



wwPDB EM Validation Summary Report ⓘ

Mar 8, 2026 – 10:47 AM UTC

PDB ID : 9GB2 / pdb_00009gb2
EMDB ID : EMD-51195
Title : Extended phiCD508 baseplate
Authors : Wilson, J.S.; Fagan, R.P.; Bullough, P.A.
Deposited on : 2024-07-29
Resolution : 3.43 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

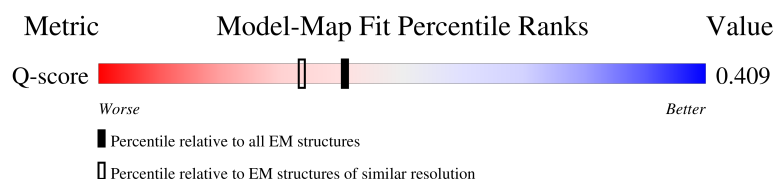
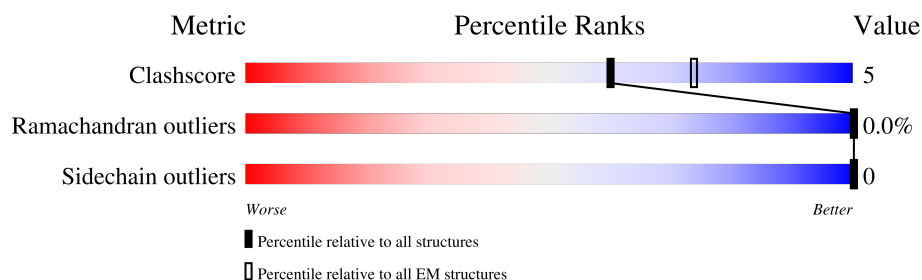
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh &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

1 Overall quality at a glance

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

The reported resolution of this entry is 3.43 Å.

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



Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	13927 (2.93 - 3.93)

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	378	38% 87% 13%
1	B	378	39% 88% 11%
1	C	378	39% 87% 13%
1	D	378	38% 87% 13%

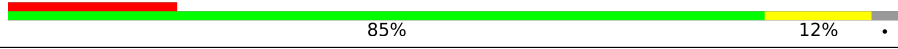
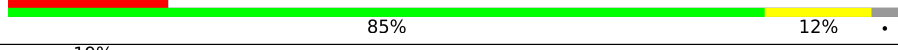
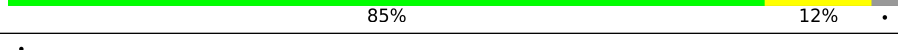
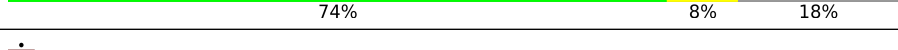
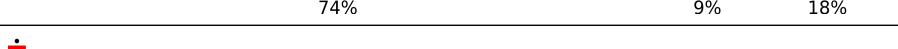
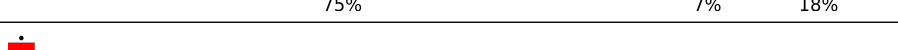
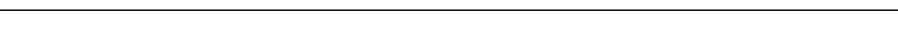
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Mol	Chain	Length	Quality of chain
1	E	378	
1	F	378	
1	G	378	
1	H	378	
1	I	378	
1	J	378	
1	K	378	
1	L	378	
2	M	209	
2	N	209	
2	O	209	
2	P	209	
2	Q	209	
2	R	209	
3	S	223	
3	T	223	
3	U	223	
3	V	223	
3	W	223	
3	X	223	
4	Y	137	
4	Z	137	
4	a	137	
4	b	137	
4	c	137	

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Mol	Chain	Length	Quality of chain
4	d	137	 19% 84% 11% 5%
5	e	473	 19% 86% 11%
5	f	473	 19% 85% 12%
5	g	473	 18% 85% 12%
5	h	473	 19% 83% 14%
5	i	473	 19% 86% 11%
5	j	473	 18% 85% 12%
6	k	152	 19% 74% 8% 18%
6	l	152	 19% 74% 9% 18%
6	m	152	 19% 75% 7% 18%
6	n	152	 19% 74% 8% 18%
6	o	152	 19% 74% 9% 18%
6	p	152	 19% 75% 7% 18%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 173292 atoms, of which 86868 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 gp65 - Triplex 1a protein.

