wwPDB EM Validation Summary Report

Oct 13, 2024 – 03:31 am BST

PDB ID	:	6HBL
EMDB ID	:	EMD-0188
Title	:	Echovirus 18 Open particle without three pentamers
Authors	:	Buchta, D.; Fuzik, T.; Hrebik, D.; Levdansky, Y.; Moravcova, J.; Plevka, P.
Deposited on	:	2018-08-10
Resolution	:	3.70 Å(reported)
Based on initial model	:	6HBH

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:

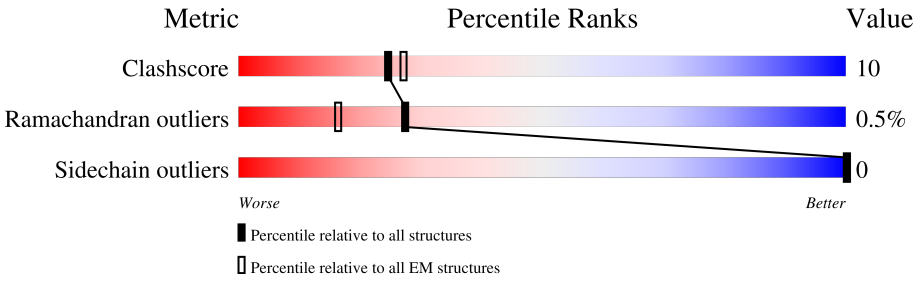
EMDB validation analysis	:	0.0.1.dev113
MolProbity	:	4.02b-467
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
MapQ	:	1.9.13
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.39

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 3.70 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	210492	15764
Ramachandran outliers	207382	16835
Sidechain outliers	206894	16415

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=5% The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion < 40%). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	287	7% 60% 18% 22%
1	D	287	11% 60% 18% 22%
1	G	287	13% 60% 18% 22%
1	J	287	25% 61% 17% 22%
1	M	287	6% 61% 17% 22%
1	P	287	14% 61% 17% 22%
1	S	287	5% 61% 17% 22%
1	V	287	41% 61% 17% 22%

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Mol	Chain	Length	Quality of chain
1	Y	287	65% 13% 22%
1	b	287	78% 22%
1	e	287	78% 22%
1	h	287	78% 22%
1	k	287	24% 78% 22%
1	n	287	78% 22%
1	q	287	78% 22%
2	B	260	38% 70% 22% 8%
2	E	260	22% 68% 24% 8%
2	H	260	9% 68% 23% 8%
2	K	260	33% 69% 23% 8%
2	N	260	7% 68% 24% 8%
2	Q	260	11% 67% 25% 8%
2	T	260	69% 23% 8%
2	W	260	74% 69% 22% 8%
2	Z	260	74% 18% 8%
2	c	260	90% 8%
2	f	260	90% 8%
2	i	260	90% 8%
2	l	260	41% 90% 8%
2	o	260	90% 8%
2	r	260	90% 8%
3	C	239	16% 72% 19% 9%
3	F	239	29% 72% 19% 9%
3	I	239	9% 72% 19% 9%

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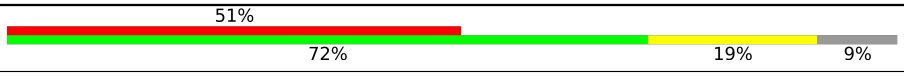
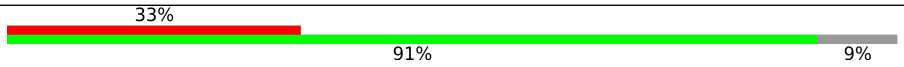
Mol	Chain	Length	Quality of chain
3	L	239	
3	O	239	
3	R	239	
3	U	239	
3	X	239	
3	a	239	
3	d	239	
3	g	239	
3	j	239	
3	m	239	
3	p	239	
3	s	239	

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 77940 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 Echovirus 18 capsid protein 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	224	Total 1741	C 1117	N 305	O 305	S 14	0	0
1	q	224	Total 1741	C 1117	N 305	O 305	S 14	0	0
1	n	224	Total 1741	C 1117	N 305	O 305	S 14	0	0
1	k	224	Total 1741	C 1117	N 305	O 305	S 14	0	0
1	h	224	Total 1741	C 1117	N 305	O 305	S 14	0	0
1	e	224	Total 1741	C 1117	N 305	O 305	S 14	0	0
1	b	224	Total 1741	C 1117	N 305	O 305	S 14	0	0
1	Y	224	Total 1741	C 1117	N 305	O 305	S 14	0	0
1	V	224	Total 1741	C 1117	N 305	O 305	S 14	0	0
1	S	224	Total 1741	C 1117	N 305	O 305	S 14	0	0
1	P	224	Total 1741	C 1117	N 305	O 305	S 14	0	0
1	M	224	Total 1741	C 1117	N 305	O 305	S 14	0	0
1	J	224	Total 1741	C 1117	N 305	O 305	S 14	0	0
1	G	224	Total 1741	C 1117	N 305	O 305	S 14	0	0
1	D	224	Total 1741	C 1117	N 305	O 305	S 14	0	0

- Molecule 2 is a protein called Echovirus 18 capsid protein 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	r	238	Total	C	N	O	S	0	0
			1821	1172	302	332	15		
2	o	238	Total	C	N	O	S	0	0
			1821	1172	302	332	15		
2	l	238	Total	C	N	O	S	0	0
			1821	1172	302	332	15		
2	i	238	Total	C	N	O	S	0	0
			1821	1172	302	332	15		
2	f	238	Total	C	N	O	S	0	0
			1821	1172	302	332	15		
2	c	238	Total	C	N	O	S	0	0
			1821	1172	302	332	15		
2	Z	238	Total	C	N	O	S	0	0
			1821	1172	302	332	15		
2	W	238	Total	C	N	O	S	0	0
			1821	1172	302	332	15		
2	T	238	Total	C	N	O	S	0	0
			1821	1172	302	332	15		
2	Q	238	Total	C	N	O	S	0	0
			1821	1172	302	332	15		
2	N	238	Total	C	N	O	S	0	0
			1821	1172	302	332	15		
2	K	238	Total	C	N	O	S	0	0
			1821	1172	302	332	15		
2	H	238	Total	C	N	O	S	0	0
			1821	1172	302	332	15		
2	E	238	Total	C	N	O	S	0	0
			1821	1172	302	332	15		
2	B	238	Total	C	N	O	S	0	0
			1821	1172	302	332	15		

- Molecule 3 is a protein called Echovirus 18 capsid protein 3.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	s	218	Total	C	N	O	S	0	0
			1634	1045	265	308	16		
3	p	218	Total	C	N	O	S	0	0
			1634	1045	265	308	16		
3	m	218	Total	C	N	O	S	0	0
			1634	1045	265	308	16		
3	j	218	Total	C	N	O	S	0	0
			1634	1045	265	308	16		
3	g	218	Total	C	N	O	S	0	0
			1634	1045	265	308	16		

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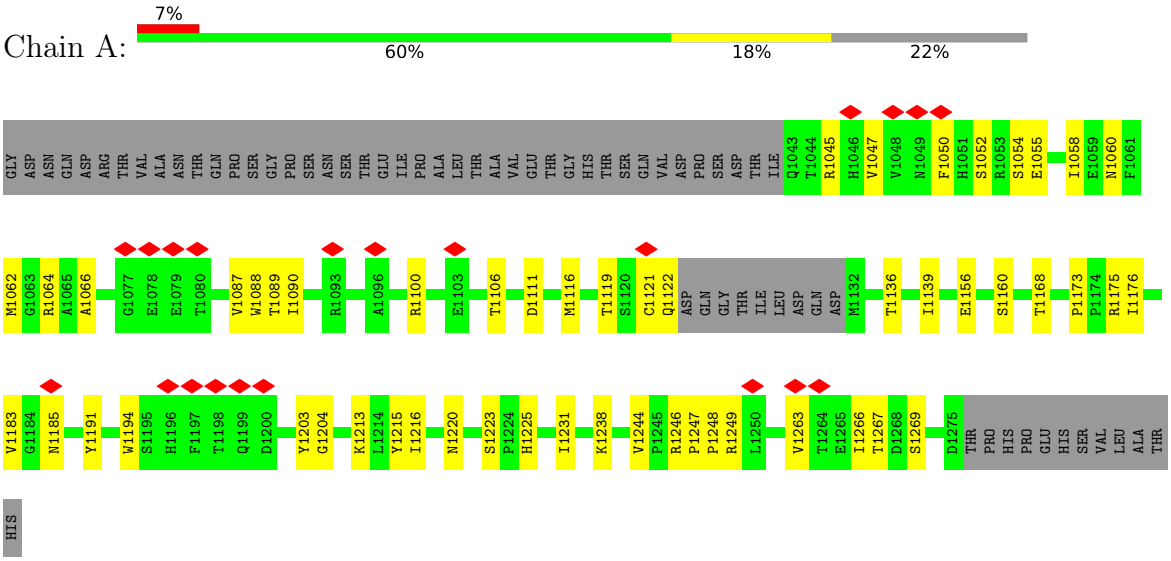
Mol	Chain	Residues	Atoms					AltConf	Trace
3	d	218	Total	C	N	O	S	0	0
			1634	1045	265	308	16		
3	a	218	Total	C	N	O	S	0	0
			1634	1045	265	308	16		
3	X	218	Total	C	N	O	S	0	0
			1634	1045	265	308	16		
3	U	218	Total	C	N	O	S	0	0
			1634	1045	265	308	16		
3	R	218	Total	C	N	O	S	0	0
			1634	1045	265	308	16		
3	O	218	Total	C	N	O	S	0	0
			1634	1045	265	308	16		
3	L	218	Total	C	N	O	S	0	0
			1634	1045	265	308	16		
3	I	218	Total	C	N	O	S	0	0
			1634	1045	265	308	16		
3	F	218	Total	C	N	O	S	0	0
			1634	1045	265	308	16		
3	C	218	Total	C	N	O	S	0	0
			1634	1045	265	308	16		

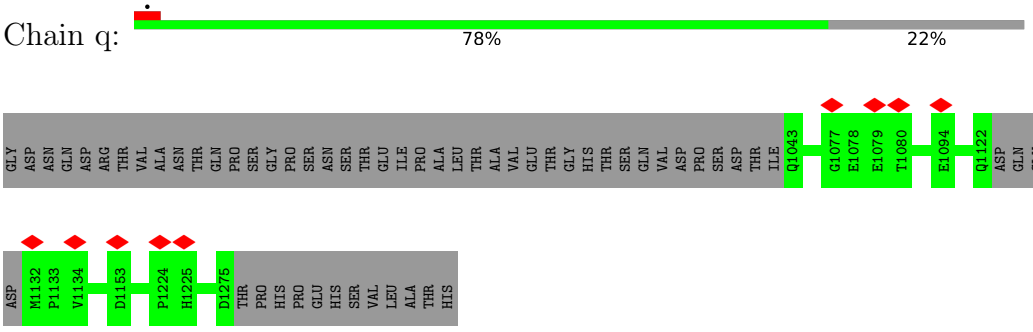
3 Residue-property plots (i)

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.

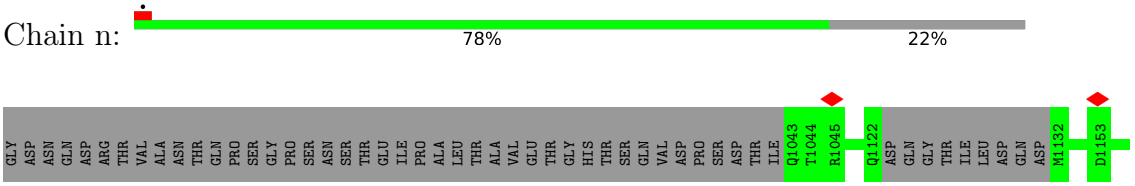
• Molecule 1: Echovirus 18 capsid protein 1

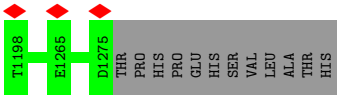


• Molecule 1: Echovirus 18 capsid protein 1



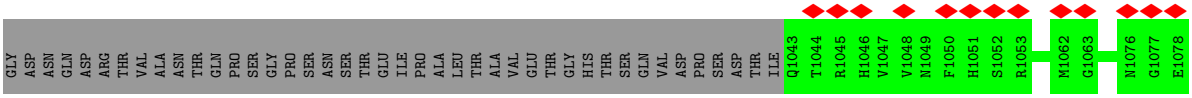
• Molecule 1: Echovirus 18 capsid protein 1

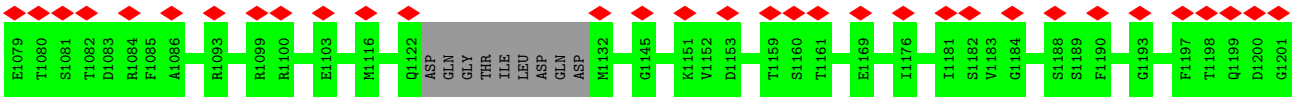


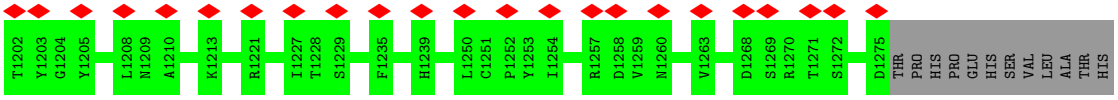


• Molecule 1: Echovirus 18 capsid protein 1

Chain k:

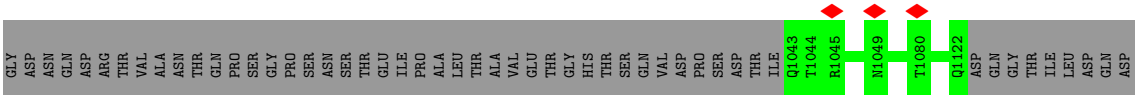


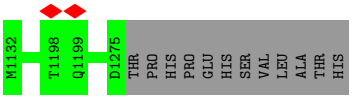




• Molecule 1: Echovirus 18 capsid protein 1

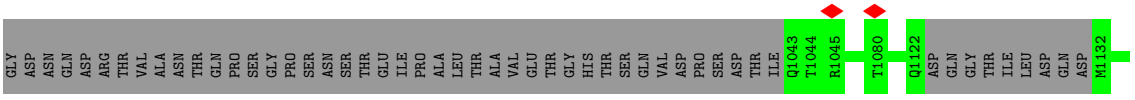
Chain h:

