	V	145	6% 61% 38%
4	b	145	6% 61% 38%
4	n	145	6% 59% 38%
5	E	181	8% 89% 8%
5	K	181	21% 89% 8%
5	Q	181	8% 89% 8%
5	W	181	8% 89% 8%
5	c	181	44% 89% 8%
5	o	181	18% 89% 8%
6	F	157	15% 92% 5%
6	L	157	7% 92% 5%
6	R	157	7% 92% 5%
6	X	157	11% 92% 5%
6	d	157	8% 92% 5%
6	p	157	8% 92% 5%

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Mol	Chain	Length	Quality of chain
7	e	90	 19% 100%
7	f	90	 11% 100%
7	g	90	 19% 100%
7	h	90	 16% 100%
7	i	90	 16% 100%
7	j	90	 26% 100%

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 15210 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 ATP-dependent Clp protease ATP-binding subunit.

Mol	Chain	Residues	Atoms			AltConf	Trace
			Total	C	N		
1	A	82	246	164	82	0	0
1	G	82	246	164	82	0	0
1	M	82	246	164	82	0	0
1	S	82	246	164	82	0	0
1	Y	82	246	164	82	0	0
1	k	82	246	164	82	0	0

- Molecule 2 is a protein called ATP-dependent Clp protease ATP-binding subunit.

Mol	Chain	Residues	Atoms			AltConf	Trace
			Total	C	N		
2	B	210	630	420	210	0	0
2	H	210	630	420	210	0	0
2	N	210	630	420	210	0	0
2	T	210	630	420	210	0	0
2	Z	210	630	420	210	0	0
2	l	210	630	420	210	0	0

- Molecule 3 is a protein called Class III stress response-related ATPase, AAA+ superfamily.

Mol	Chain	Residues	Atoms			AltConf	Trace
			Total	C	N		
3	C	55	165	110	55	0	0

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Mol	Chain	Residues	Atoms			AltConf	Trace
3	I	55	Total 165	C 110	N 55	0	0
3	O	55	Total 165	C 110	N 55	0	0
3	U	55	Total 165	C 110	N 55	0	0
3	a	55	Total 165	C 110	N 55	0	0
3	m	55	Total 165	C 110	N 55	0	0

- Molecule 4 is a protein called ATP-dependent Clp protease ATP-binding subunit ClpC.

Mol	Chain	Residues	Atoms			AltConf	Trace
4	D	90	Total 270	C 180	N 90	0	0
4	J	90	Total 270	C 180	N 90	0	0
4	P	90	Total 270	C 180	N 90	0	0
4	V	90	Total 270	C 180	N 90	0	0
4	b	90	Total 270	C 180	N 90	0	0
4	n	90	Total 270	C 180	N 90	0	0

- Molecule 5 is a protein called ATP-dependent Clp protease ATP-binding subunit ClpC.

Mol	Chain	Residues	Atoms			AltConf	Trace
5	E	166	Total 498	C 332	N 166	0	0
5	K	166	Total 498	C 332	N 166	0	0
5	Q	166	Total 498	C 332	N 166	0	0
5	W	166	Total 498	C 332	N 166	0	0
5	c	166	Total 498	C 332	N 166	0	0
5	o	166	Total 498	C 332	N 166	0	0

- Molecule 6 is a protein called ATP-dependent Clp protease ATP-binding subunit ClpC.

Mol	Chain	Residues	Atoms			AltConf	Trace
6	F	152	Total 456	C 304	N 152	0	0
6	L	152	Total 456	C 304	N 152	0	0
6	R	152	Total 456	C 304	N 152	0	0
6	X	152	Total 456	C 304	N 152	0	0
6	d	152	Total 456	C 304	N 152	0	0
6	p	152	Total 456	C 304	N 152	0	0

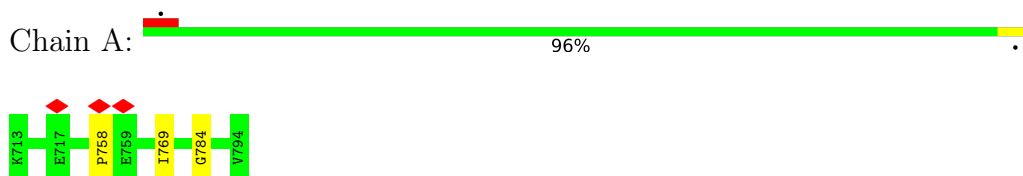
- Molecule 7 is a protein called Adapter protein MecA.

Mol	Chain	Residues	Atoms			AltConf	Trace
7	e	90	Total 270	C 180	N 90	0	0
7	f	90	Total 270	C 180	N 90	0	0
7	g	90	Total 270	C 180	N 90	0	0
7	h	90	Total 270	C 180	N 90	0	0
7	i	90	Total 270	C 180	N 90	0	0
7	j	90	Total 270	C 180	N 90	0	0

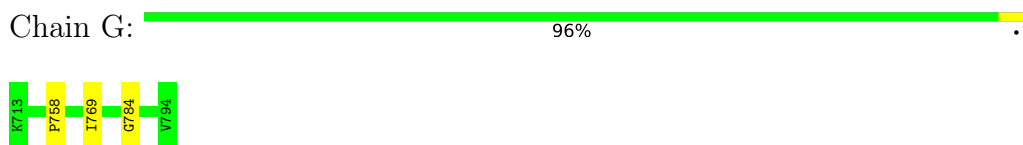
3 Residue-property plots [i](#)

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.

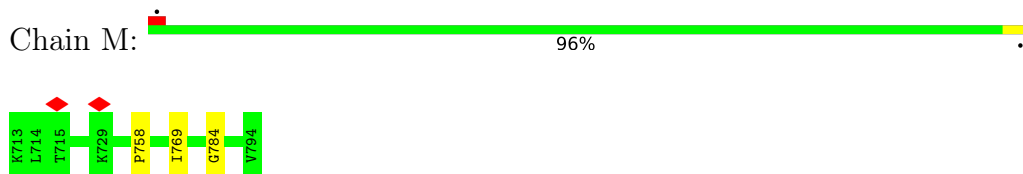
- Molecule 1: ATP-dependent Clp protease ATP-binding subunit



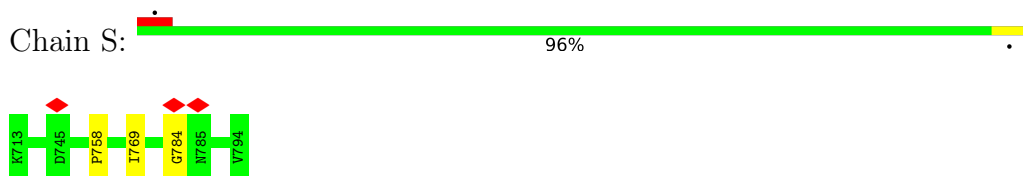
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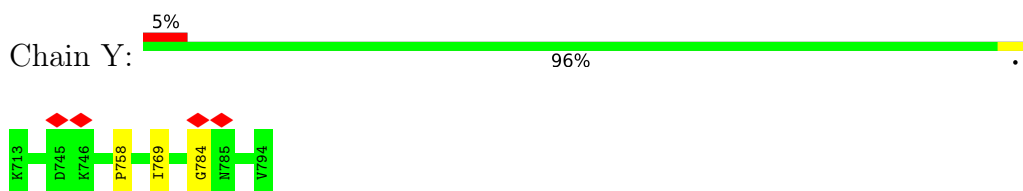
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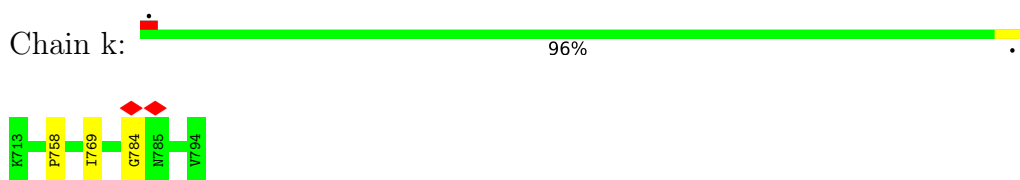
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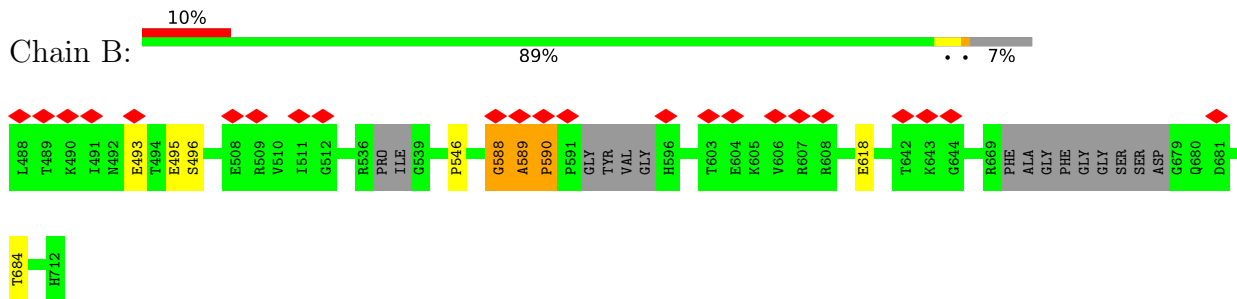
- Molecule 1: ATP-dependent Clp protease ATP-binding subunit



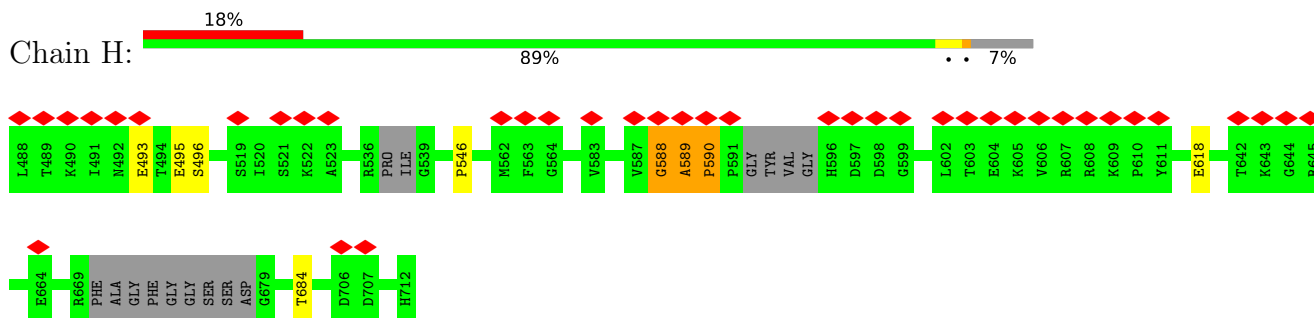
- Molecule 1: ATP-dependent Clp protease ATP-binding subunit