Mol	Chain	Residues	Atoms						AltConf	Trace
1	A	377	Total	C	H	N	O	S	0	0
			5809	1836	2899	473	597	4		
1	B	377	Total	C	H	N	O	S	0	0
			5809	1836	2899	473	597	4		
1	C	377	Total	C	H	N	O	S	0	0
			5809	1836	2899	473	597	4		
1	D	377	Total	C	H	N	O	S	0	0
			5809	1836	2899	473	597	4		
1	E	377	Total	C	H	N	O	S	0	0
			5809	1836	2899	473	597	4		
1	F	377	Total	C	H	N	O	S	0	0
			5809	1836	2899	473	597	4		
1	G	374	Total	C	H	N	O	S	0	0
			5777	1824	2885	470	594	4		
1	H	374	Total	C	H	N	O	S	0	0
			5777	1824	2885	470	594	4		
1	I	374	Total	C	H	N	O	S	0	0
			5777	1824	2885	470	594	4		
1	J	374	Total	C	H	N	O	S	0	0
			5777	1824	2885	470	594	4		
1	K	374	Total	C	H	N	O	S	0	0
			5777	1824	2885	470	594	4		
1	L	374	Total	C	H	N	O	S	0	0
			5777	1824	2885	470	594	4		

- Molecule 2 is a protein called gp66 - Triplex 2 protein.

Mol	Chain	Residues	Atoms						AltConf	Trace
2	M	157	Total	C	H	N	O	S	0	0
			2559	811	1295	205	241	7		
2	N	157	Total	C	H	N	O	S	0	0
			2559	811	1295	205	241	7		
2	O	157	Total	C	H	N	O	S	0	0
			2559	811	1295	205	241	7		

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Mol	Chain	Residues	Atoms						AltConf	Trace
2	P	157	Total	C	H	N	O	S	0	0
			2559	811	1295	205	241	7		
2	Q	157	Total	C	H	N	O	S	0	0
			2559	811	1295	205	241	7		
2	R	157	Total	C	H	N	O	S	0	0
			2559	811	1295	205	241	7		

- Molecule 3 is a protein called gp61 - Tail tube initiator.

Mol	Chain	Residues	Atoms						AltConf	Trace
3	S	206	Total	C	H	N	O	S	0	0
			3411	1076	1736	283	315	1		
3	T	206	Total	C	H	N	O	S	0	0
			3411	1076	1736	283	315	1		
3	U	206	Total	C	H	N	O	S	0	0
			3411	1076	1736	283	315	1		
3	V	206	Total	C	H	N	O	S	0	0
			3411	1076	1736	283	315	1		
3	W	206	Total	C	H	N	O	S	0	0
			3411	1076	1736	283	315	1		
3	X	206	Total	C	H	N	O	S	0	0
			3411	1076	1736	283	315	1		

- Molecule 4 is a protein called gp56 - Tail tube protein.

Mol	Chain	Residues	Atoms						AltConf	Trace
4	Y	130	Total	C	H	N	O	S	0	0
			2056	652	1027	167	207	3		
4	Z	130	Total	C	H	N	O	S	0	0
			2056	652	1027	167	207	3		
4	a	130	Total	C	H	N	O	S	0	0
			2056	652	1027	167	207	3		
4	b	130	Total	C	H	N	O	S	0	0
			2056	652	1027	167	207	3		
4	c	130	Total	C	H	N	O	S	0	0
			2056	652	1027	167	207	3		
4	d	130	Total	C	H	N	O	S	0	0
			2056	652	1027	167	207	3		

- Molecule 5 is a protein called gp55 - Tail sheath protein.