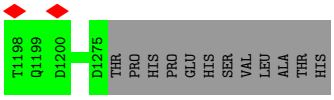




• Molecule 1: Echovirus 18 capsid protein 1

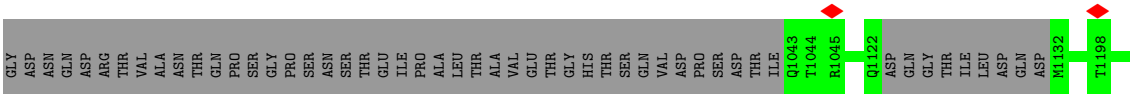
Chain e:

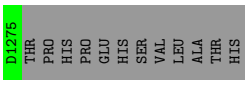




• Molecule 1: Echovirus 18 capsid protein 1

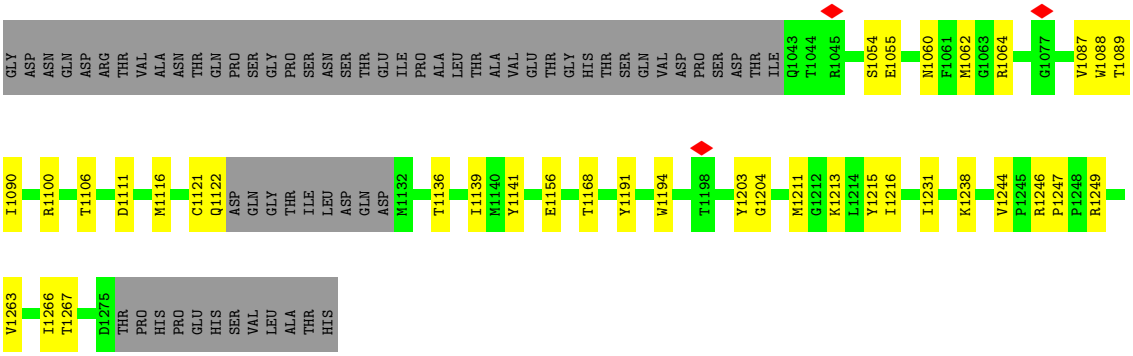
Chain b:





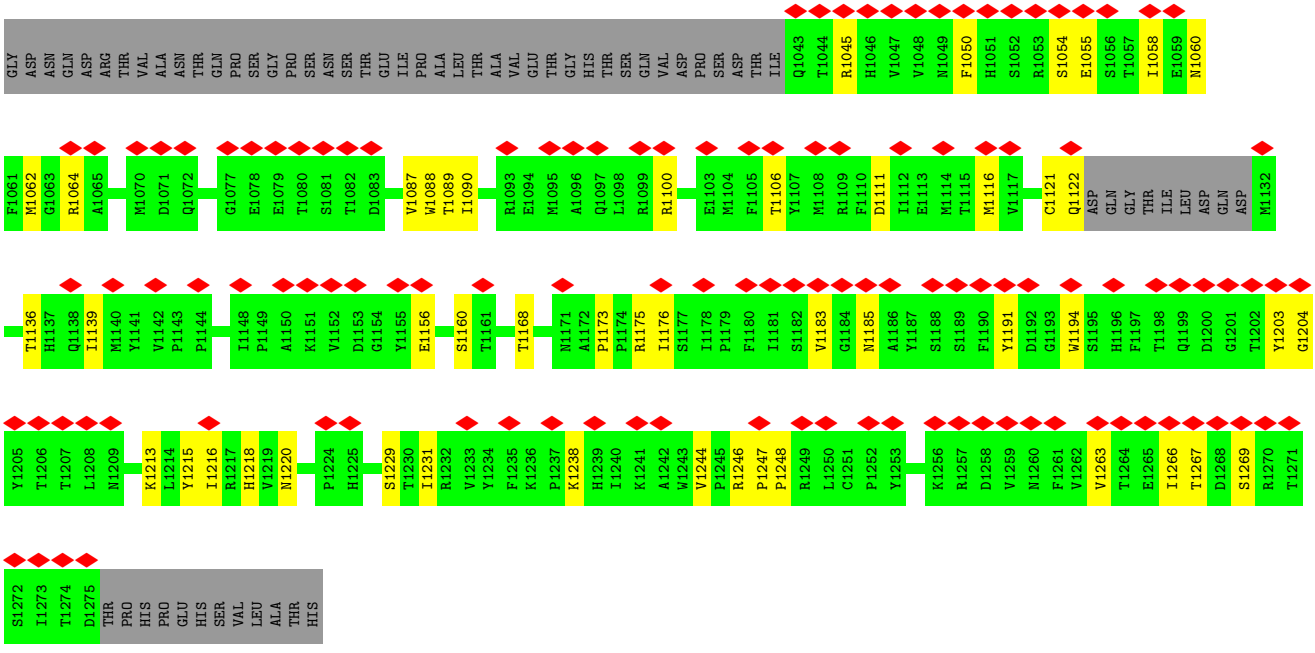
• Molecule 1: Echovirus 18 capsid protein 1

Chain Y:



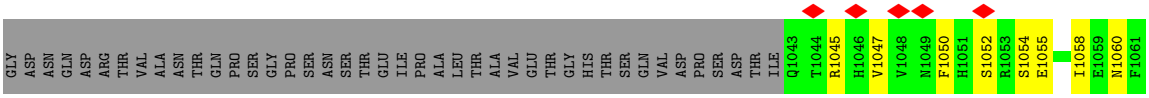
• Molecule 1: Echovirus 18 capsid protein 1

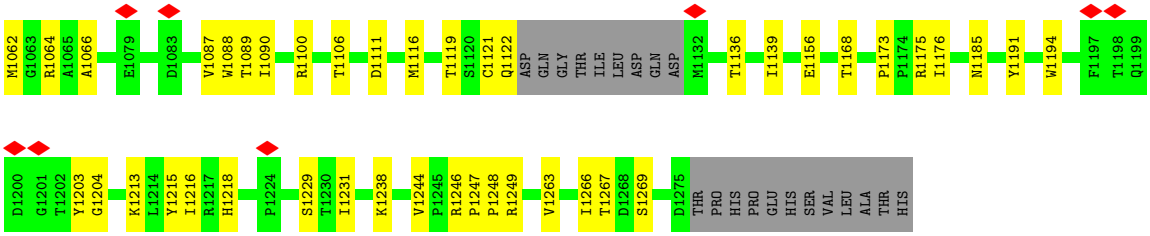
Chain V:



• Molecule 1: Echovirus 18 capsid protein 1

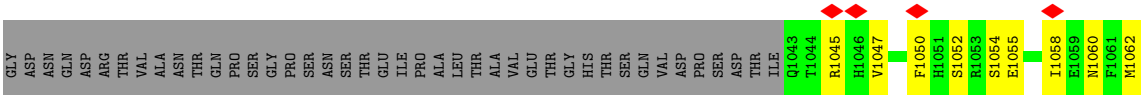
Chain S:

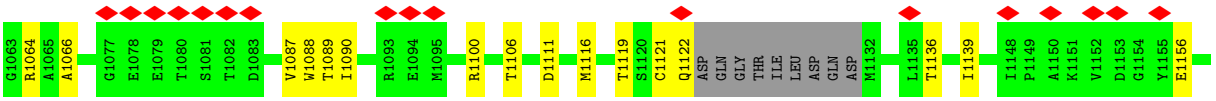


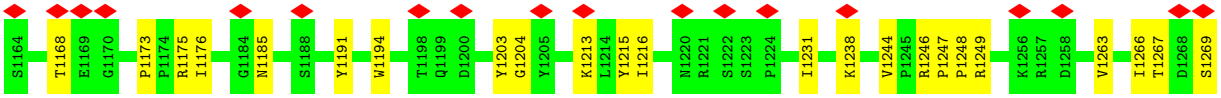


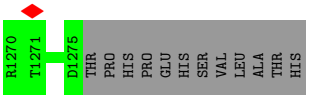
• Molecule 1: Echovirus 18 capsid protein 1

Chain P:



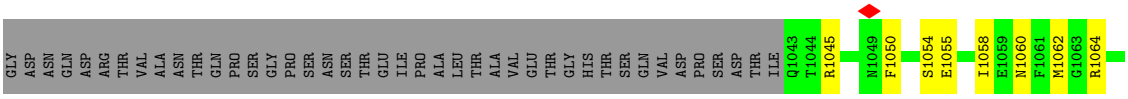


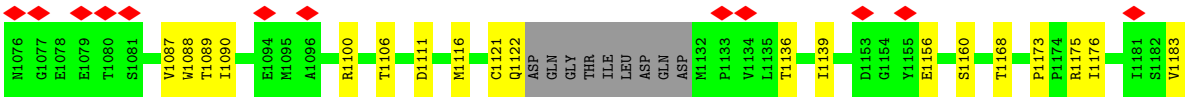


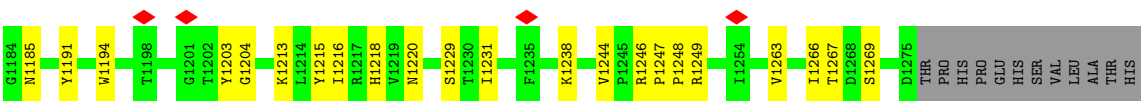


• Molecule 1: Echovirus 18 capsid protein 1

Chain M:

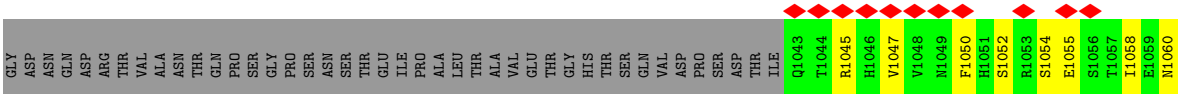


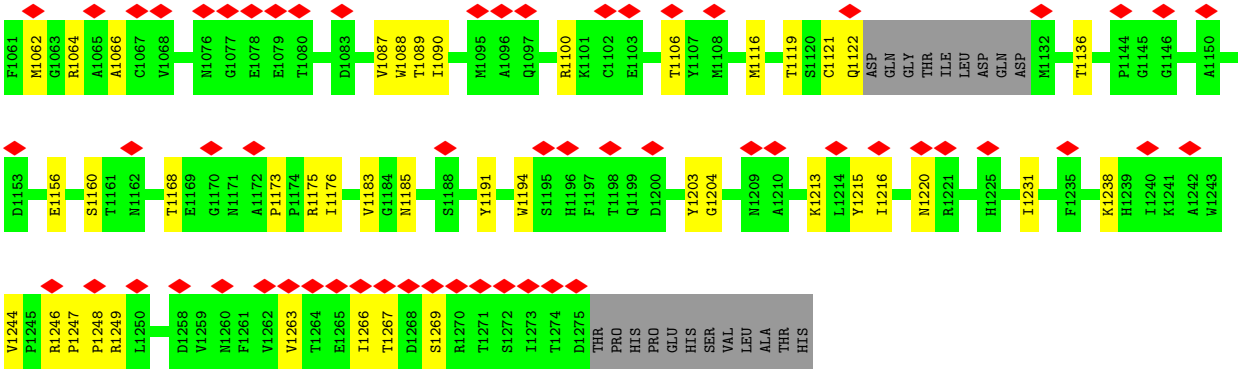




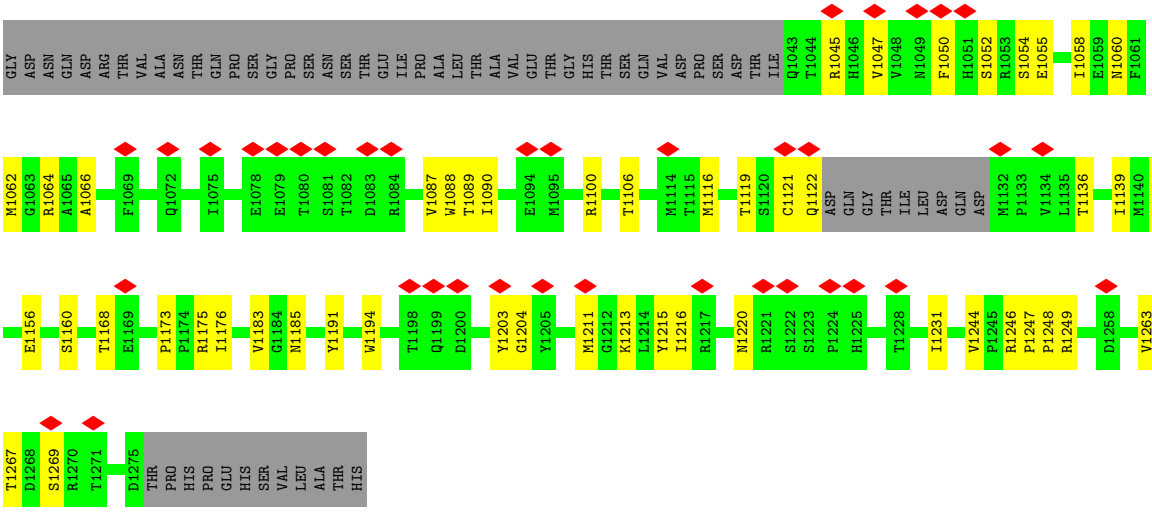
• Molecule 1: Echovirus 18 capsid protein 1

Chain J:

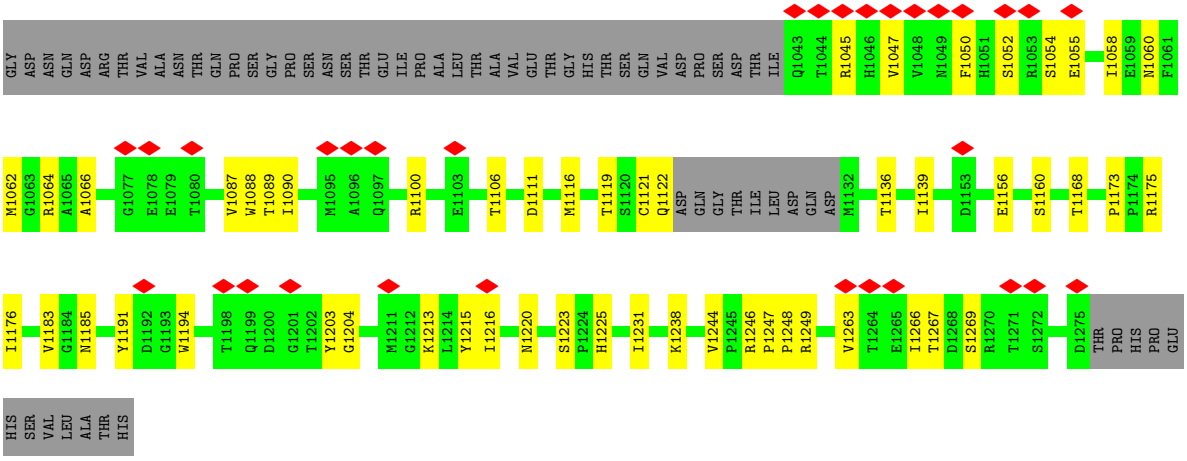




• Molecule 1: Echovirus 18 capsid protein 1

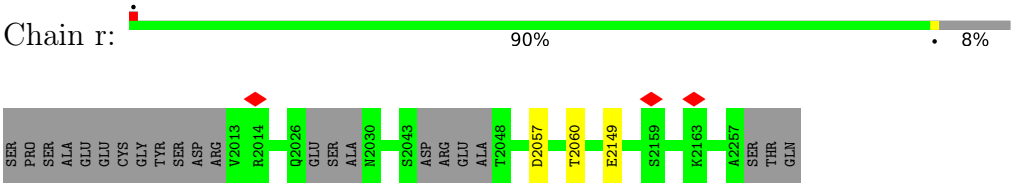


• Molecule 1: Echovirus 18 capsid protein 1



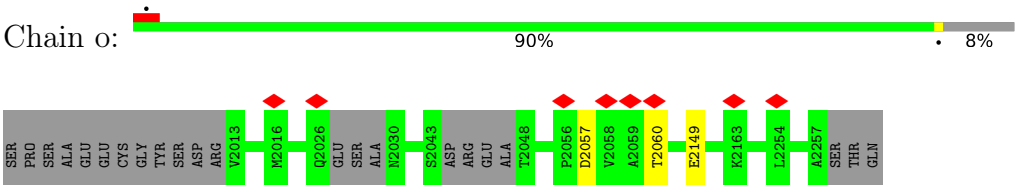
• Molecule 2: Echovirus 18 capsid protein 2

Chain r:



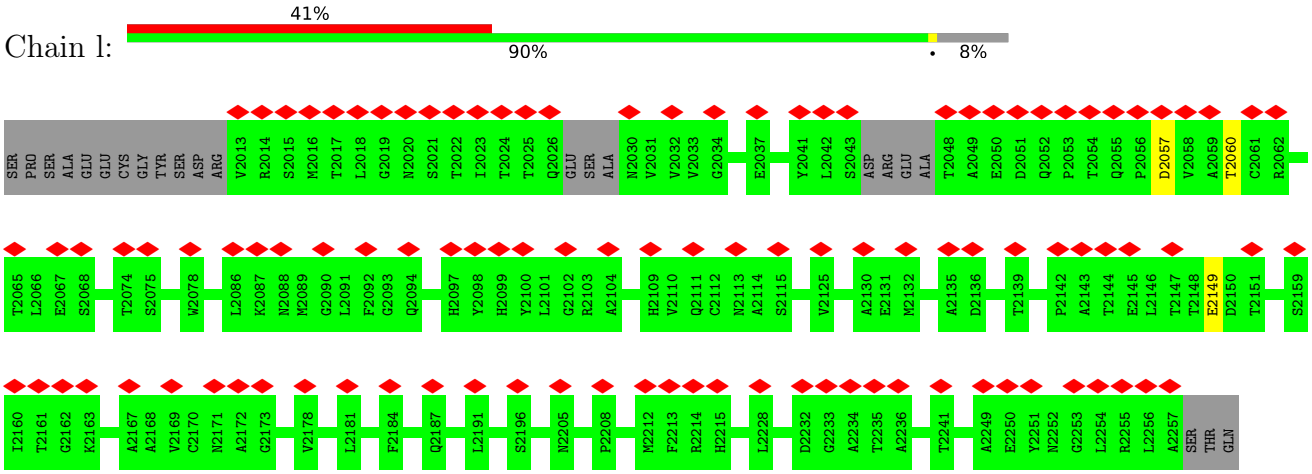
• Molecule 2: Echovirus 18 capsid protein 2

Chain o:



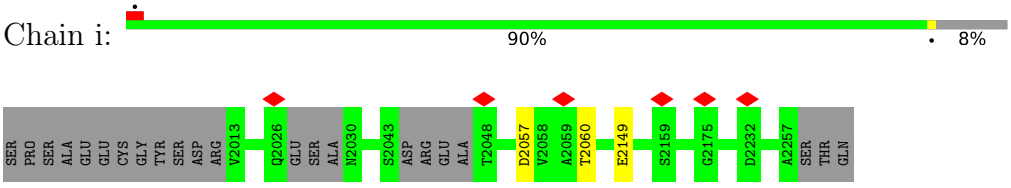
• Molecule 2: Echovirus 18 capsid protein 2

Chain l:



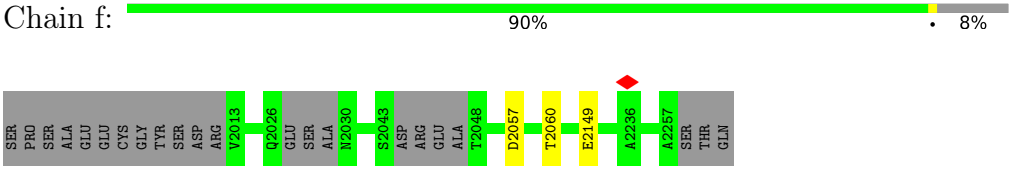
• Molecule 2: Echovirus 18 capsid protein 2

Chain i:



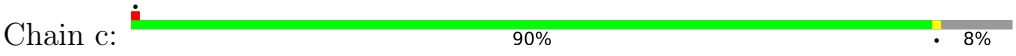
• Molecule 2: Echovirus 18 capsid protein 2

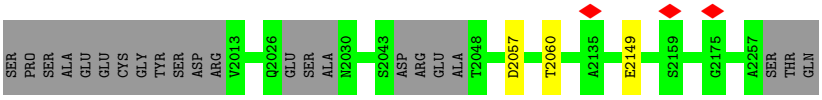
Chain f:



• Molecule 2: Echovirus 18 capsid protein 2

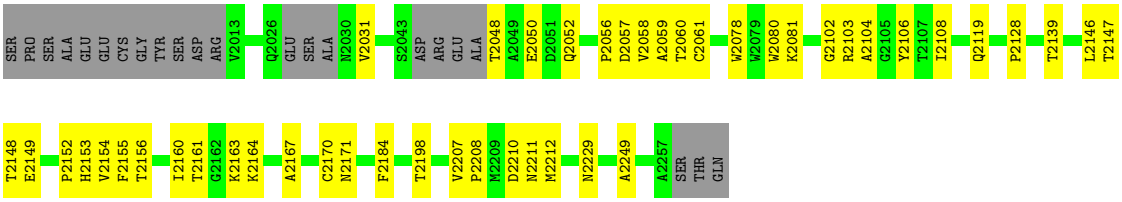
Chain c:





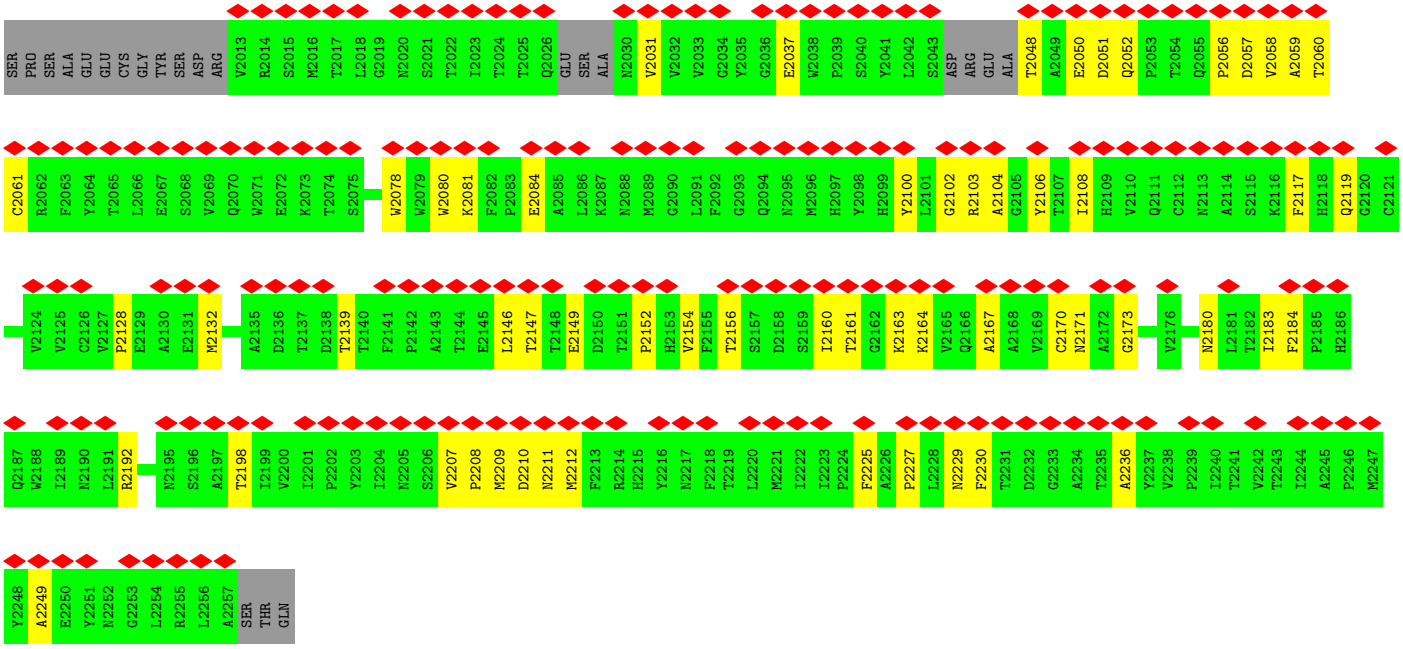
• Molecule 2: Echovirus 18 capsid protein 2

Chain Z:



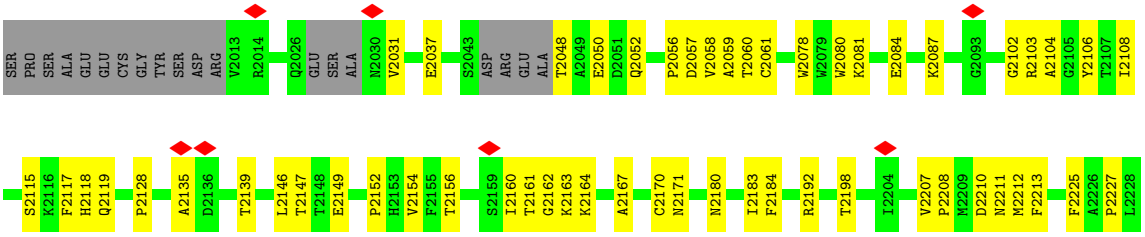
• Molecule 2: Echovirus 18 capsid protein 2

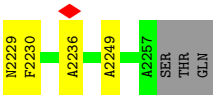
Chain W:



• Molecule 2: Echovirus 18 capsid protein 2

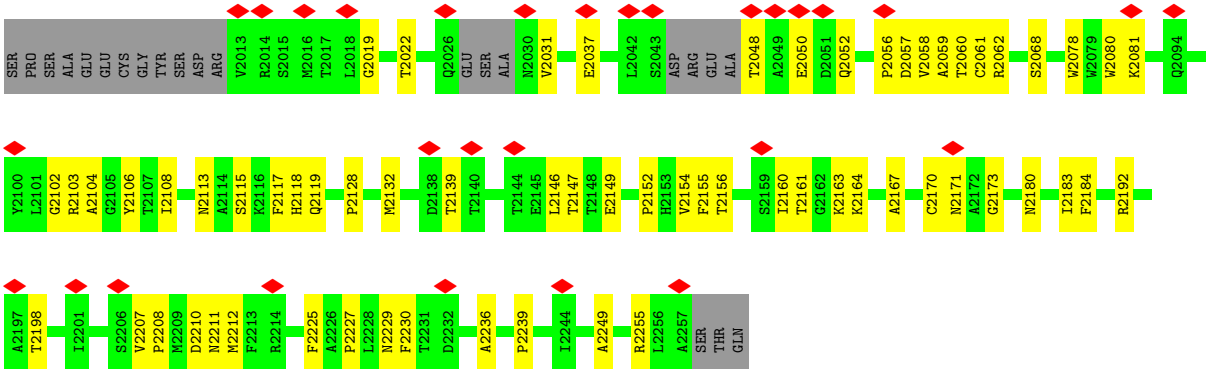
Chain T:





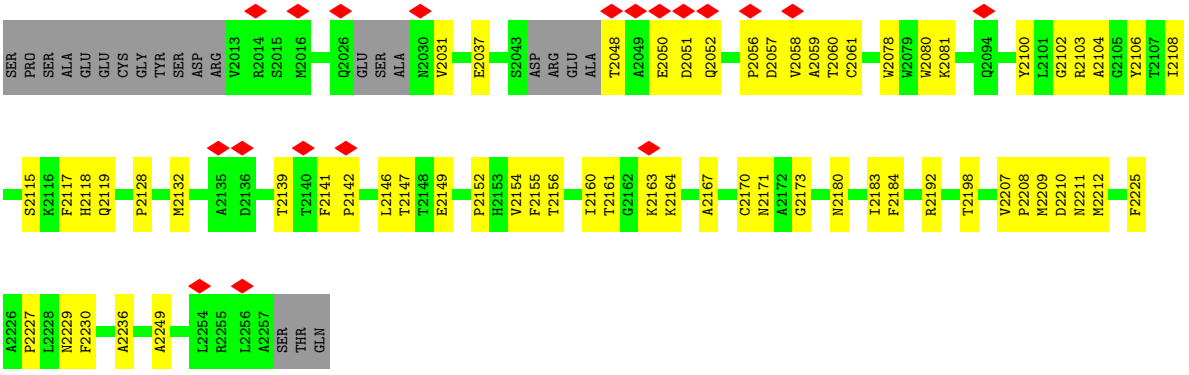
• Molecule 2: Echovirus 18 capsid protein 2