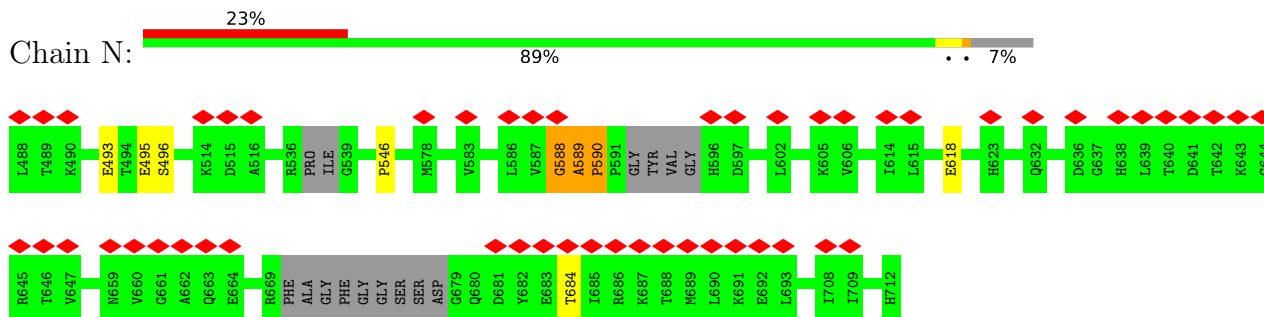
- Molecule 2: ATP-dependent Clp protease ATP-binding subunit



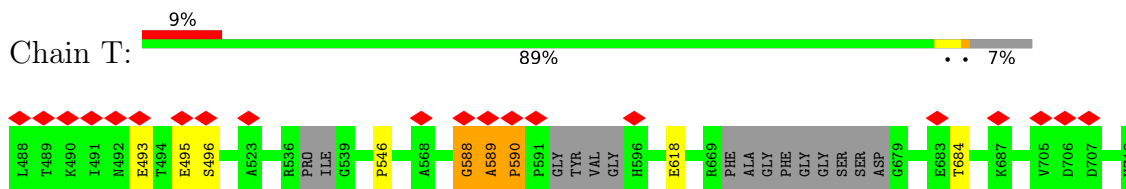
- Molecule 2: ATP-dependent Clp protease ATP-binding subunit



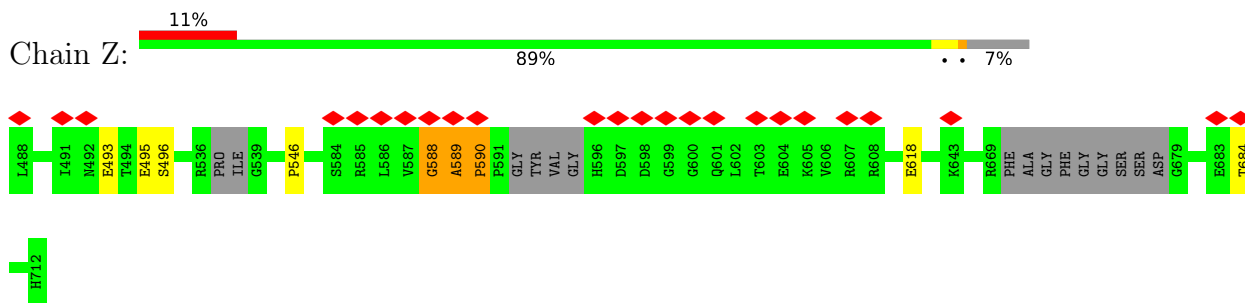
- Molecule 2: ATP-dependent Clp protease ATP-binding subunit



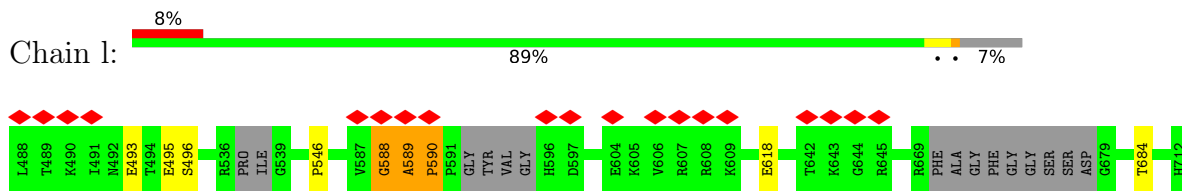
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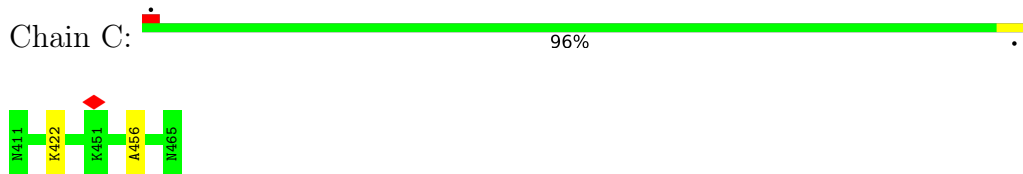
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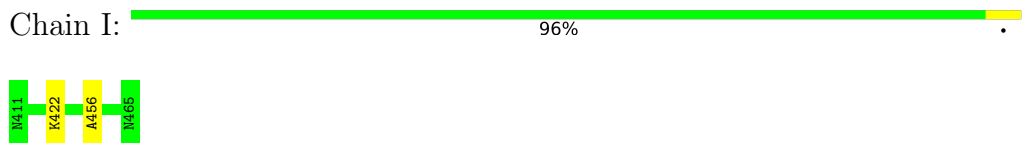
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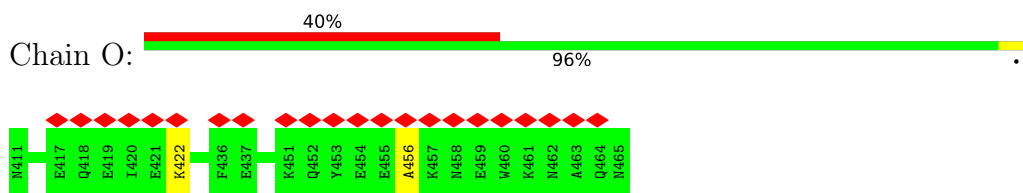
• Molecule 3: Class III stress response-related ATPase, AAA+ superfamily



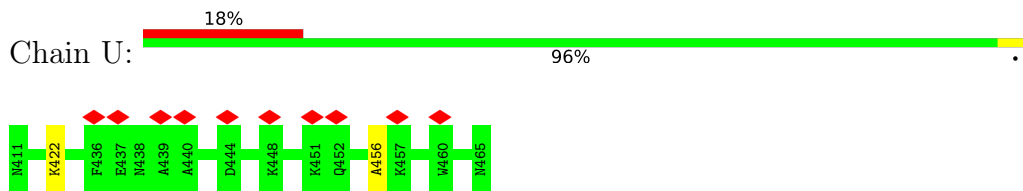
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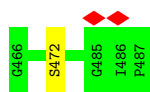


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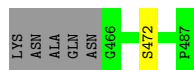
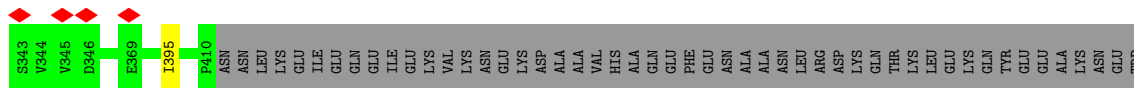


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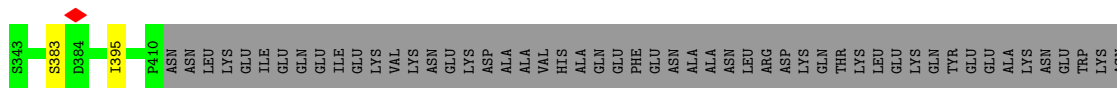




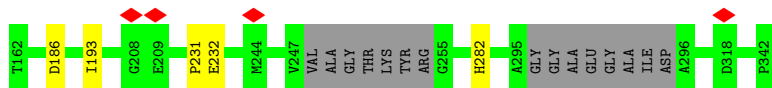
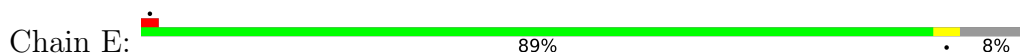
• Molecule 4: ATP-dependent Clp protease ATP-binding subunit ClpC



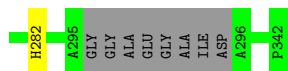
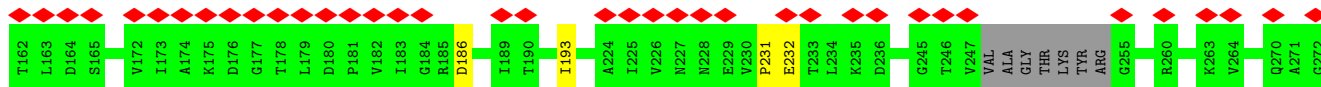
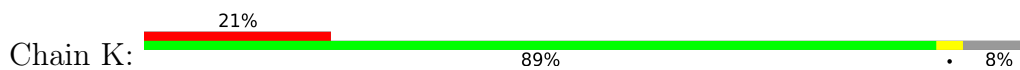
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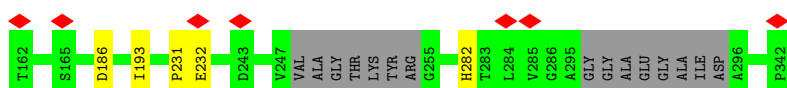
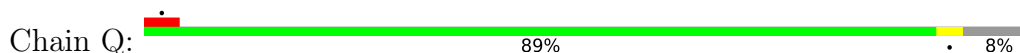
• Molecule 5: ATP-dependent Clp protease ATP-binding subunit ClpC