Mol	Chain	Residues	Atoms						AltConf	Trace
5	e	460	Total	C	H	N	O	S	0	0
			7233	2299	3618	585	721	10		
5	f	460	Total	C	H	N	O	S	0	0
			7233	2299	3618	585	721	10		
5	g	460	Total	C	H	N	O	S	0	0
			7233	2299	3618	585	721	10		
5	h	460	Total	C	H	N	O	S	0	0
			7233	2299	3618	585	721	10		
5	i	460	Total	C	H	N	O	S	0	0
			7233	2299	3618	585	721	10		
5	j	460	Total	C	H	N	O	S	0	0
			7233	2299	3618	585	721	10		

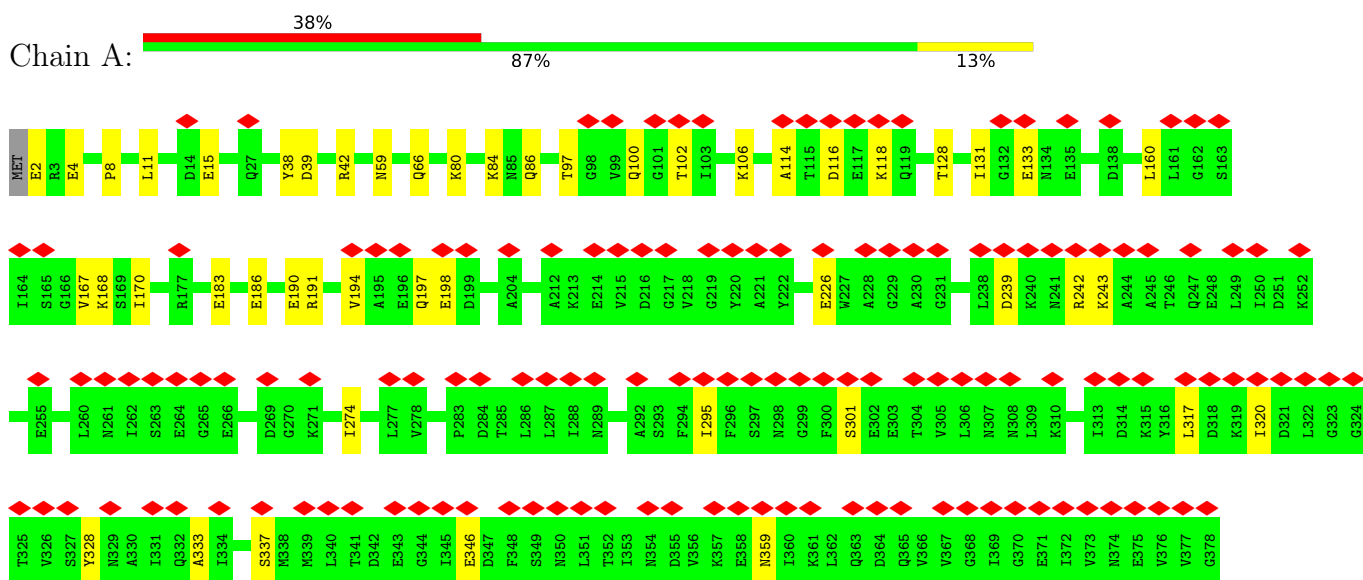
- Molecule 6 is a protein called gp64 - Sheath initiator.

Mol	Chain	Residues	Atoms						AltConf	Trace
6	k	125	Total	C	H	N	O	S	0	0
			2037	658	1018	162	197	2		
6	l	125	Total	C	H	N	O	S	0	0
			2037	658	1018	162	197	2		
6	m	125	Total	C	H	N	O	S	0	0
			2037	658	1018	162	197	2		
6	n	125	Total	C	H	N	O	S	0	0
			2037	658	1018	162	197	2		
6	o	125	Total	C	H	N	O	S	0	0
			2037	658	1018	162	197	2		
6	p	125	Total	C	H	N	O	S	0	0
			2037	658	1018	162	197	2		