Chain Q:



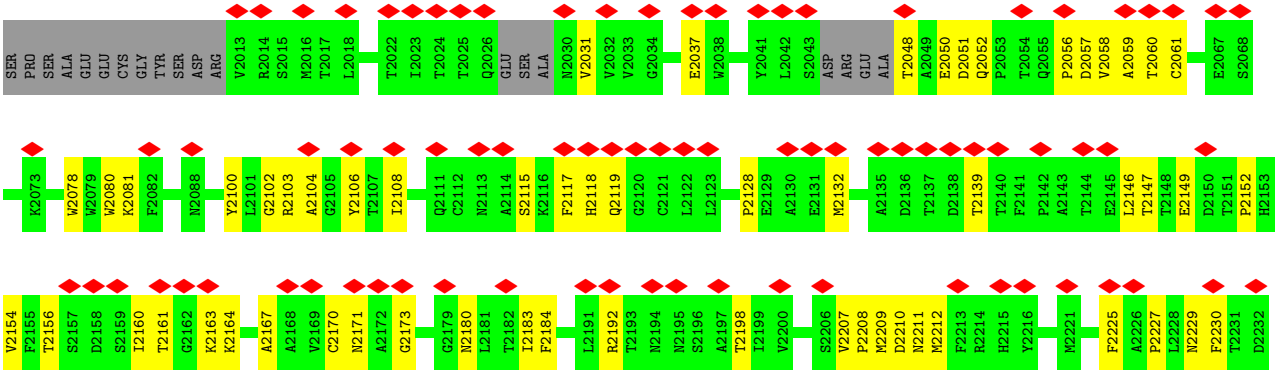
• Molecule 2: Echovirus 18 capsid protein 2

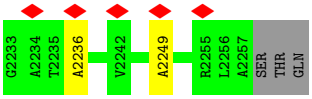
Chain N:



• Molecule 2: Echovirus 18 capsid protein 2

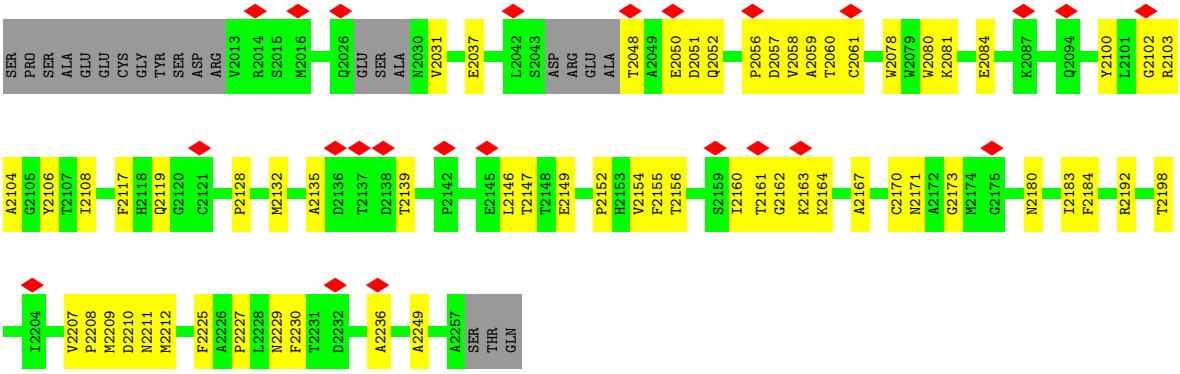
Chain K:





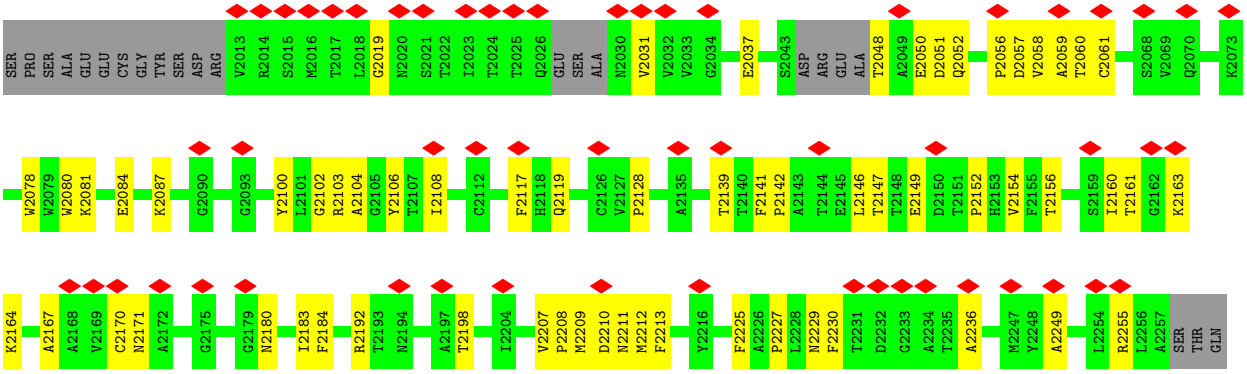
• Molecule 2: Echovirus 18 capsid protein 2

Chain H:



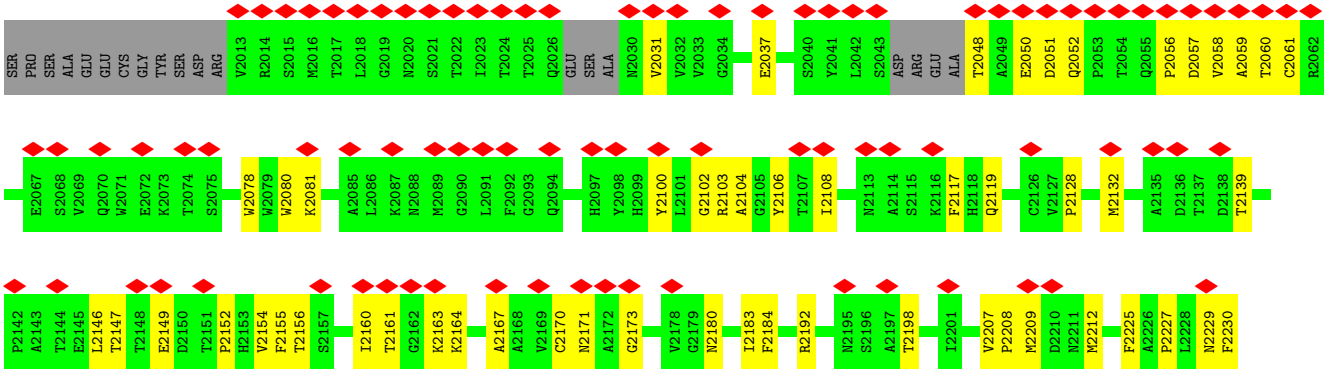
• Molecule 2: Echovirus 18 capsid protein 2

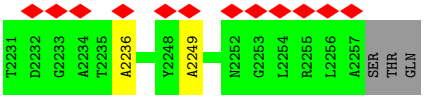
Chain E:



• Molecule 2: Echovirus 18 capsid protein 2

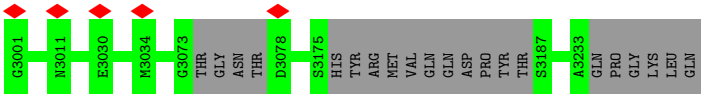
Chain B:





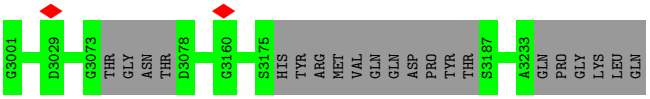
• Molecule 3: Echovirus 18 capsid protein 3

Chain s:



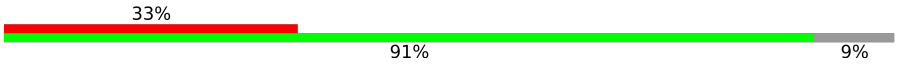
• Molecule 3: Echovirus 18 capsid protein 3

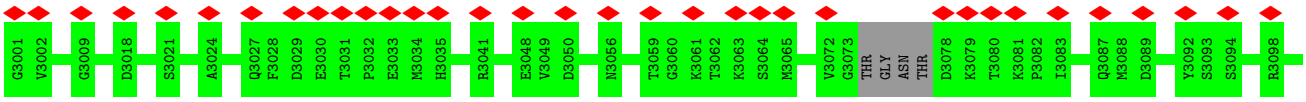
Chain p:

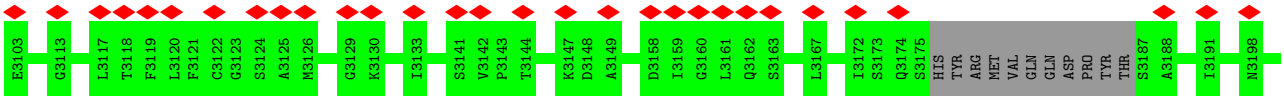


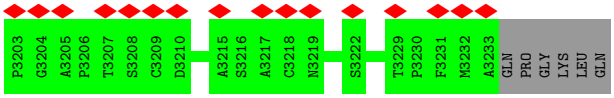
• Molecule 3: Echovirus 18 capsid protein 3

Chain m:



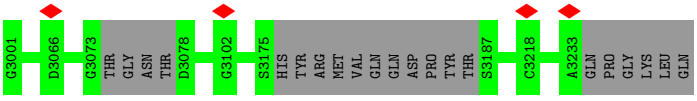






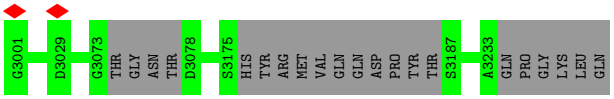
• Molecule 3: Echovirus 18 capsid protein 3

Chain j:



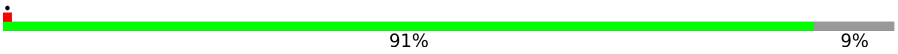
• Molecule 3: Echovirus 18 capsid protein 3

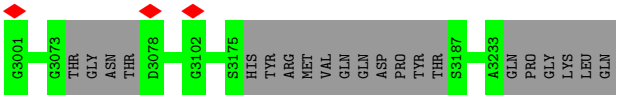
Chain g:



• Molecule 3: Echovirus 18 capsid protein 3

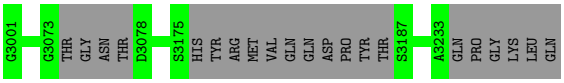
Chain d:





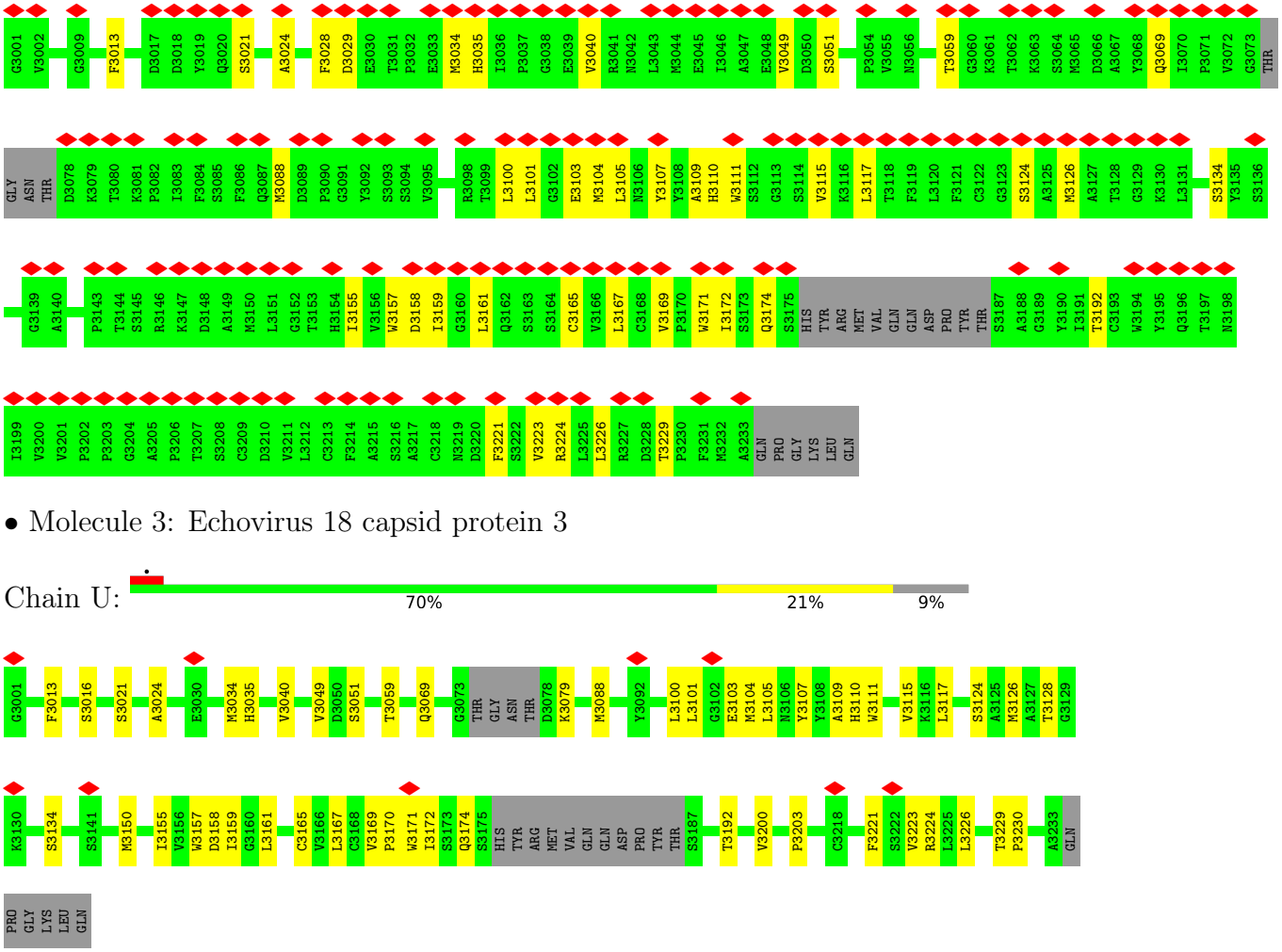
• Molecule 3: Echovirus 18 capsid protein 3