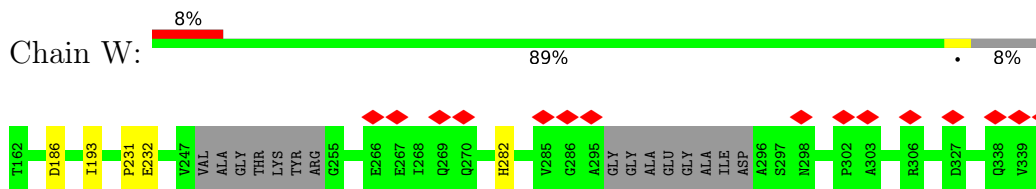
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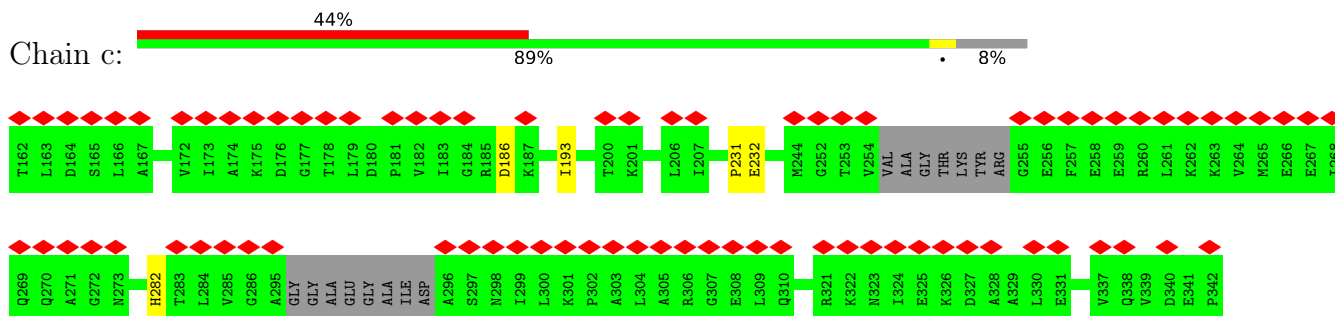
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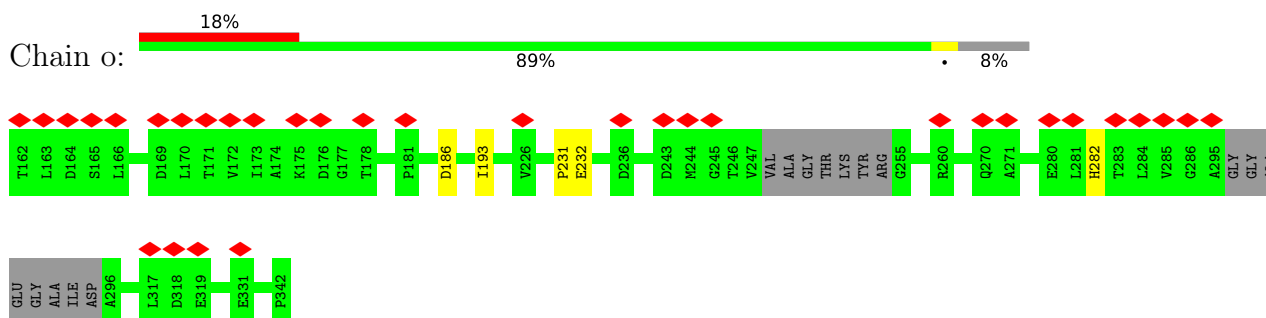
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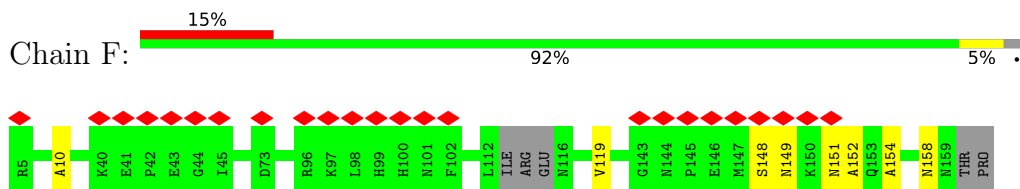
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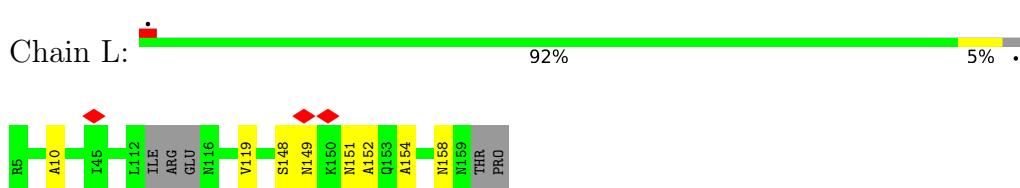
- Molecule 5: ATP-dependent Clp protease ATP-binding subunit ClpC



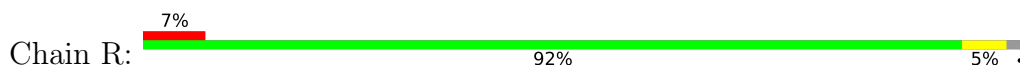
- Molecule 6: ATP-dependent Clp protease ATP-binding subunit ClpC

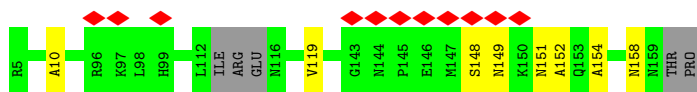


- Molecule 6: ATP-dependent Clp protease ATP-binding subunit ClpC

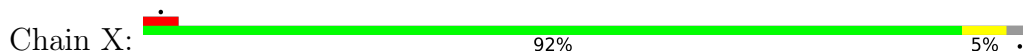


- Molecule 6: ATP-dependent Clp protease ATP-binding subunit ClpC

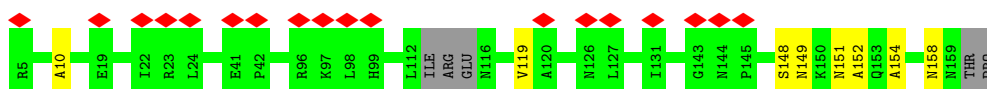
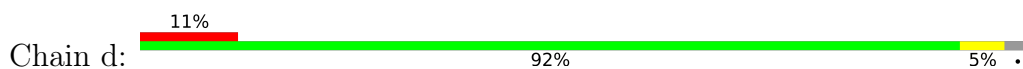




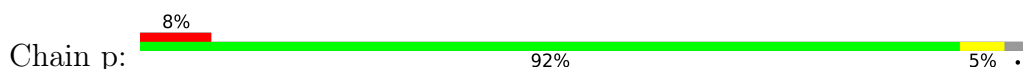
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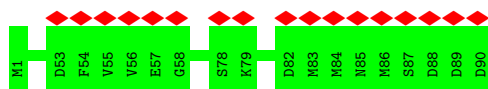
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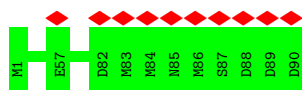
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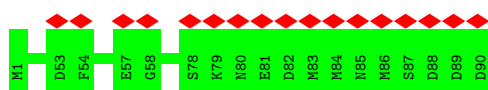
- Molecule 7: Adapter protein MecA



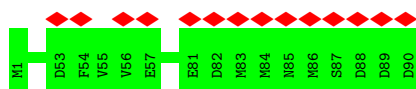
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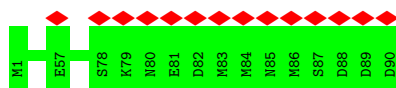
- Molecule 7: Adapter protein MecA



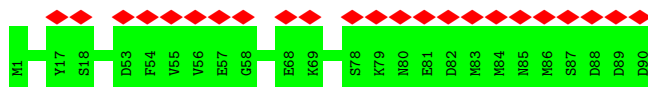
- Molecule 7: Adapter protein MecA



- Molecule 7: Adapter protein MecA



- Molecule 7: Adapter protein MecA



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	26000	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 ($e^-/\text{\AA}^2$)	1.25	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	3000	Depositor
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.069	Depositor
Minimum map value	-0.022	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.004	Depositor
Recommended contour level	0.01	Depositor
Map size (\AA)	353.6, 353.6, 353.6	wwPDB
Map dimensions	260, 260, 260	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (\AA)	1.36, 1.36, 1.36	Depositor

5 Model quality i

5.1 Standard geometry i

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.45	0/244	0.67	0/242
1	G	0.45	0/244	0.67	0/242
1	M	0.45	0/244	0.67	0/242
1	S	0.45	0/244	0.67	0/242
1	Y	0.45	0/244	0.67	0/242
1	k	0.45	0/244	0.67	0/242
2	B	0.63	1/626 (0.2%)	0.81	3/622 (0.5%)
2	H	0.63	1/626 (0.2%)	0.81	3/622 (0.5%)
2	N	0.63	1/626 (0.2%)	0.81	3/622 (0.5%)
2	T	0.63	1/626 (0.2%)	0.81	3/622 (0.5%)
2	Z	0.63	1/626 (0.2%)	0.81	3/622 (0.5%)
2	l	0.63	1/626 (0.2%)	0.81	3/622 (0.5%)
3	C	0.22	0/162	0.49	0/159
3	I	0.22	0/162	0.49	0/159
3	O	0.22	0/162	0.49	0/159
3	U	0.22	0/162	0.49	0/159
3	a	0.22	0/162	0.49	0/159
3	m	0.22	0/162	0.49	0/159
4	D	0.46	0/268	0.73	0/266
4	J	0.46	0/268	0.73	0/266
4	P	0.46	0/268	0.73	0/266
4	V	0.46	0/268	0.73	0/266
4	b	0.46	0/268	0.73	0/266
4	n	0.46	0/268	0.73	0/266
5	E	0.46	0/494	0.74	0/490
5	K	0.46	0/494	0.74	0/490
5	Q	0.46	0/494	0.74	0/490
5	W	0.46	0/494	0.74	0/490
5	c	0.46	0/494	0.74	0/490
5	o	0.46	0/494	0.74	0/490
6	F	0.29	0/452	0.73	0/448
6	L	0.28	0/452	0.73	0/448
6	R	0.28	0/452	0.73	0/448
6	X	0.29	0/452	0.73	0/448

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
6	d	0.28	0/452	0.73	0/448
6	p	0.28	0/452	0.73	0/448
7	e	0.38	0/269	0.87	0/268
7	f	0.38	0/269	0.88	0/268
7	g	0.38	0/269	0.87	0/268
7	h	0.39	0/269	0.87	0/268
7	i	0.39	0/269	0.87	0/268
7	j	0.39	0/269	0.87	0/268
All	All	0.47	6/15090 (0.0%)	0.75	18/14970 (0.1%)

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
2	B	0	1
2	H	0	1
2	N	0	1
2	T	0	1
2	Z	0	1
2	l	0	1
4	n	0	1
5	E	0	2
5	K	0	2
5	Q	0	2
5	W	0	2
5	c	0	2
5	o	0	2
All	All	0	19

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	N	590	PRO	CA-C	-10.01	1.32	1.52
2	l	590	PRO	CA-C	-10.00	1.32	1.52
2	T	590	PRO	CA-C	-9.98	1.32	1.52
2	Z	590	PRO	CA-C	-9.97	1.32	1.52
2	H	590	PRO	CA-C	-9.97	1.32	1.52
2	B	590	PRO	CA-C	-9.95	1.32	1.52

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	l	589	ALA	N-CA-C	-8.39	88.34	111.00
2	H	589	ALA	N-CA-C	-8.39	88.36	111.00
2	Z	589	ALA	N-CA-C	-8.38	88.38	111.00
2	N	589	ALA	N-CA-C	-8.38	88.39	111.00
2	T	589	ALA	N-CA-C	-8.37	88.39	111.00
2	B	589	ALA	N-CA-C	-8.36	88.43	111.00
2	T	590	PRO	CA-C-N	6.35	134.88	117.10
2	l	590	PRO	CA-C-N	6.35	134.87	117.10
2	Z	590	PRO	CA-C-N	6.34	134.86	117.10
2	B	590	PRO	CA-C-N	6.33	134.84	117.10
2	N	590	PRO	CA-C-N	6.33	134.81	117.10
2	H	590	PRO	CA-C-N	6.32	134.79	117.10
2	H	588	GLY	C-N-CA	-6.17	106.28	121.70
2	Z	588	GLY	C-N-CA	-6.15	106.32	121.70
2	l	588	GLY	C-N-CA	-6.15	106.33	121.70
2	B	588	GLY	C-N-CA	-6.14	106.35	121.70
2	N	588	GLY	C-N-CA	-6.14	106.36	121.70
2	T	588	GLY	C-N-CA	-6.13	106.38	121.70

There are no chirality outliers.