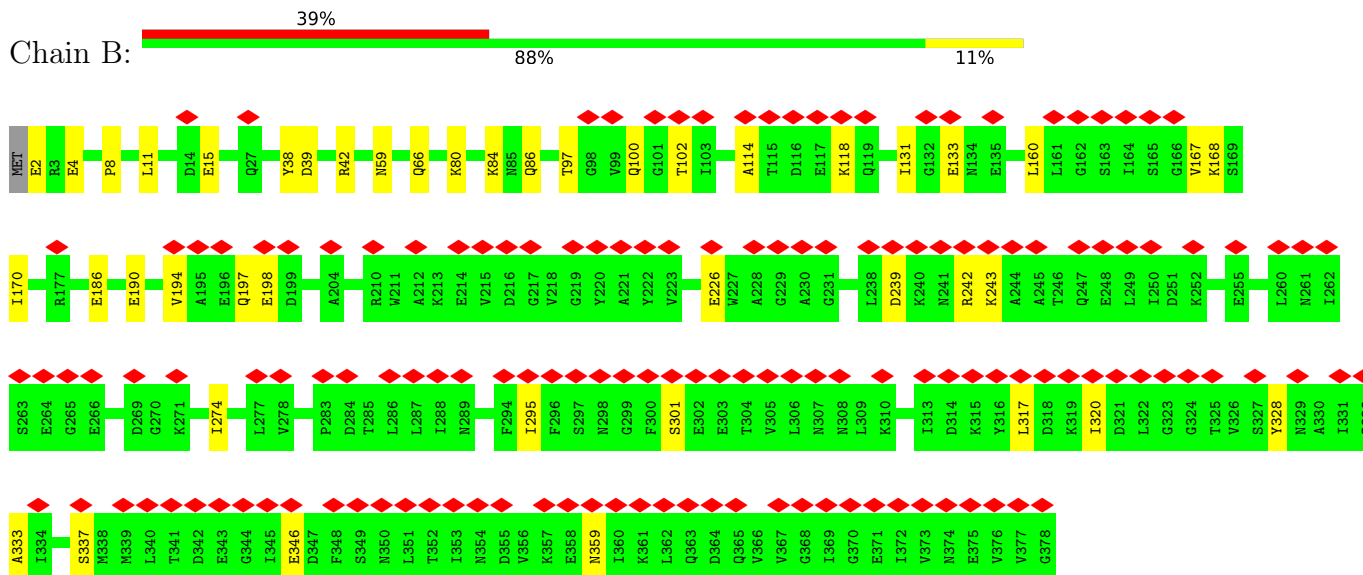
3 Residue-property plots [i](#)

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.