Chain a:



• Molecule 3: Echovirus 18 capsid protein 3

Chain X:

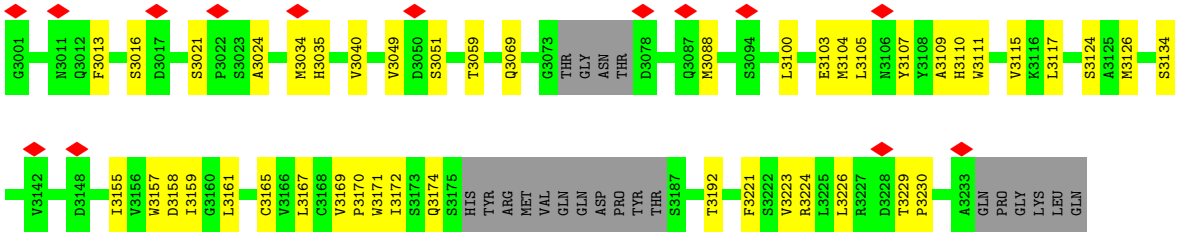


• Molecule 3: Echovirus 18 capsid protein 3

Chain U:

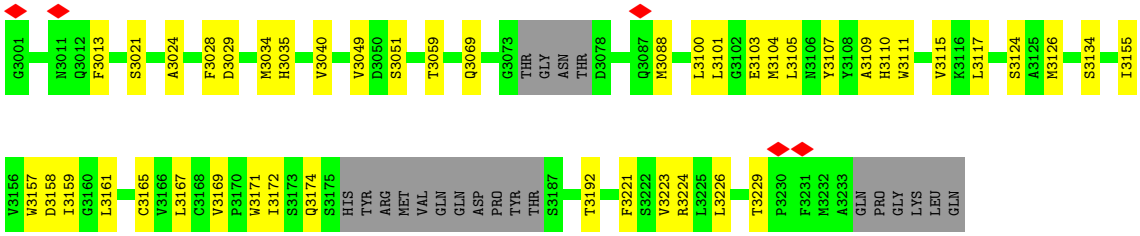
• Molecule 3: Echovirus 18 capsid protein 3

Chain R:



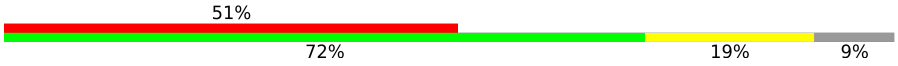
• Molecule 3: Echovirus 18 capsid protein 3

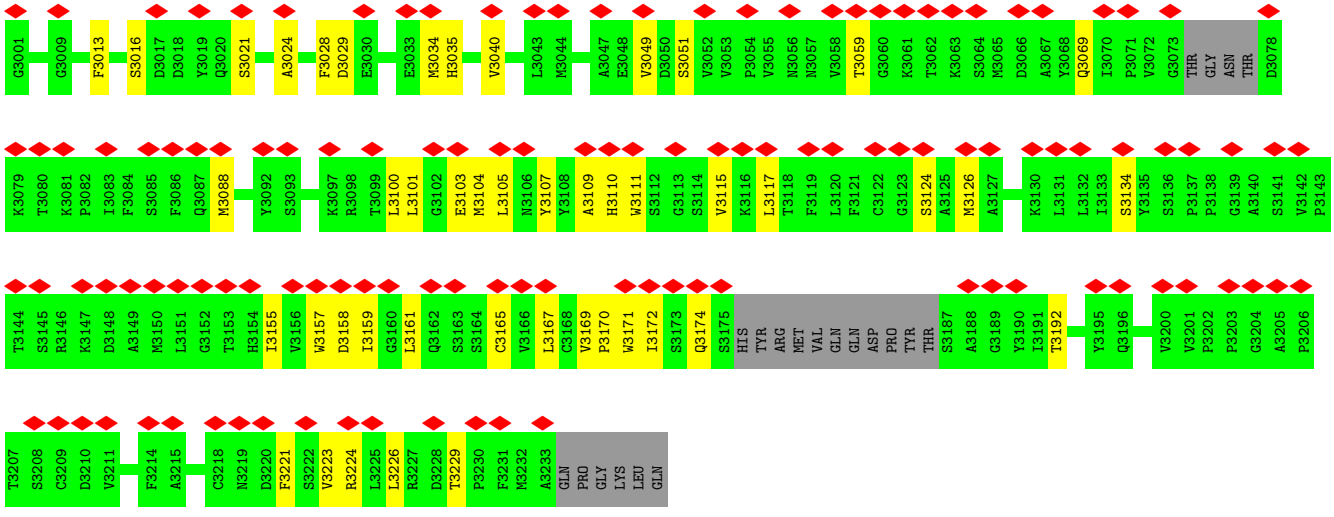
Chain O:



• Molecule 3: Echovirus 18 capsid protein 3

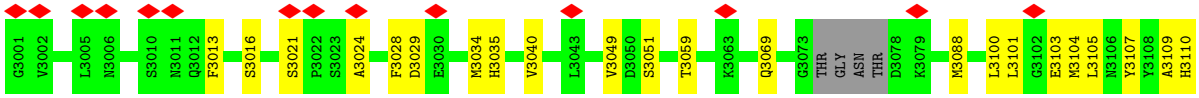
Chain L:

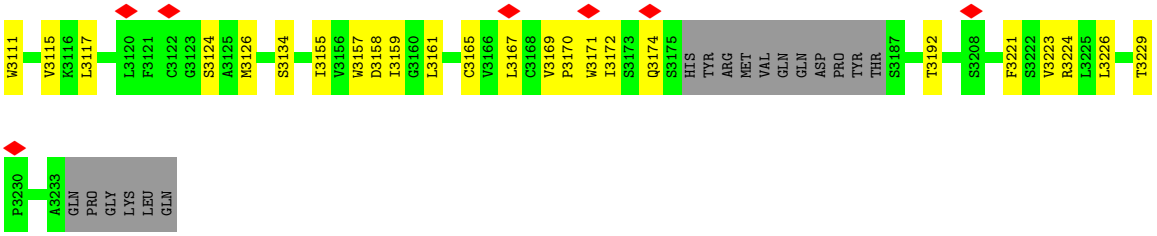




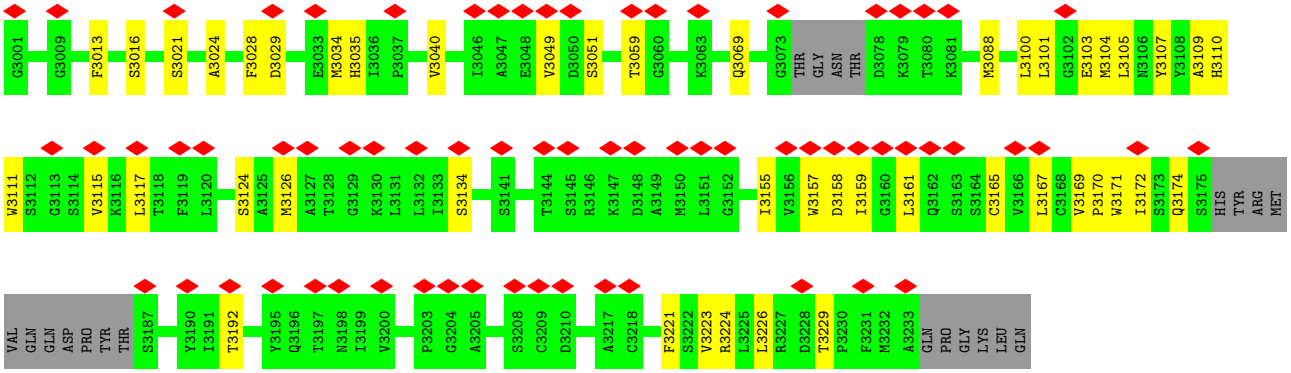
• Molecule 3: Echovirus 18 capsid protein 3

Chain I:

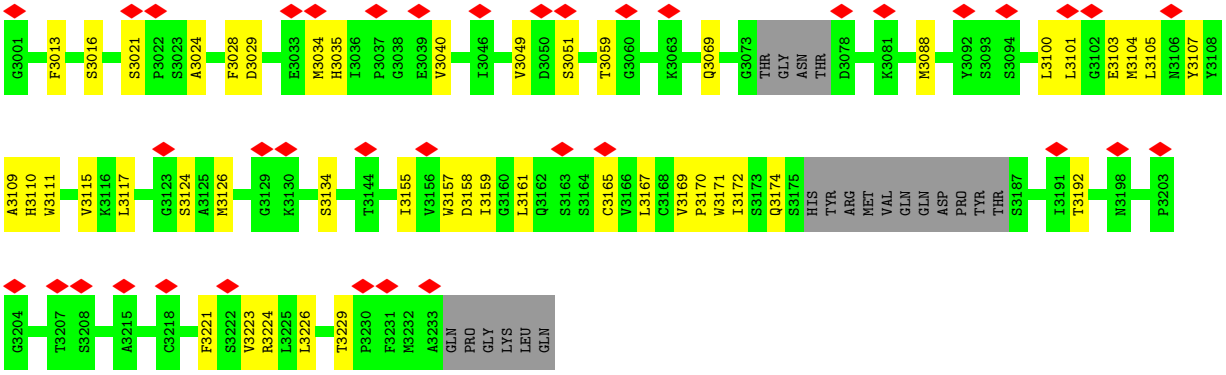




• Molecule 3: Echovirus 18 capsid protein 3



• Molecule 3: Echovirus 18 capsid protein 3



4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C3	Depositor
Number of particles used	13635	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING ONLY	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/Å^2$)	45.2	Depositor
Minimum defocus (nm)	651	Depositor
Maximum defocus (nm)	3282	Depositor
Magnification	79575	Depositor
Image detector	FEI FALCON III (4k x 4k)	Depositor
Maximum map value	0.568	Depositor
Minimum map value	-0.289	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.016	Depositor
Recommended contour level	0.066	Depositor
Map size (Å)	543.232, 543.232, 543.232	wwPDB
Map dimensions	512, 512, 512	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.061, 1.061, 1.061	Depositor

5 Model quality (i)

5.1 Standard geometry (i)

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	#$\|Z\|$ >5	RMSZ	#$\|Z\|$ >5
1	A	0.27	0/1791	0.47	0/2442
1	D	0.27	0/1791	0.47	0/2442
1	G	0.27	0/1791	0.47	0/2442
1	J	0.27	0/1791	0.47	0/2442
1	M	0.27	0/1791	0.47	0/2442
1	P	0.27	0/1791	0.47	0/2442
1	S	0.27	0/1791	0.47	0/2442
1	V	0.27	0/1791	0.47	0/2442
1	Y	0.27	0/1791	0.47	0/2442
1	b	0.27	0/1791	0.47	0/2442
1	e	0.27	0/1791	0.47	0/2442
1	h	0.27	0/1791	0.47	0/2442
1	k	0.27	0/1791	0.47	0/2442
1	n	0.27	0/1791	0.47	0/2442
1	q	0.27	0/1791	0.47	0/2442
2	B	0.27	0/1874	0.53	0/2567
2	E	0.27	0/1874	0.53	0/2567
2	H	0.27	0/1874	0.53	0/2567
2	K	0.26	0/1874	0.53	0/2567
2	N	0.27	0/1874	0.53	0/2567
2	Q	0.27	0/1874	0.53	0/2567
2	T	0.27	0/1874	0.53	0/2567
2	W	0.27	0/1874	0.53	0/2567
2	Z	0.27	0/1874	0.53	0/2567
2	c	0.27	0/1874	0.53	0/2567
2	f	0.27	0/1874	0.53	0/2567
2	i	0.27	0/1874	0.53	0/2567
2	l	0.27	0/1874	0.53	0/2567
2	o	0.26	0/1874	0.53	0/2567
2	r	0.26	0/1874	0.53	0/2567
3	C	0.26	0/1676	0.48	0/2287
3	F	0.26	0/1676	0.48	0/2287
3	I	0.26	0/1676	0.48	0/2287
3	L	0.26	0/1676	0.48	0/2287

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\#\|Z\| > 5$	RMSZ	$\#\|Z\| > 5$
3	O	0.26	0/1676	0.48	0/2287
3	R	0.26	0/1676	0.48	0/2287
3	U	0.26	0/1676	0.48	0/2287
3	X	0.26	0/1676	0.48	0/2287
3	a	0.26	0/1676	0.48	0/2287
3	d	0.26	0/1676	0.48	0/2287
3	g	0.26	0/1676	0.48	0/2287
3	j	0.26	0/1676	0.48	0/2287
3	m	0.26	0/1676	0.48	0/2287
3	p	0.26	0/1676	0.48	0/2287
3	s	0.26	0/1676	0.48	0/2287
All	All	0.27	0/80115	0.49	0/109440

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1741	0	1659	46	0
1	D	1741	0	1659	46	0
1	G	1741	0	1659	45	0
1	J	1741	0	1659	44	0
1	M	1741	0	1659	41	0
1	P	1741	0	1659	42	0
1	S	1741	0	1659	43	0
1	V	1741	0	1659	40	0
1	Y	1741	0	1659	26	0
1	b	1741	0	1659	0	0
1	e	1741	0	1659	0	0
1	h	1741	0	1659	0	0
1	k	1741	0	1659	0	0
1	n	1741	0	1659	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	q	1741	0	1659	0	0
2	B	1821	0	1739	42	0
2	E	1821	0	1739	47	0
2	H	1821	0	1739	45	0
2	K	1821	0	1739	43	0
2	N	1821	0	1739	46	0
2	Q	1821	0	1739	48	0
2	T	1821	0	1739	41	0
2	W	1821	0	1739	43	0
2	Z	1821	0	1739	27	0
2	c	1821	0	1739	0	0
2	f	1821	0	1739	0	0
2	i	1821	0	1739	0	0
2	l	1821	0	1739	0	0
2	o	1821	0	1739	0	0
2	r	1821	0	1739	0	0
3	C	1634	0	1596	49	0
3	F	1634	0	1596	48	0
3	I	1634	0	1596	49	0
3	L	1634	0	1596	49	0
3	O	1634	0	1596	43	0
3	R	1634	0	1596	44	0
3	U	1634	0	1596	54	0
3	X	1634	0	1596	43	0
3	a	1634	0	1596	0	0
3	d	1634	0	1596	0	0
3	g	1634	0	1596	0	0
3	j	1634	0	1596	0	0
3	m	1634	0	1596	0	0
3	p	1634	0	1596	0	0
3	s	1634	0	1596	0	0
All	All	77940	0	74910	862	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 10.