All (19) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	618	GLU	Peptide
5	E	231	PRO	Peptide
5	E	282	HIS	Peptide
2	H	618	GLU	Peptide
5	K	231	PRO	Peptide
5	K	282	HIS	Peptide
2	N	618	GLU	Peptide
5	Q	231	PRO	Peptide
5	Q	282	HIS	Peptide
2	T	618	GLU	Peptide
5	W	231	PRO	Peptide
5	W	282	HIS	Peptide
2	Z	618	GLU	Peptide
5	c	231	PRO	Peptide
5	c	282	HIS	Peptide
2	l	618	GLU	Peptide
4	n	383	SER	Peptide
5	o	231	PRO	Peptide
5	o	282	HIS	Peptide

5.2 Too-close contacts [i](#)

Due to software issues we are unable to calculate clashes - this section is therefore empty.

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	78/82 (95%)	55 (70%)	20 (26%)	3 (4%)	3	24
1	G	78/82 (95%)	55 (70%)	20 (26%)	3 (4%)	3	24
1	M	78/82 (95%)	55 (70%)	20 (26%)	3 (4%)	3	24
1	S	78/82 (95%)	55 (70%)	20 (26%)	3 (4%)	3	24
1	Y	78/82 (95%)	55 (70%)	20 (26%)	3 (4%)	3	24
1	k	78/82 (95%)	55 (70%)	20 (26%)	3 (4%)	3	24
2	B	202/225 (90%)	162 (80%)	32 (16%)	8 (4%)	3	23
2	H	202/225 (90%)	162 (80%)	32 (16%)	8 (4%)	3	23
2	N	202/225 (90%)	162 (80%)	32 (16%)	8 (4%)	3	23
2	T	202/225 (90%)	162 (80%)	32 (16%)	8 (4%)	3	23
2	Z	202/225 (90%)	162 (80%)	32 (16%)	8 (4%)	3	23
2	l	202/225 (90%)	162 (80%)	32 (16%)	8 (4%)	3	23
3	C	50/55 (91%)	45 (90%)	3 (6%)	2 (4%)	3	23
3	I	50/55 (91%)	45 (90%)	3 (6%)	2 (4%)	3	23
3	O	50/55 (91%)	45 (90%)	3 (6%)	2 (4%)	3	23
3	U	50/55 (91%)	45 (90%)	3 (6%)	2 (4%)	3	23
3	a	50/55 (91%)	45 (90%)	3 (6%)	2 (4%)	3	23
3	m	50/55 (91%)	45 (90%)	3 (6%)	2 (4%)	3	23
4	D	86/145 (59%)	68 (79%)	16 (19%)	2 (2%)	6	34
4	J	86/145 (59%)	68 (79%)	16 (19%)	2 (2%)	6	34
4	P	86/145 (59%)	68 (79%)	15 (17%)	3 (4%)	3	25

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
4	V	86/145 (59%)	68 (79%)	16 (19%)	2 (2%)	6	34
4	b	86/145 (59%)	68 (79%)	16 (19%)	2 (2%)	6	34
4	n	86/145 (59%)	68 (79%)	15 (17%)	3 (4%)	3	25
5	E	158/181 (87%)	131 (83%)	24 (15%)	3 (2%)	8	38
5	K	158/181 (87%)	131 (83%)	24 (15%)	3 (2%)	8	38
5	Q	158/181 (87%)	131 (83%)	24 (15%)	3 (2%)	8	38
5	W	158/181 (87%)	131 (83%)	24 (15%)	3 (2%)	8	38
5	c	158/181 (87%)	131 (83%)	24 (15%)	3 (2%)	8	38
5	o	158/181 (87%)	131 (83%)	24 (15%)	3 (2%)	8	38
6	F	144/157 (92%)	116 (81%)	20 (14%)	8 (6%)	2	19
6	L	144/157 (92%)	116 (81%)	20 (14%)	8 (6%)	2	19
6	R	144/157 (92%)	116 (81%)	20 (14%)	8 (6%)	2	19
6	X	144/157 (92%)	116 (81%)	20 (14%)	8 (6%)	2	19
6	d	144/157 (92%)	116 (81%)	20 (14%)	8 (6%)	2	19
6	p	144/157 (92%)	116 (81%)	20 (14%)	8 (6%)	2	19
7	e	88/90 (98%)	80 (91%)	8 (9%)	0	100	100
7	f	88/90 (98%)	80 (91%)	8 (9%)	0	100	100
7	g	88/90 (98%)	80 (91%)	8 (9%)	0	100	100
7	h	88/90 (98%)	80 (91%)	8 (9%)	0	100	100
7	i	88/90 (98%)	80 (91%)	8 (9%)	0	100	100
7	j	88/90 (98%)	80 (91%)	8 (9%)	0	100	100
All	All	4836/5610 (86%)	3942 (82%)	736 (15%)	158 (3%)	6	26

All (158) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	493	GLU
2	B	495	GLU
2	B	589	ALA
2	B	590	PRO
6	F	10	ALA
6	F	148	SER
6	F	149	ASN
6	F	152	ALA
6	F	154	ALA

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Mol	Chain	Res	Type
2	H	493	GLU
2	H	495	GLU
2	H	589	ALA
2	H	590	PRO
6	L	10	ALA
6	L	148	SER
6	L	149	ASN
6	L	152	ALA
6	L	154	ALA
2	N	493	GLU
2	N	495	GLU
2	N	589	ALA
2	N	590	PRO
6	R	10	ALA
6	R	148	SER
6	R	149	ASN
6	R	152	ALA
6	R	154	ALA
2	T	493	GLU
2	T	495	GLU
2	T	589	ALA
2	T	590	PRO
6	X	10	ALA
6	X	148	SER
6	X	149	ASN
6	X	152	ALA
6	X	154	ALA
2	Z	493	GLU
2	Z	495	GLU
2	Z	589	ALA
2	Z	590	PRO
6	d	10	ALA
6	d	148	SER
6	d	149	ASN
6	d	152	ALA
6	d	154	ALA
2	l	493	GLU
2	l	495	GLU
2	l	589	ALA
2	l	590	PRO
6	p	10	ALA
6	p	148	SER

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Mol	Chain	Res	Type
6	p	149	ASN
6	p	152	ALA
6	p	154	ALA
6	F	119	VAL
6	F	151	ASN
6	L	119	VAL
6	L	151	ASN
6	R	119	VAL
6	R	151	ASN
6	X	119	VAL
6	X	151	ASN
6	d	119	VAL
6	d	151	ASN
6	p	119	VAL
6	p	151	ASN
2	B	496	SER
6	F	158	ASN
2	H	496	SER
6	L	158	ASN
2	N	496	SER
6	R	158	ASN
2	T	496	SER
6	X	158	ASN
2	Z	496	SER
6	d	158	ASN
2	l	496	SER
6	p	158	ASN
1	A	758	PRO
2	B	546	PRO
2	B	684	THR
3	C	422	LYS
3	C	456	ALA
1	G	758	PRO
2	H	546	PRO
2	H	684	THR
3	I	422	LYS
3	I	456	ALA
5	K	232	GLU
1	M	758	PRO
2	N	546	PRO
2	N	684	THR
3	O	422	LYS

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Mol	Chain	Res	Type
3	O	456	ALA
5	Q	232	GLU
1	S	758	PRO
2	T	546	PRO
2	T	684	THR
3	U	422	LYS
3	U	456	ALA
5	W	232	GLU
1	Y	758	PRO
2	Z	546	PRO
2	Z	684	THR
3	a	422	LYS
3	a	456	ALA
5	c	232	GLU
1	k	758	PRO
2	l	546	PRO
2	l	684	THR
3	m	422	LYS
3	m	456	ALA
5	o	232	GLU
4	D	472	SER
5	E	186	ASP
5	E	232	GLU
4	J	472	SER
5	K	186	ASP
4	P	472	SER
4	P	477	ALA
5	Q	186	ASP
4	V	472	SER
5	W	186	ASP
4	b	472	SER
5	c	186	ASP
4	n	472	SER
4	n	477	ALA
5	o	186	ASP
2	B	588	GLY
2	H	588	GLY
2	N	588	GLY
2	T	588	GLY
2	Z	588	GLY
2	l	588	GLY
1	A	769	ILE

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Mol	Chain	Res	Type
4	D	395	ILE
1	G	769	ILE
4	J	395	ILE
1	M	769	ILE
4	P	395	ILE
1	S	769	ILE
4	V	395	ILE
1	Y	769	ILE
4	b	395	ILE
1	k	769	ILE
4	n	395	ILE
1	A	784	GLY
5	E	193	ILE
1	G	784	GLY
5	K	193	ILE
1	M	784	GLY
5	Q	193	ILE
1	S	784	GLY
5	W	193	ILE
1	Y	784	GLY
5	c	193	ILE
1	k	784	GLY
5	o	193	ILE

5.3.2 Protein sidechains [i](#)

There are no protein residues with a non-rotameric sidechain to report in this entry.