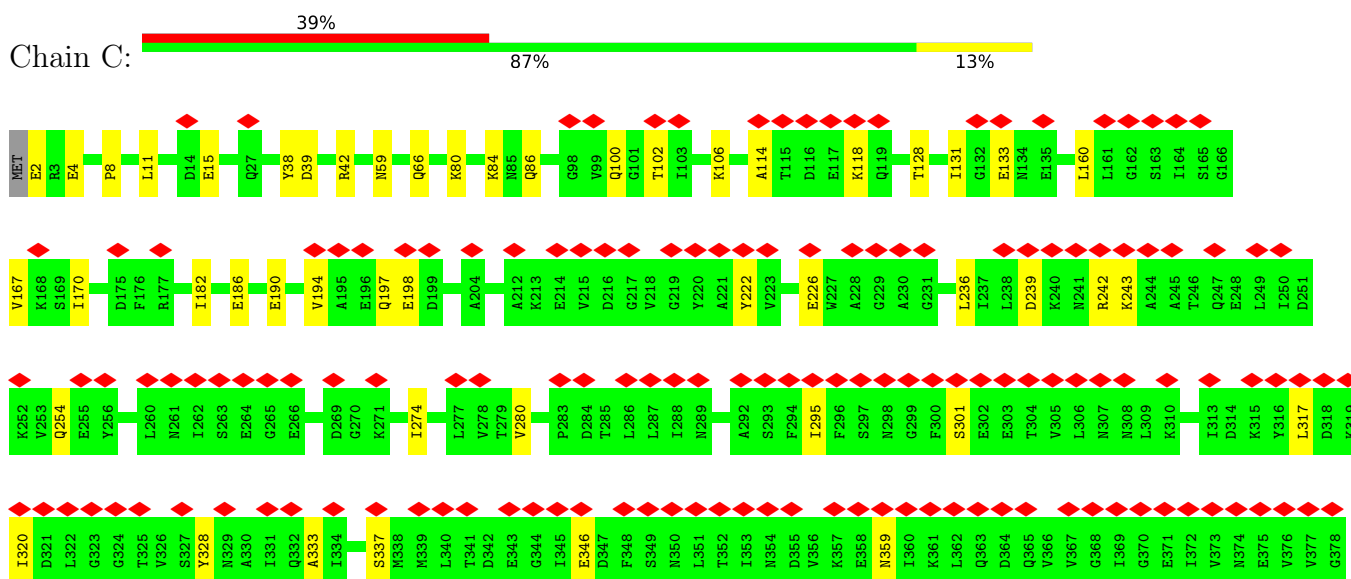
- Molecule 1: gp65 - Triplex 1a protein



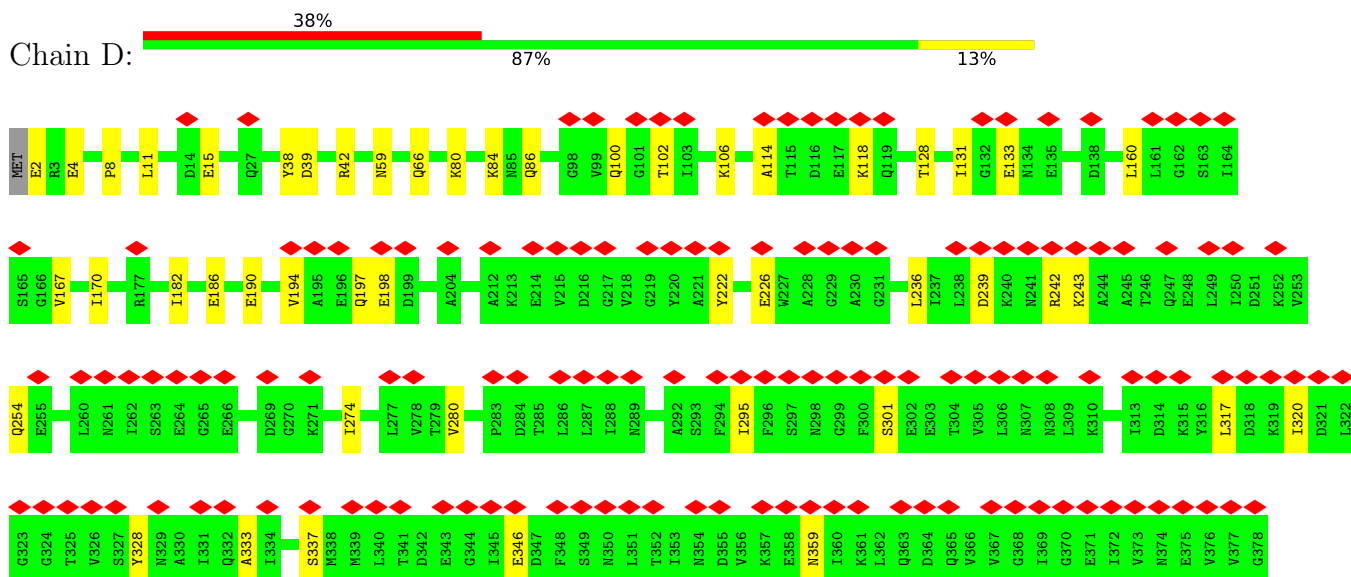
- Molecule 1: gp65 - Triplex 1a protein



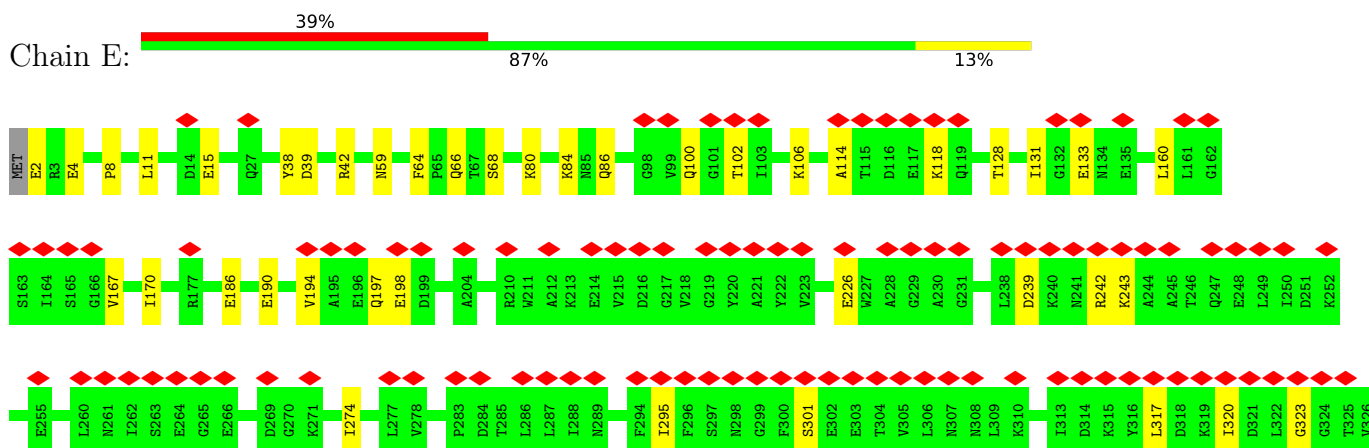