The worst 5 of 862 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:V:1216:ILE:HD13	1:V:1231:ILE:HD13	1.56	0.88
1:A:1216:ILE:HD13	1:A:1231:ILE:HD13	1.56	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:S:1216:ILE:HD13	1:S:1231:ILE:HD13	1.56	0.87
1:G:1216:ILE:HD13	1:G:1231:ILE:HD13	1.56	0.87
1:D:1216:ILE:HD13	1:D:1231:ILE:HD13	1.56	0.87

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	220/287 (77%)	200 (91%)	20 (9%)	0	100	100
1	D	220/287 (77%)	200 (91%)	20 (9%)	0	100	100
1	G	220/287 (77%)	200 (91%)	20 (9%)	0	100	100
1	J	220/287 (77%)	200 (91%)	20 (9%)	0	100	100
1	M	220/287 (77%)	199 (90%)	21 (10%)	0	100	100
1	P	220/287 (77%)	200 (91%)	20 (9%)	0	100	100
1	S	220/287 (77%)	200 (91%)	20 (9%)	0	100	100
1	V	220/287 (77%)	200 (91%)	20 (9%)	0	100	100
1	Y	220/287 (77%)	200 (91%)	20 (9%)	0	100	100
1	b	220/287 (77%)	200 (91%)	20 (9%)	0	100	100
1	e	220/287 (77%)	200 (91%)	20 (9%)	0	100	100
1	h	220/287 (77%)	199 (90%)	21 (10%)	0	100	100
1	k	220/287 (77%)	200 (91%)	20 (9%)	0	100	100
1	n	220/287 (77%)	200 (91%)	20 (9%)	0	100	100
1	q	220/287 (77%)	199 (90%)	21 (10%)	0	100	100
2	B	232/260 (89%)	204 (88%)	25 (11%)	3 (1%)	10	40
2	E	232/260 (89%)	204 (88%)	25 (11%)	3 (1%)	10	40

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	H	232/260 (89%)	204 (88%)	25 (11%)	3 (1%)	10	40
2	K	232/260 (89%)	203 (88%)	26 (11%)	3 (1%)	10	40
2	N	232/260 (89%)	204 (88%)	25 (11%)	3 (1%)	10	40
2	Q	232/260 (89%)	204 (88%)	25 (11%)	3 (1%)	10	40
2	T	232/260 (89%)	204 (88%)	25 (11%)	3 (1%)	10	40
2	W	232/260 (89%)	204 (88%)	25 (11%)	3 (1%)	10	40
2	Z	232/260 (89%)	204 (88%)	25 (11%)	3 (1%)	10	40
2	c	232/260 (89%)	204 (88%)	25 (11%)	3 (1%)	10	40
2	f	232/260 (89%)	204 (88%)	25 (11%)	3 (1%)	10	40
2	i	232/260 (89%)	204 (88%)	25 (11%)	3 (1%)	10	40
2	l	232/260 (89%)	204 (88%)	25 (11%)	3 (1%)	10	40
2	o	232/260 (89%)	204 (88%)	25 (11%)	3 (1%)	10	40
2	r	232/260 (89%)	203 (88%)	26 (11%)	3 (1%)	10	40
3	C	212/239 (89%)	197 (93%)	15 (7%)	0	100	100
3	F	212/239 (89%)	197 (93%)	15 (7%)	0	100	100
3	I	212/239 (89%)	197 (93%)	15 (7%)	0	100	100
3	L	212/239 (89%)	197 (93%)	15 (7%)	0	100	100
3	O	212/239 (89%)	197 (93%)	15 (7%)	0	100	100
3	R	212/239 (89%)	197 (93%)	15 (7%)	0	100	100
3	U	212/239 (89%)	197 (93%)	15 (7%)	0	100	100
3	X	212/239 (89%)	197 (93%)	15 (7%)	0	100	100
3	a	212/239 (89%)	197 (93%)	15 (7%)	0	100	100
3	d	212/239 (89%)	197 (93%)	15 (7%)	0	100	100
3	g	212/239 (89%)	197 (93%)	15 (7%)	0	100	100
3	j	212/239 (89%)	197 (93%)	15 (7%)	0	100	100
3	m	212/239 (89%)	197 (93%)	15 (7%)	0	100	100
3	p	212/239 (89%)	197 (93%)	15 (7%)	0	100	100
3	s	212/239 (89%)	197 (93%)	15 (7%)	0	100	100
All	All	9960/11790 (84%)	9010 (90%)	905 (9%)	45 (0%)	27	57

5 of 45 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	r	2060	THR
2	r	2149	GLU
2	o	2060	THR
2	o	2149	GLU
2	l	2060	THR

5.3.2 Protein sidechains (i)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	182/259 (70%)	182 (100%)	0	100	100
1	D	182/259 (70%)	182 (100%)	0	100	100
1	G	182/259 (70%)	182 (100%)	0	100	100
1	J	182/259 (70%)	182 (100%)	0	100	100
1	M	182/259 (70%)	182 (100%)	0	100	100
1	P	182/259 (70%)	182 (100%)	0	100	100
1	S	182/259 (70%)	182 (100%)	0	100	100
1	V	182/259 (70%)	182 (100%)	0	100	100
1	Y	182/259 (70%)	182 (100%)	0	100	100
1	b	182/259 (70%)	182 (100%)	0	100	100
1	e	182/259 (70%)	182 (100%)	0	100	100
1	h	182/259 (70%)	182 (100%)	0	100	100
1	k	182/259 (70%)	182 (100%)	0	100	100
1	n	182/259 (70%)	182 (100%)	0	100	100
1	q	182/259 (70%)	182 (100%)	0	100	100
2	B	193/221 (87%)	193 (100%)	0	100	100
2	E	193/221 (87%)	193 (100%)	0	100	100
2	H	193/221 (87%)	193 (100%)	0	100	100
2	K	193/221 (87%)	193 (100%)	0	100	100
2	N	193/221 (87%)	193 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	Q	193/221 (87%)	193 (100%)	0	100	100
2	T	193/221 (87%)	193 (100%)	0	100	100
2	W	193/221 (87%)	193 (100%)	0	100	100
2	Z	193/221 (87%)	193 (100%)	0	100	100
2	c	193/221 (87%)	193 (100%)	0	100	100
2	f	193/221 (87%)	193 (100%)	0	100	100
2	i	193/221 (87%)	193 (100%)	0	100	100
2	l	193/221 (87%)	193 (100%)	0	100	100
2	o	193/221 (87%)	193 (100%)	0	100	100
2	r	193/221 (87%)	193 (100%)	0	100	100
3	C	184/209 (88%)	184 (100%)	0	100	100
3	F	184/209 (88%)	184 (100%)	0	100	100
3	I	184/209 (88%)	184 (100%)	0	100	100
3	L	184/209 (88%)	184 (100%)	0	100	100
3	O	184/209 (88%)	184 (100%)	0	100	100
3	R	184/209 (88%)	184 (100%)	0	100	100
3	U	184/209 (88%)	184 (100%)	0	100	100
3	X	184/209 (88%)	184 (100%)	0	100	100
3	a	184/209 (88%)	184 (100%)	0	100	100
3	d	184/209 (88%)	184 (100%)	0	100	100
3	g	184/209 (88%)	184 (100%)	0	100	100
3	j	184/209 (88%)	184 (100%)	0	100	100
3	m	184/209 (88%)	184 (100%)	0	100	100
3	p	184/209 (88%)	184 (100%)	0	100	100
3	s	184/209 (88%)	184 (100%)	0	100	100
All	All	8385/10335 (81%)	8385 (100%)	0	100	100

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

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 51 such sidechains are listed below:

Mol	Chain	Res	Type
2	T	2099	HIS

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Mol	Chain	Res	Type
3	R	3110	HIS
3	C	3110	HIS
3	U	3110	HIS
2	Q	2026	GLN

5.3.3 RNA

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains

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

5.5 Carbohydrates

There are no oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

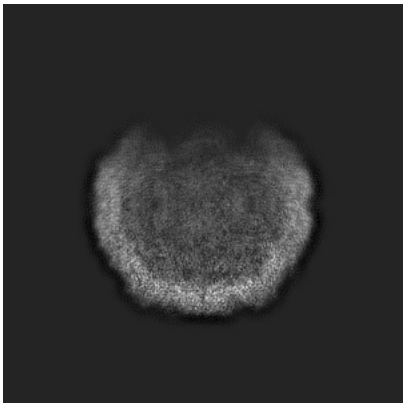
6 Map visualisation

This section contains visualisations of the EMDB entry EMD-0188. These allow visual inspection of the internal detail of the map and identification of artifacts.

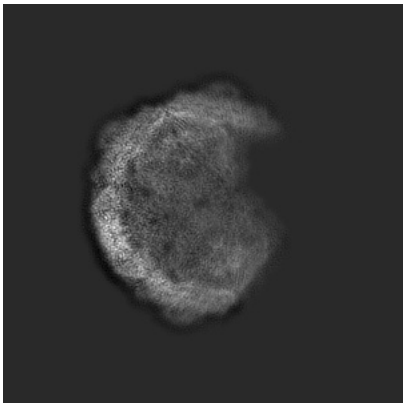
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections

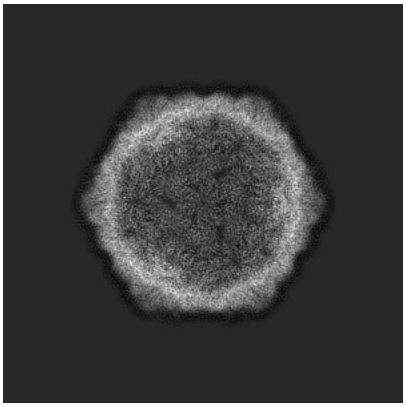
6.1.1 Primary map



X



Y



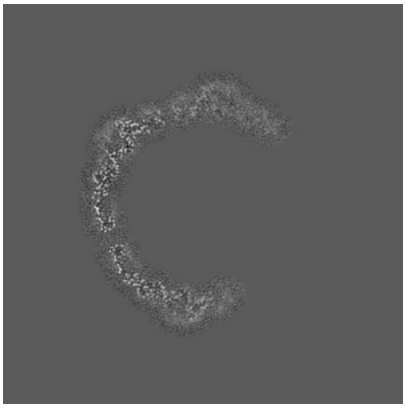
Z

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

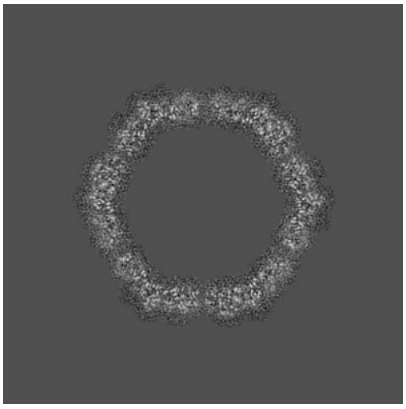
6.2 Central slices

6.2.1 Primary map

X Index: 256



Y Index: 256

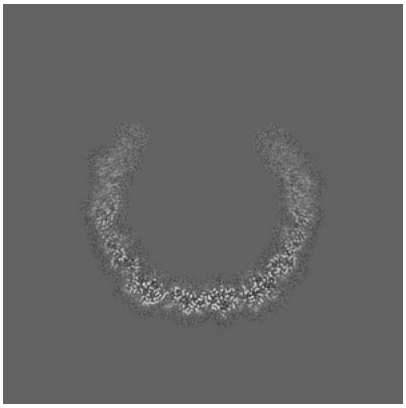


Z Index: 256

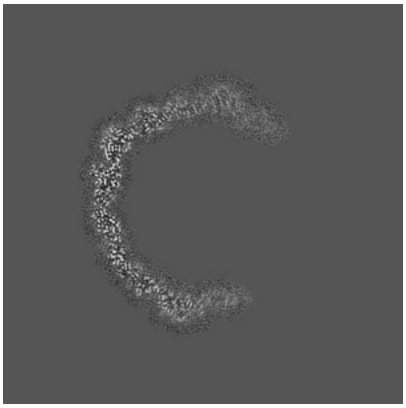
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices (i)

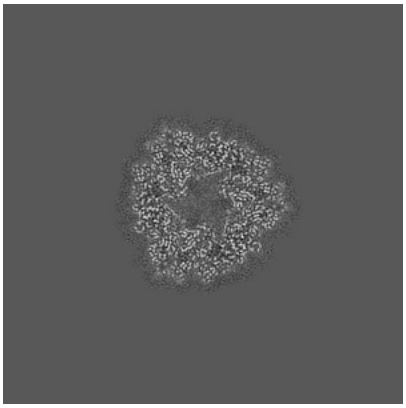
6.3.1 Primary map



X Index: 228



Y Index: 270

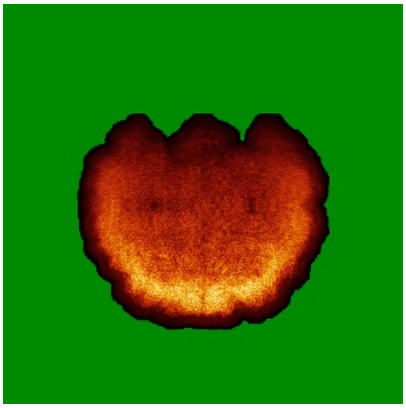


Z Index: 146

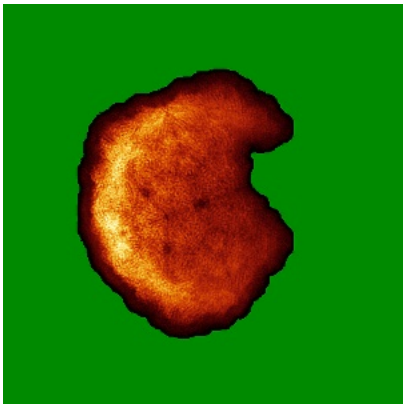
The images above show the largest variance slices of the map in three orthogonal directions.