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
3	C	2
3	I	2
3	O	2
3	U	2
3	a	2
3	m	2
6	X	2
6	F	2
6	L	2
6	R	2
6	d	2
6	p	2
1	A	1
1	G	1
1	M	1
1	S	1
1	Y	1
1	k	1
5	E	1
5	K	1
5	Q	1
5	W	1
5	c	1
5	o	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	C	464:GLN	C	465:ASN	N	12.34
1	I	464:GLN	C	465:ASN	N	12.34

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Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	O	464:GLN	C	465:ASN	N	12.34
1	U	464:GLN	C	465:ASN	N	12.34
1	a	464:GLN	C	465:ASN	N	12.34
1	m	464:GLN	C	465:ASN	N	12.34
1	X	69:GLY	C	70:HIS	N	7.19
1	F	69:GLY	C	70:HIS	N	7.18
1	L	69:GLY	C	70:HIS	N	7.18
1	R	69:GLY	C	70:HIS	N	7.18
1	d	69:GLY	C	70:HIS	N	7.18
1	p	69:GLY	C	70:HIS	N	7.18
1	C	437:GLU	C	438:ASN	N	6.46
1	I	437:GLU	C	438:ASN	N	6.46
1	O	437:GLU	C	438:ASN	N	6.46
1	U	437:GLU	C	438:ASN	N	6.46
1	a	437:GLU	C	438:ASN	N	6.46
1	m	437:GLU	C	438:ASN	N	6.46
1	F	45:ILE	C	46:ALA	N	6.00
1	L	45:ILE	C	46:ALA	N	6.00
1	R	45:ILE	C	46:ALA	N	6.00
1	X	45:ILE	C	46:ALA	N	6.00
1	d	45:ILE	C	46:ALA	N	6.00
1	p	45:ILE	C	46:ALA	N	6.00
1	A	736:GLU	C	737:GLN	N	4.81
1	G	736:GLU	C	737:GLN	N	4.81
1	M	736:GLU	C	737:GLN	N	4.81
1	S	736:GLU	C	737:GLN	N	4.81
1	Y	736:GLU	C	737:GLN	N	4.81
1	k	736:GLU	C	737:GLN	N	4.81
1	E	272:GLY	C	273:ASN	N	2.90
1	K	272:GLY	C	273:ASN	N	2.90
1	Q	272:GLY	C	273:ASN	N	2.90
1	W	272:GLY	C	273:ASN	N	2.90
1	c	272:GLY	C	273:ASN	N	2.90
1	o	272:GLY	C	273:ASN	N	2.90

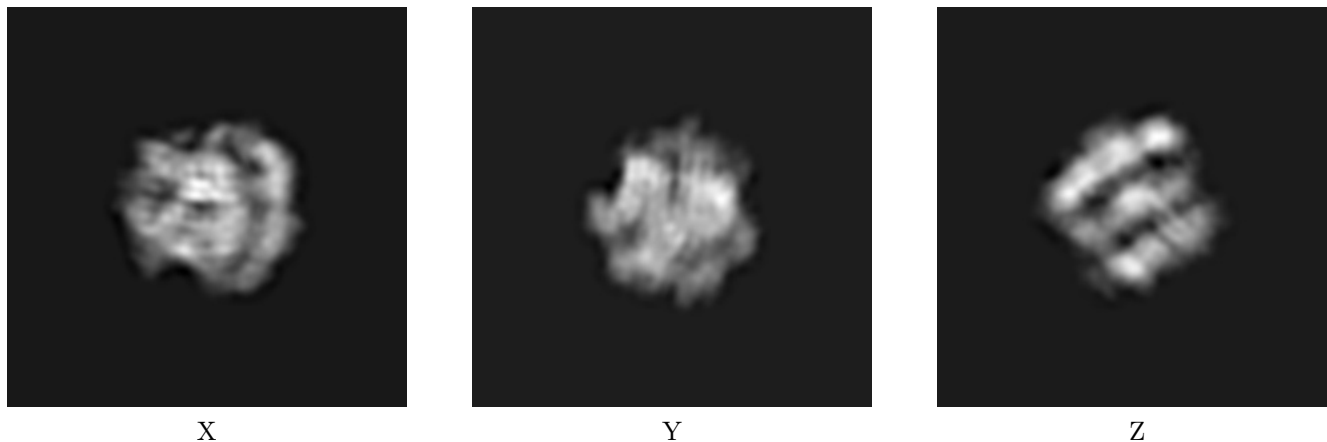
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-3897. These allow visual inspection of the internal detail of the map and identification of artifacts.

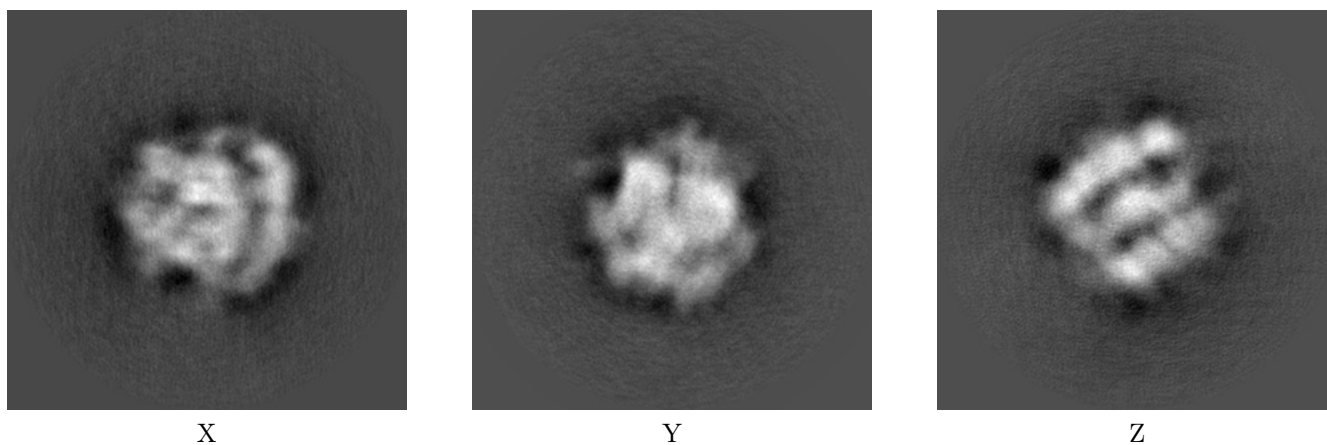
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