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

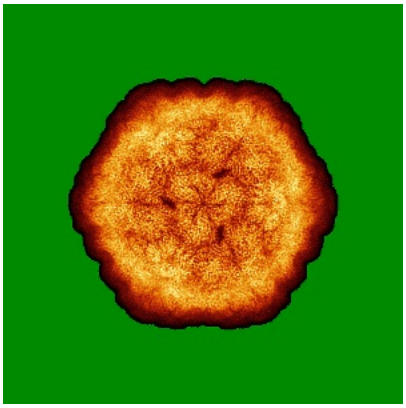
6.4.1 Primary map



X



Y



Z

The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

6.5 Orthogonal surface views (i)

6.5.1 Primary map

X Y Z

The images above show the 3D surface view of the map at the recommended contour level 0.066. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.6 Mask visualisation

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

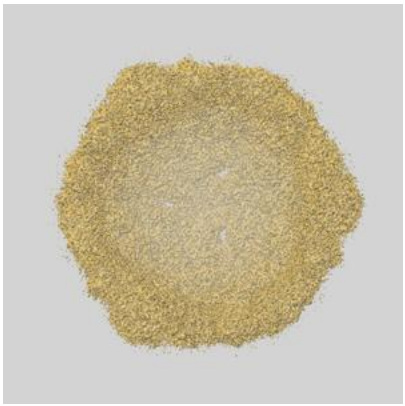
A mask typically either:

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

6.6.1 emd_0188_msk_2.map

X

Y

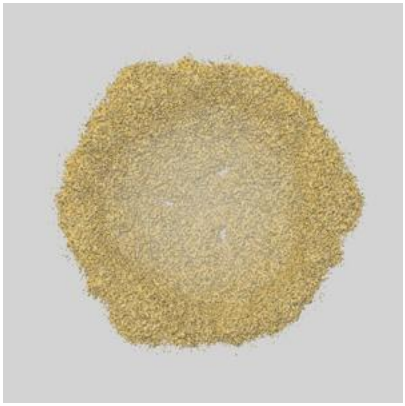


Z

6.6.2 emd_0188_msk_1.map

X

Y

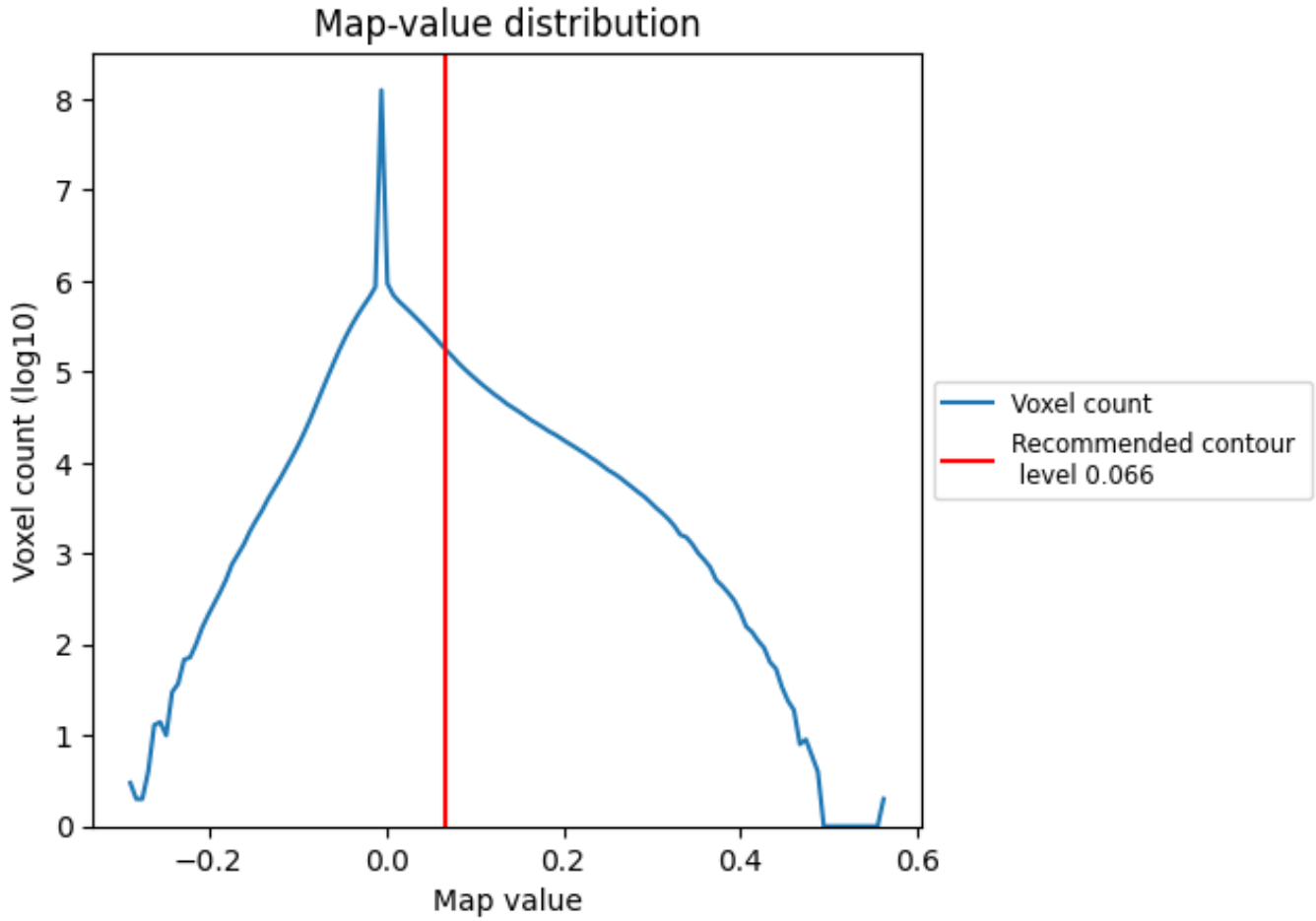


Z

7 Map analysis

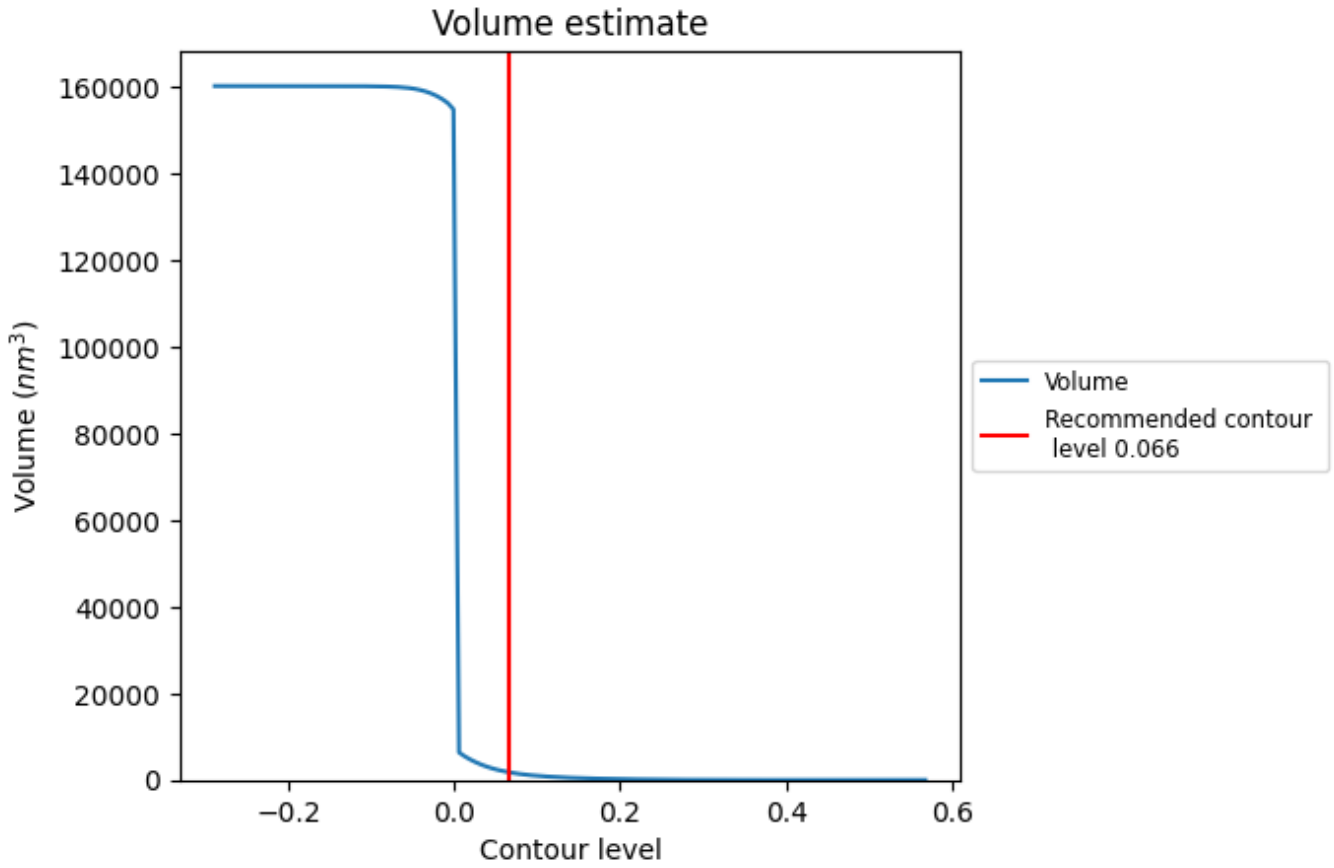
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution



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

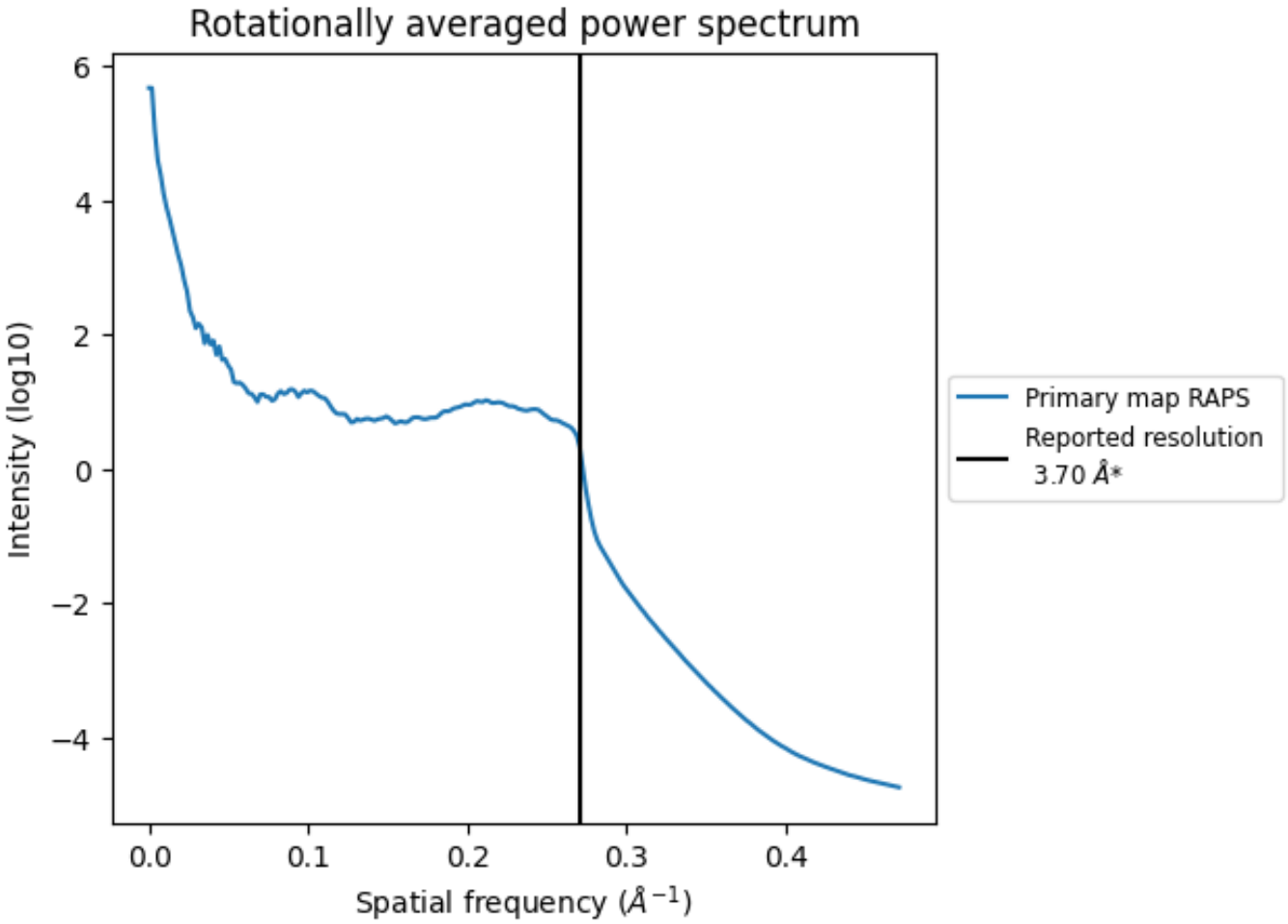
7.2 Volume estimate (i)



The volume at the recommended contour level is 1776 nm^3; this corresponds to an approximate mass of 1605 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)