6.1.1 Primary map



6.1.2 Raw map



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

6.2 Central slices [i](#)

6.2.1 Primary map



X Index: 130

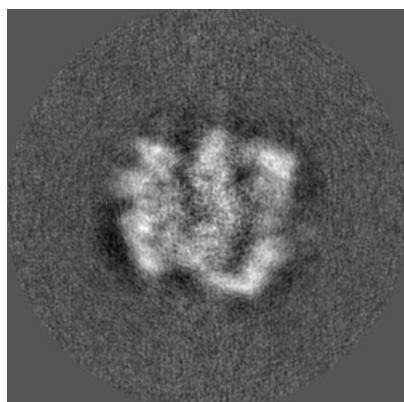


Y Index: 130

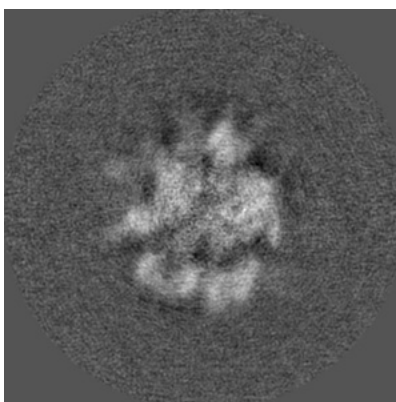


Z Index: 130

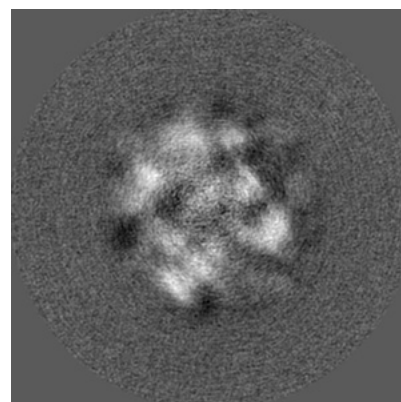
6.2.2 Raw map



X Index: 130



Y Index: 130

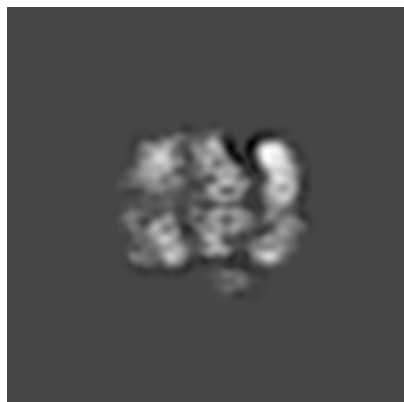


Z Index: 130

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: 140

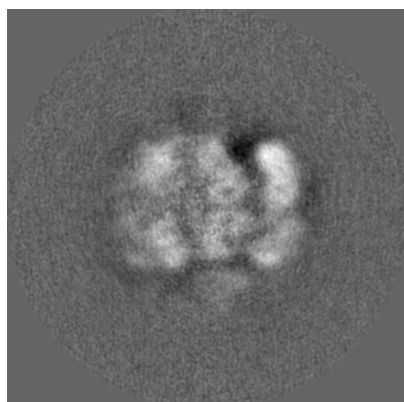


Y Index: 139

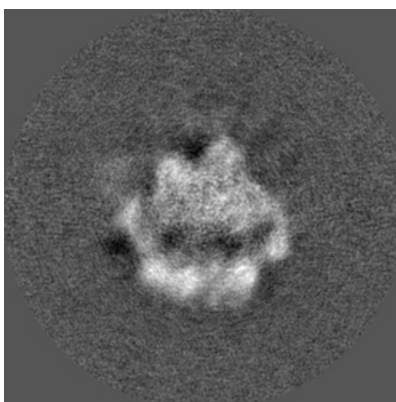


Z Index: 118

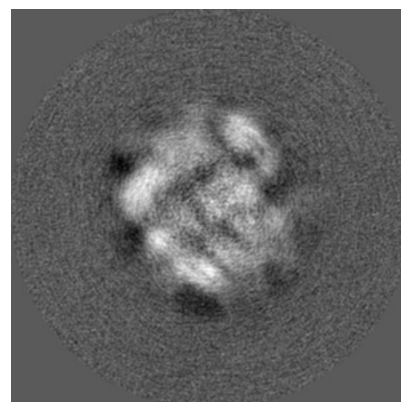
6.3.2 Raw map



X Index: 140



Y Index: 139

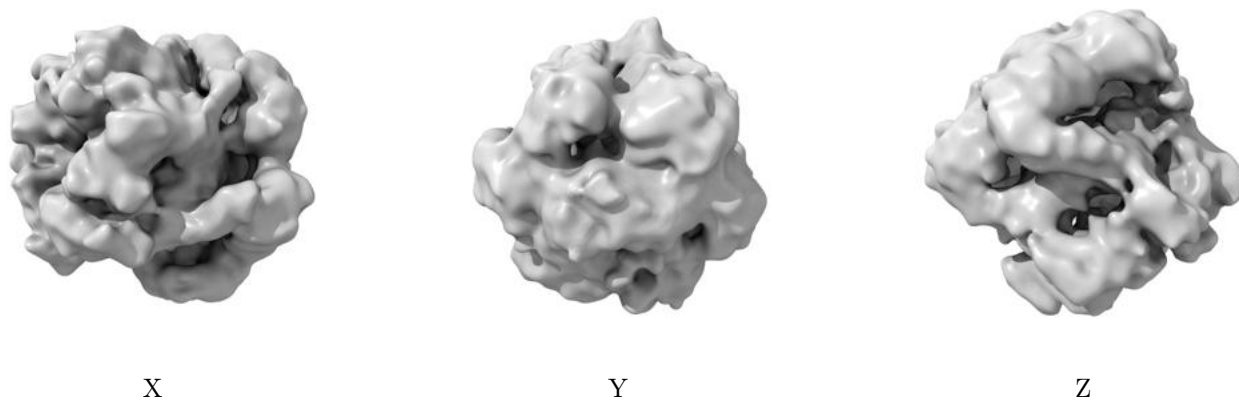


Z Index: 117

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

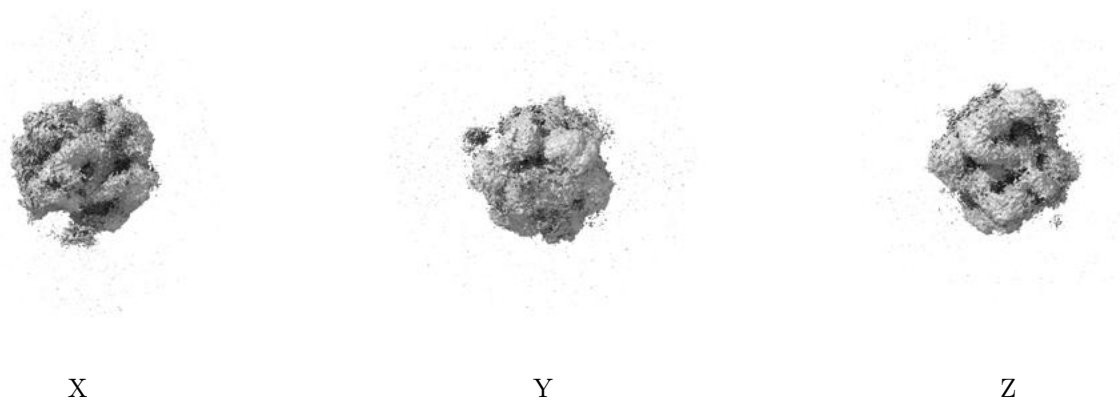
6.4 Orthogonal surface views [i](#)

6.4.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.01. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.4.2 Raw map



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.

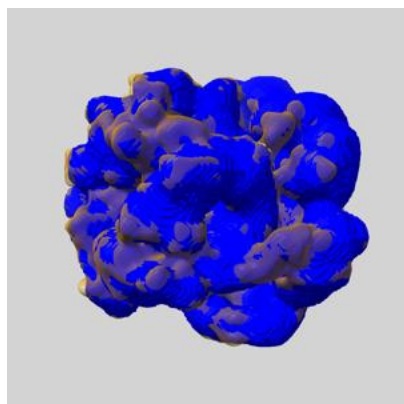
6.5 Mask visualisation [i](#)

This section shows the 3D surface view of the primary map at 50% transparency overlaid with the specified mask at 0% transparency

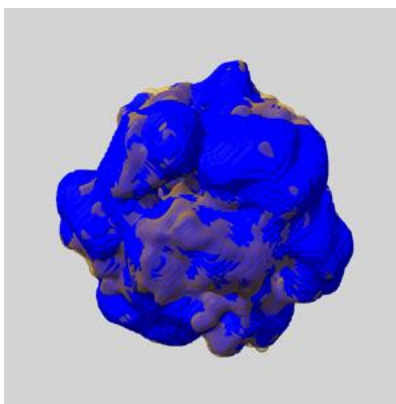
A mask typically either:

- Encompasses the whole structure
- Separates out a domain, a functional unit, a monomer or an area of interest from a larger structure

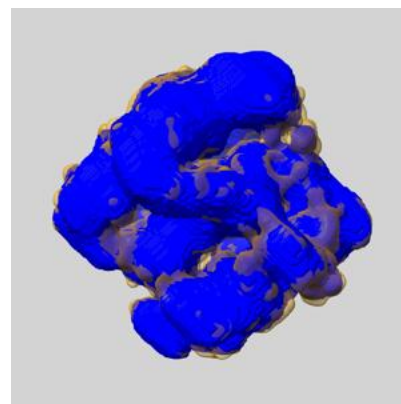
6.5.1 emd_3897_msk_1.map [i](#)