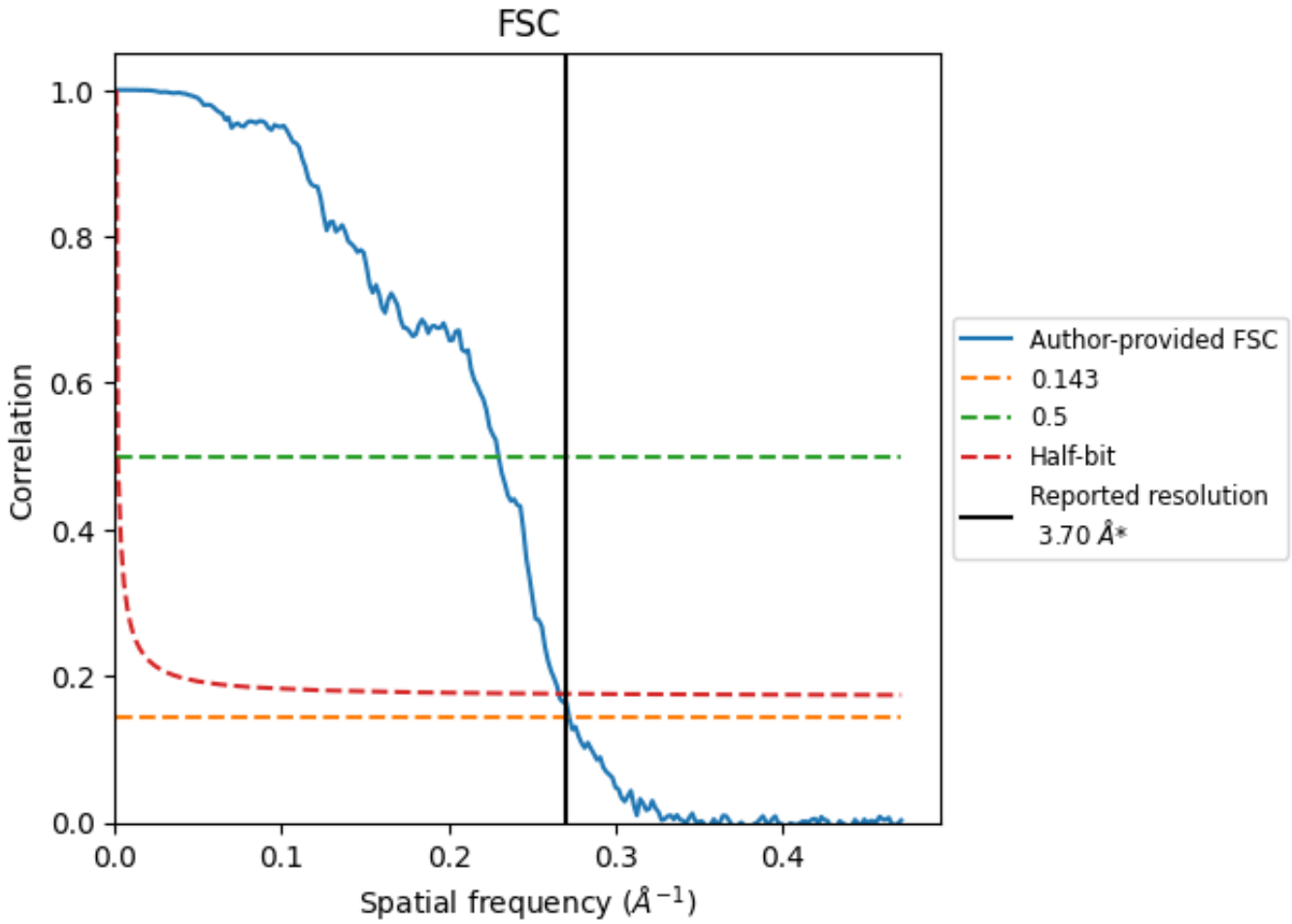


*Reported resolution corresponds to spatial frequency of 0.270 Å^{-1}

8 Fourier-Shell correlation

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



*Reported resolution corresponds to spatial frequency of 0.270 Å^{-1}

8.2 Resolution estimates (i)

Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.70	-	-
Author-provided FSC curve	3.67	4.35	3.76
Unmasked-calculated*	-	-	-

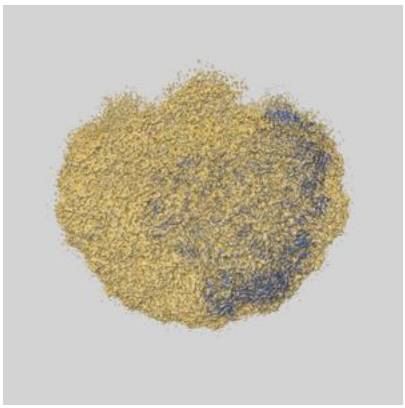
*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps.

9 Map-model fit

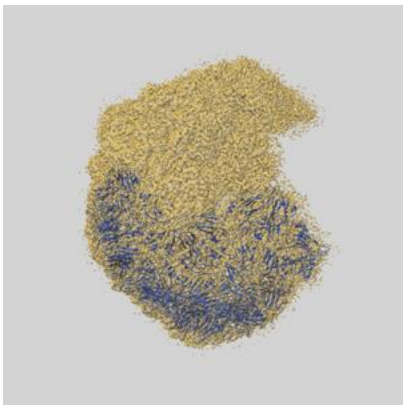
This section contains information regarding the fit between EMDB map EMD-0188 and PDB model 6HBL. Per-residue inclusion information can be found in section 3 on page 8.

9.1 Map-model overlays

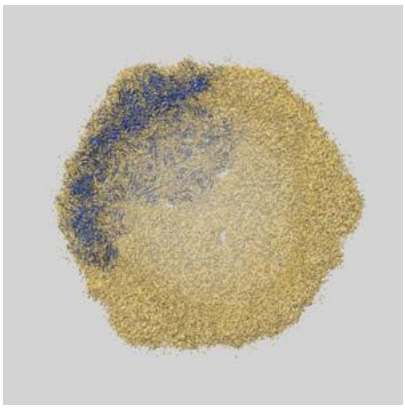
9.1.1 Map-model overlay



X

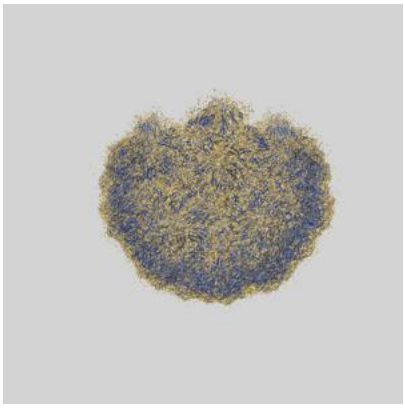


Y

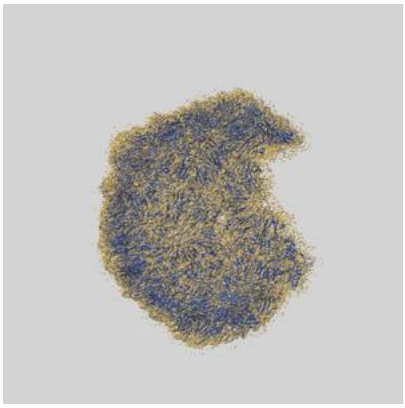


Z

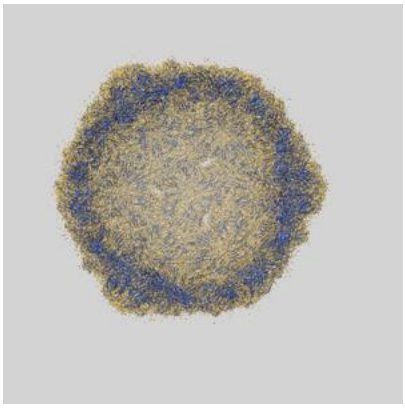
9.1.2 Map-model assembly overlay



X



Y



Z

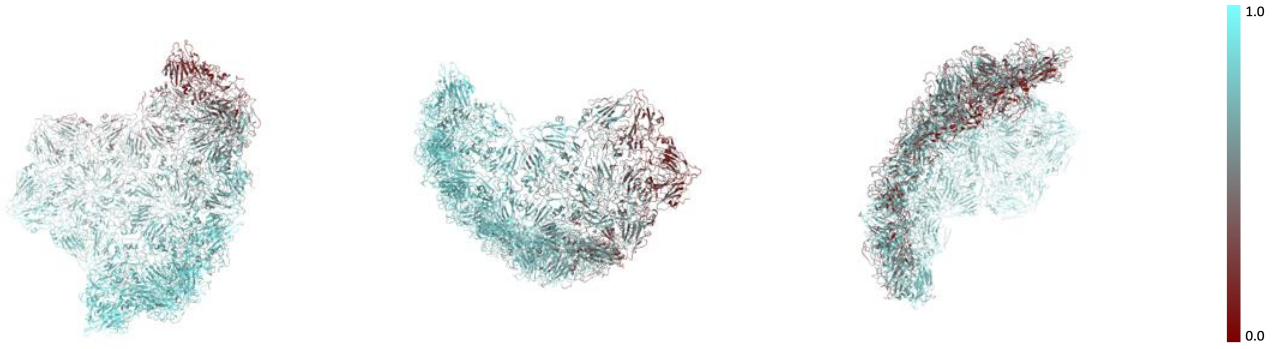
The images above show the 3D surface view of the map at the recommended contour level 0.066 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

9.2 Q-score mapped to coordinate model

X Y Z

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

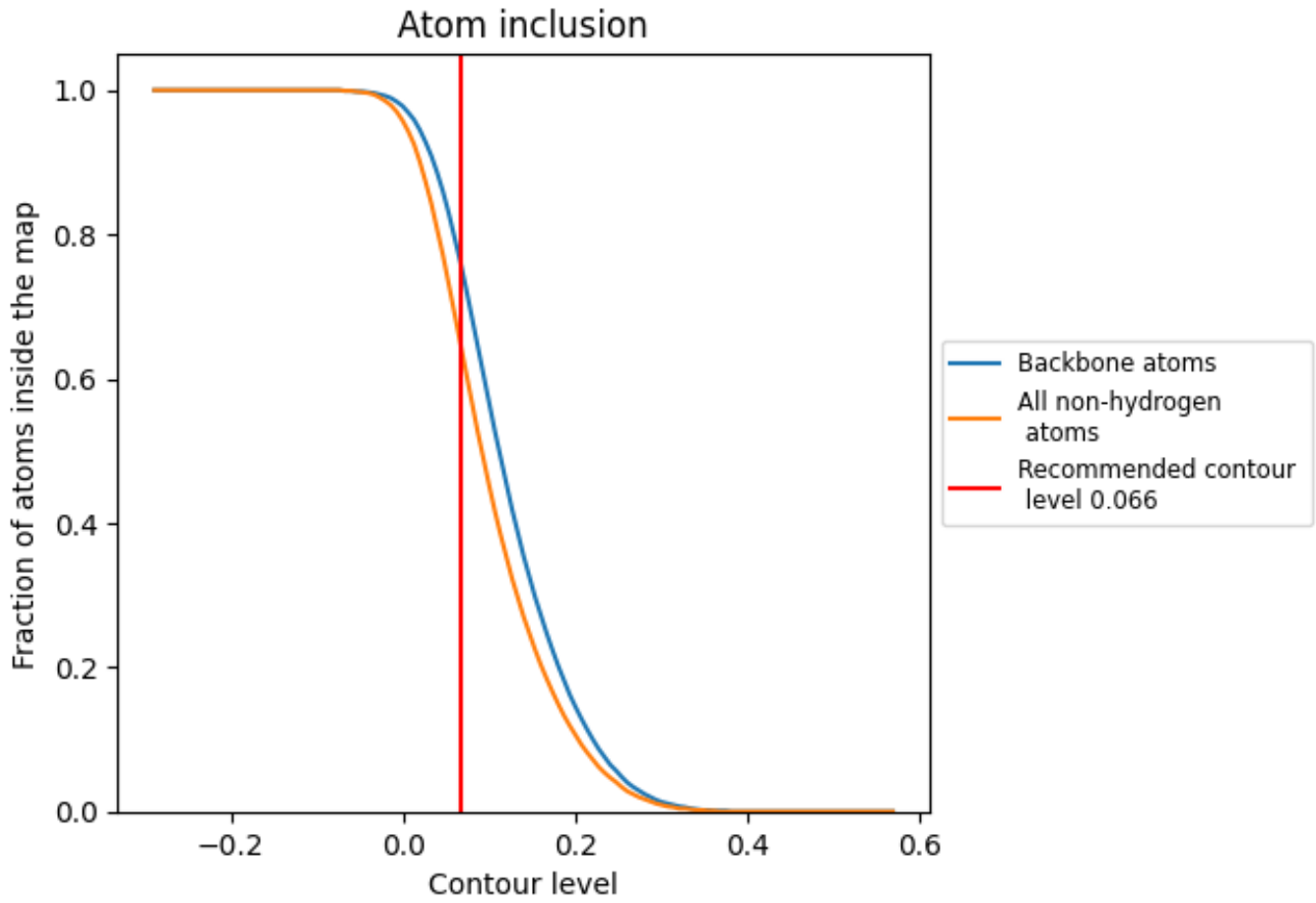
9.3 Atom inclusion mapped to coordinate model



X Y Z

The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.066).

9.4 Atom inclusion (i)



At the recommended contour level, 76% of all backbone atoms, 65% 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 (0.066) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.6510	0.2860
A	0.6550	0.2680
B	0.4690	0.2100
C	0.6000	0.2580
D	0.6390	0.2530
E	0.5700	0.2360
F	0.5210	0.1980
G	0.5930	0.2420
H	0.6470	0.2900
I	0.6550	0.2800
J	0.5110	0.1700
K	0.4800	0.1910
L	0.3940	0.1360
M	0.7180	0.3040
N	0.7030	0.3160
O	0.7430	0.3380
P	0.5780	0.2270
Q	0.6350	0.2820
R	0.6560	0.3010
S	0.7050	0.3020
T	0.7330	0.3280
U	0.7090	0.3020
V	0.4140	0.1330
W	0.2220	0.0920
X	0.2990	0.1000
Y	0.7980	0.3650
Z	0.8120	0.3780
a	0.8080	0.3900
b	0.7880	0.3630
c	0.7870	0.3770
d	0.7840	0.3810
e	0.7970	0.3650
f	0.7870	0.3680
g	0.7980	0.3880
h	0.7690	0.3500

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Chain	Atom inclusion	Q-score
i	0.7520	0.3500
j	0.7540	0.3520
k	0.5230	0.1930
l	0.4360	0.1820
m	0.4890	0.2100
n	0.7810	0.3570
o	0.7600	0.3430
p	0.7800	0.3610
q	0.7340	0.3200
r	0.7660	0.3580
s	0.7520	0.3560