X



Y

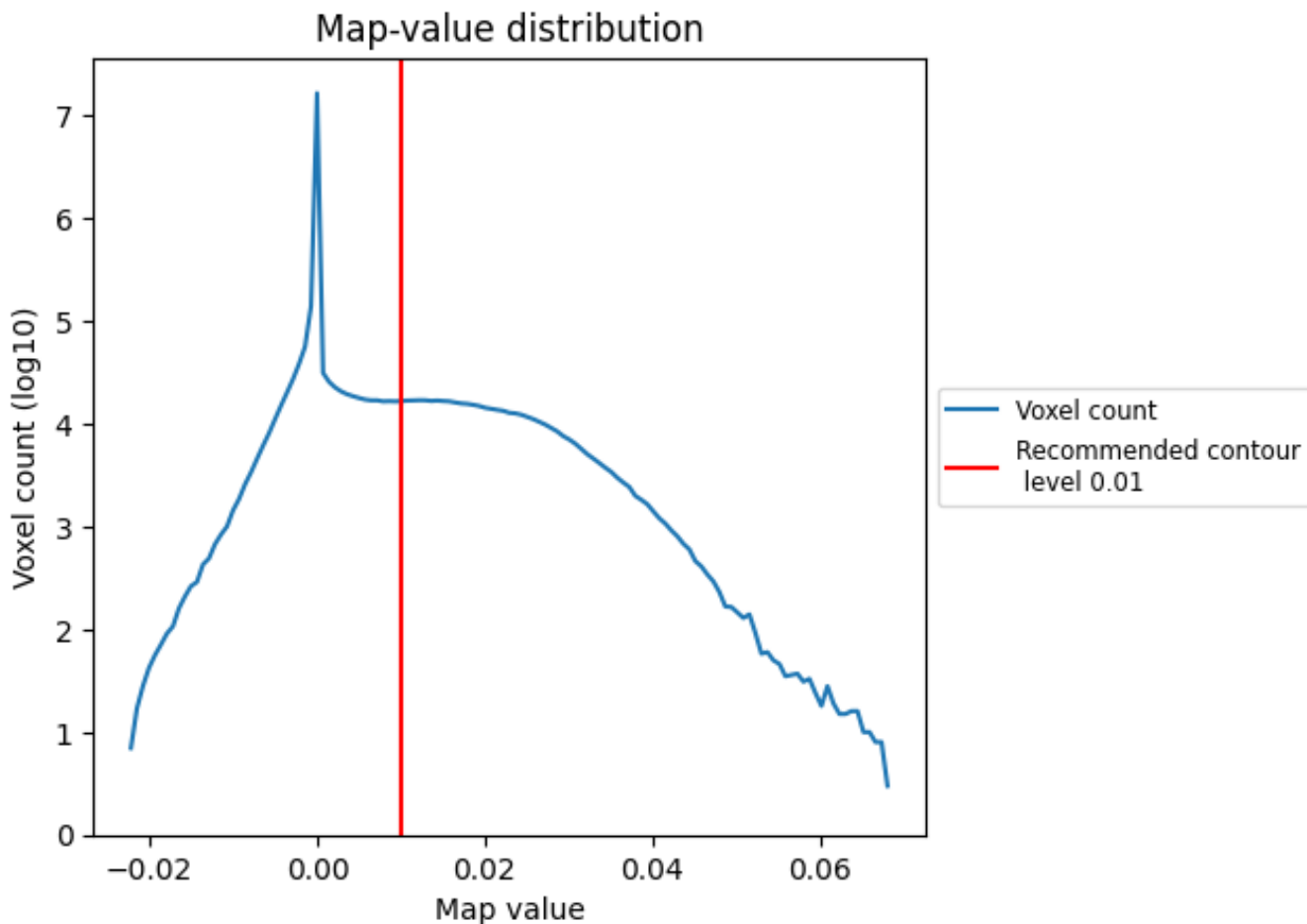


Z

7 Map analysis [i](#)

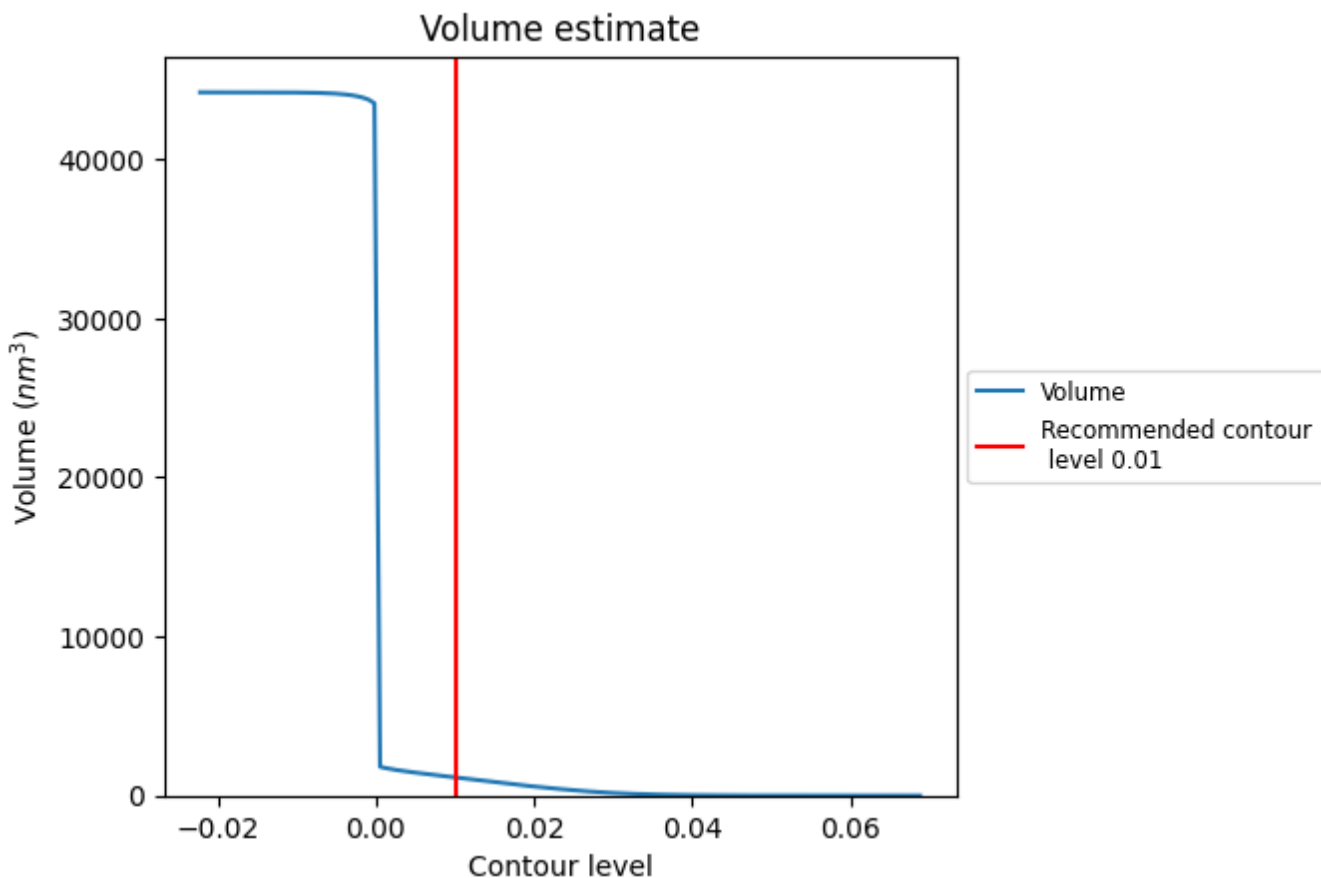
This section contains the results of statistical analysis of the map.

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.

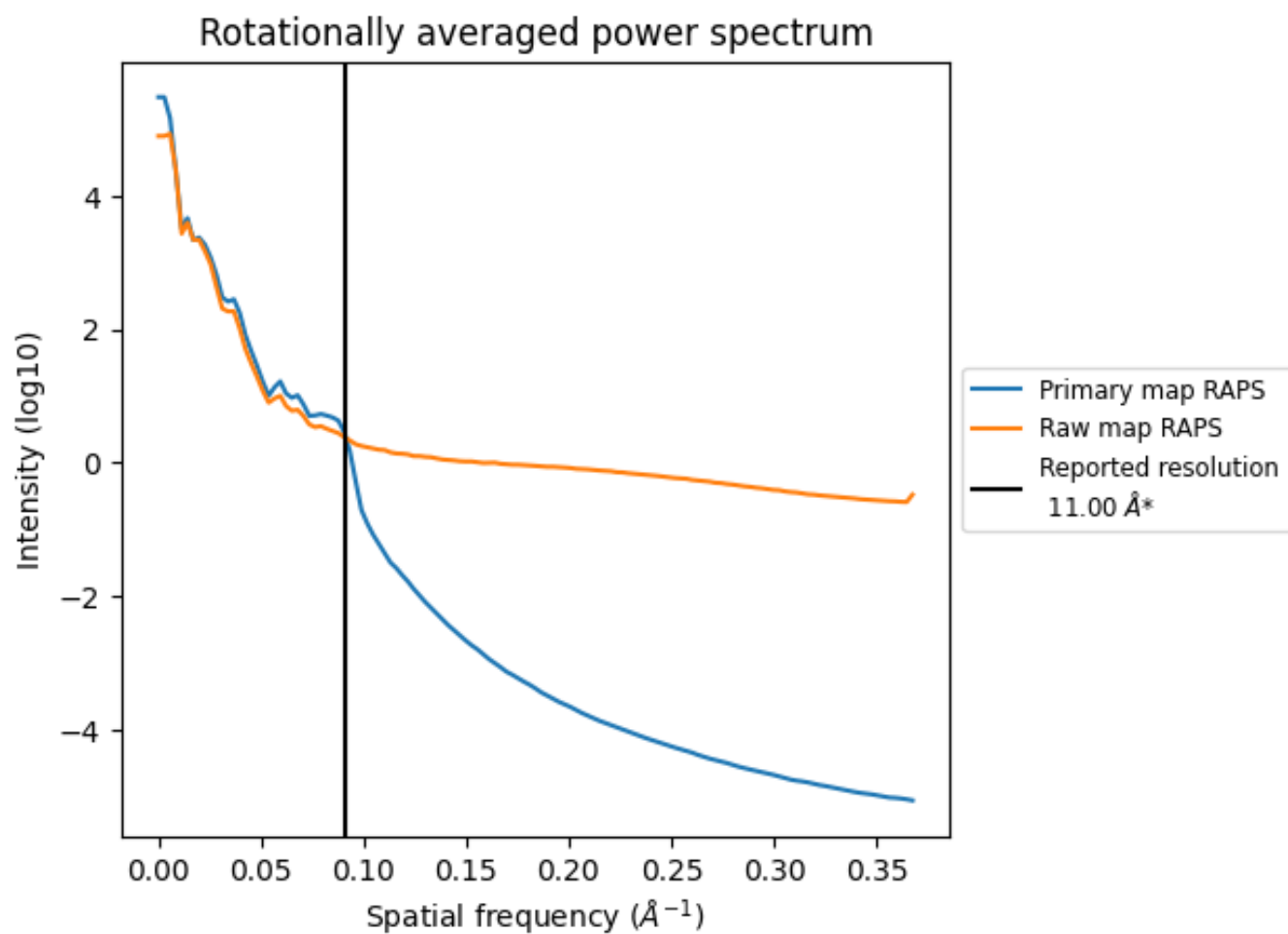
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 1143 nm^3 ; this corresponds to an approximate mass of 1033 kDa.

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

7.3 Rotationally averaged power spectrum i

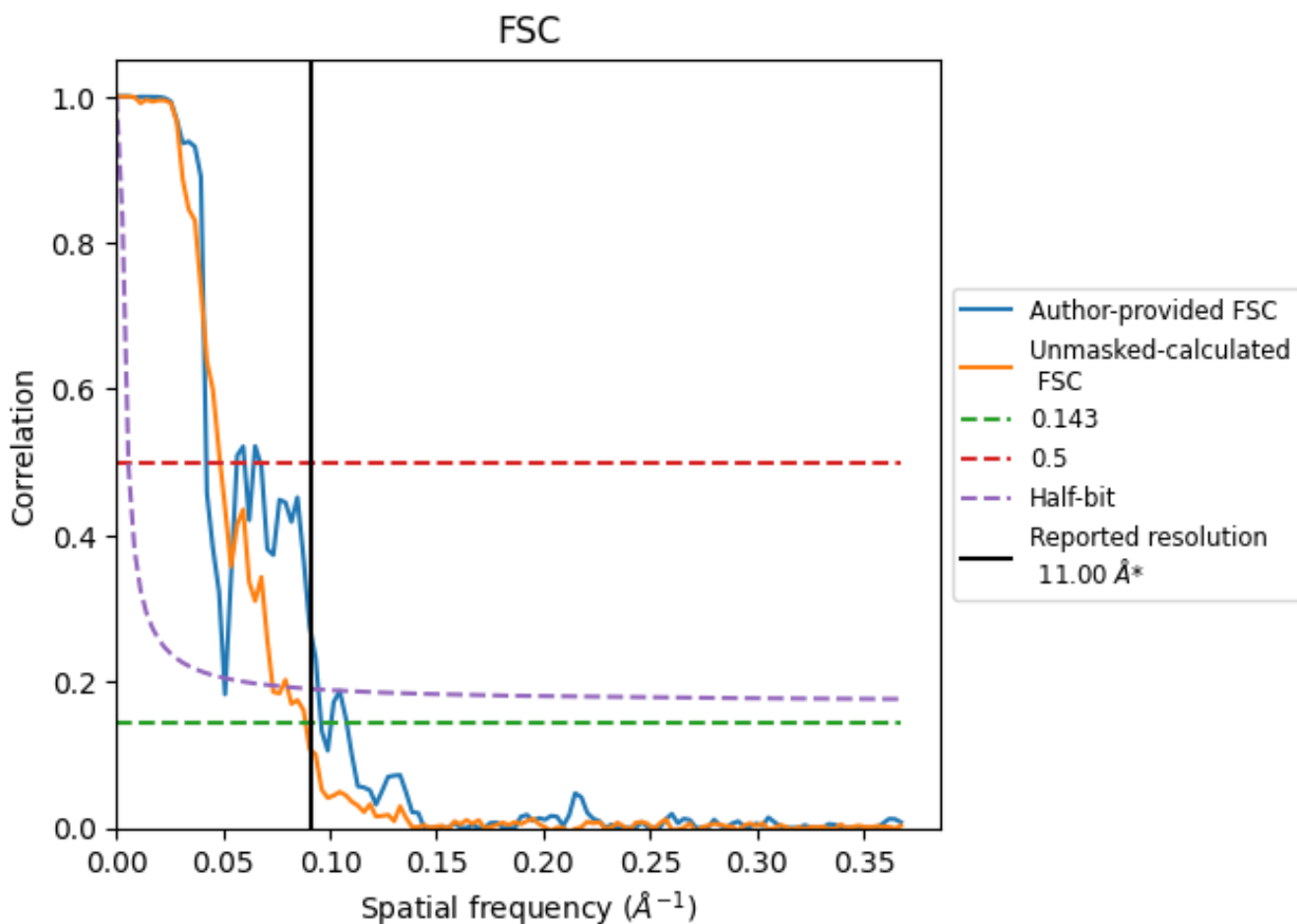


*Reported resolution corresponds to spatial frequency of 0.091 Å⁻¹

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 0.091 Å⁻¹

8.2 Resolution estimates [i](#)

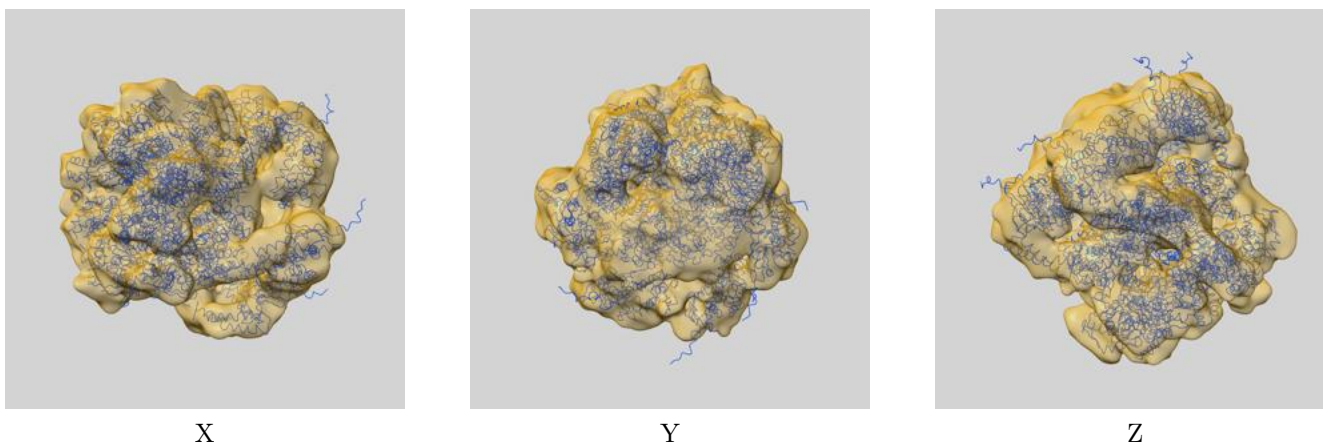
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	11.00	-	-
Author-provided FSC curve	10.44	23.75	19.80
Unmasked-calculated*	11.29	20.58	13.66

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps.

9 Map-model fit [i](#)

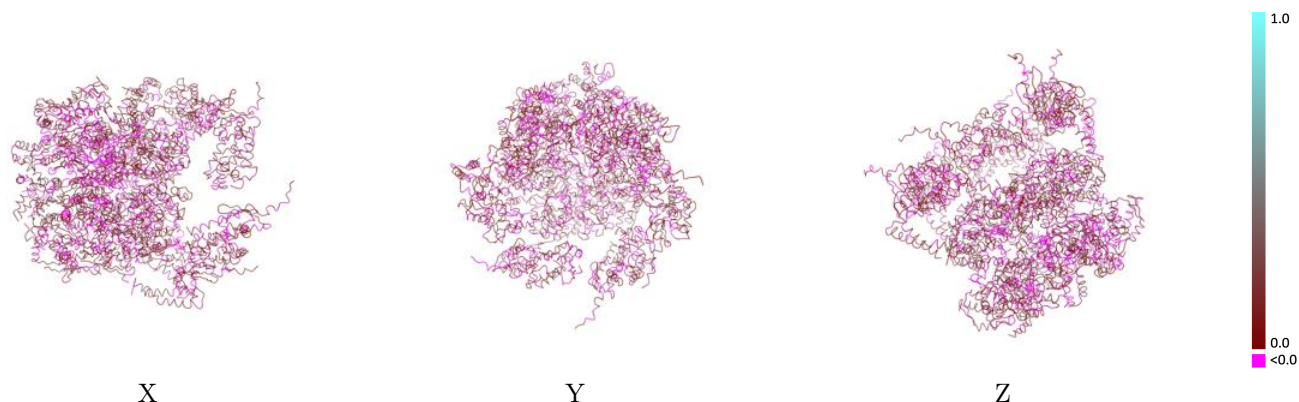
This section contains information regarding the fit between EMDB map EMD-3897 and PDB model 6EMW. Per-residue inclusion information can be found in section 3 on page 8.

9.1 Map-model overlay [i](#)



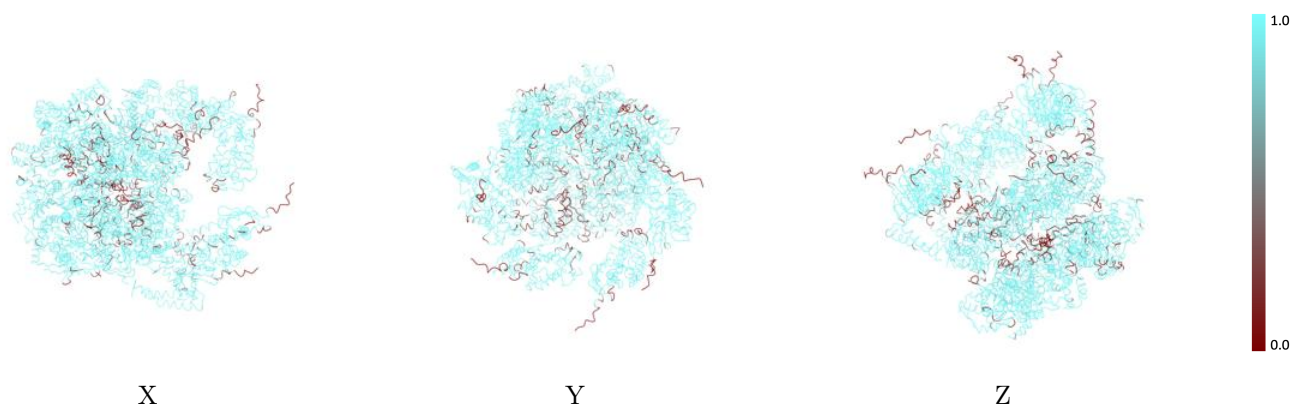
The images above show the 3D surface view of the map at the recommended contour level 0.01 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



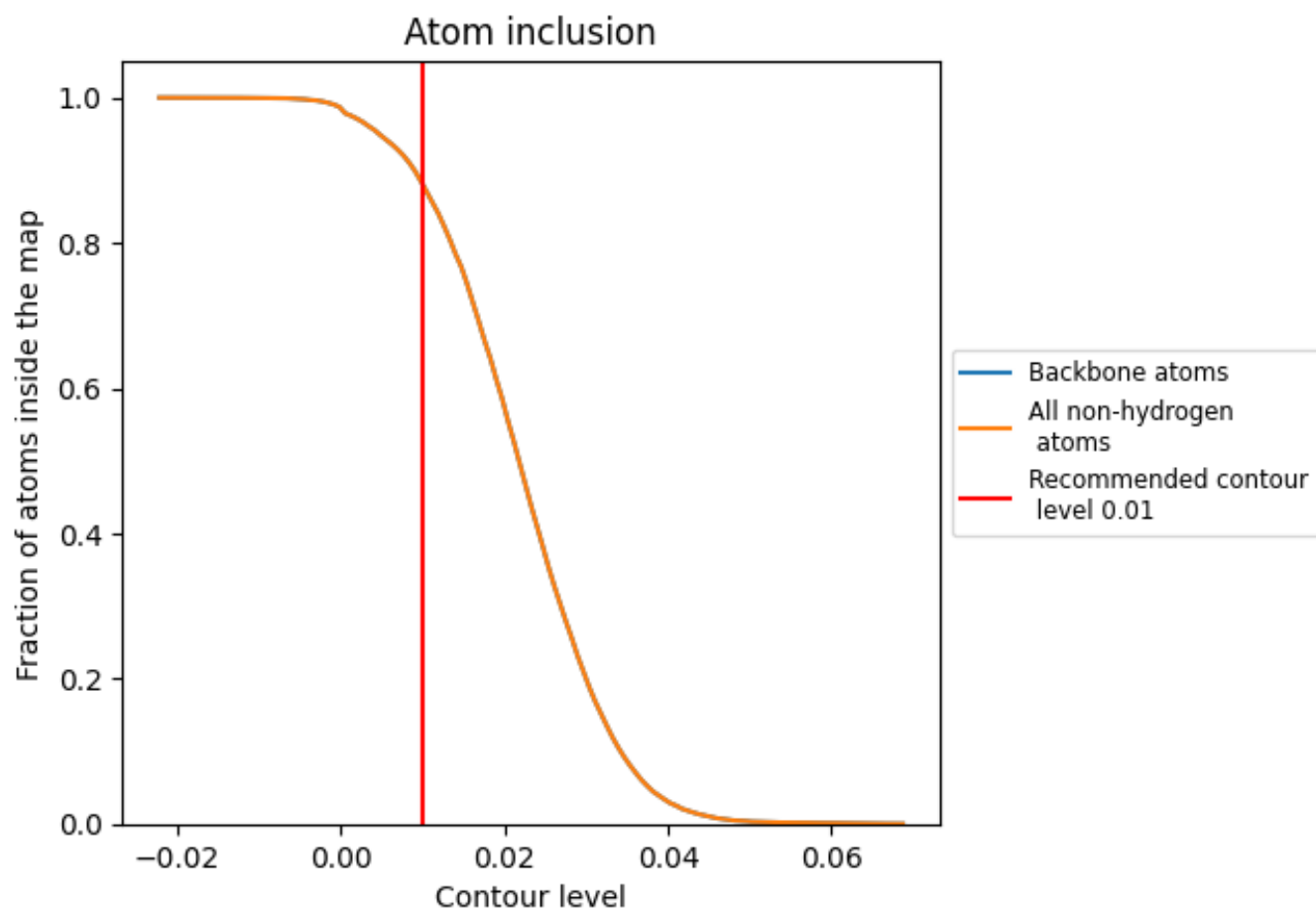
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.01).







































































9.4 Atom inclusion [i](#)



At the recommended contour level, 88% of all backbone atoms, 88% 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 (0.01) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8807	 0.0740
A	 0.9634	 0.1260
B	 0.8873	 0.0720
C	 0.9758	 0.0870
D	 0.9037	 0.1020
E	 0.9659	 0.1000
F	 0.8596	 0.0630
G	 1.0000	 0.0640
H	 0.7984	 0.0240
I	 0.9939	 0.1340
J	 0.9704	 0.1090
K	 0.7751	 0.0580
L	 0.9803	 0.1110
M	 0.9593	 0.1400
N	 0.7317	 -0.0470
O	 0.6182	 0.0190
P	 0.9778	 0.1320
Q	 0.9618	 0.0800
R	 0.9342	 0.0850
S	 0.9756	 0.0570
T	 0.9032	 0.0730
U	 0.8242	 -0.0050
V	 0.9778	 0.1450
W	 0.9056	 0.0870
X	 0.9386	 0.0920
Y	 0.9553	 0.0740
Z	 0.8921	 0.0940
a	 0.9939	 0.1290
b	 0.9630	 0.0950
c	 0.5221	 0.0560
d	 0.8991	 0.0520
e	 0.8074	 0.0470
f	 0.8852	 0.0930
g	 0.8185	 0.0800
h	 0.8333	 0.0950



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Chain	Atom inclusion	Q-score
i	 0.8444	 0.0460
j	 0.7407	 0.0090
k	 0.9715	 0.0870
l	 0.9159	 0.1100
m	 0.8727	 0.0630
n	 0.9852	 0.1120
o	 0.7892	 0.0590
p	 0.9123	 0.0550