- Molecule 1: gp65 - Triplex 1a protein

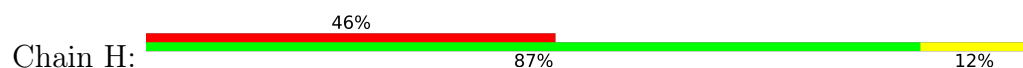


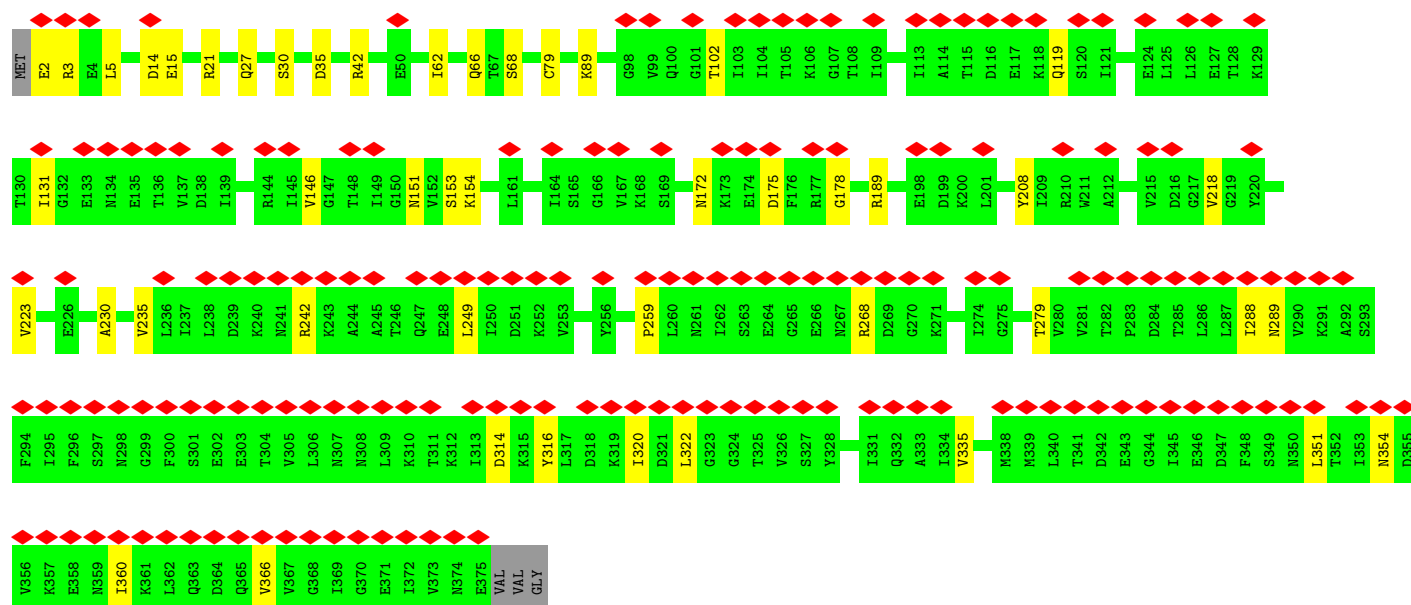
• Molecule 1: gp65 - Triplex 1a protein



• Molecule 1: gp65 - Triplex 1a protein

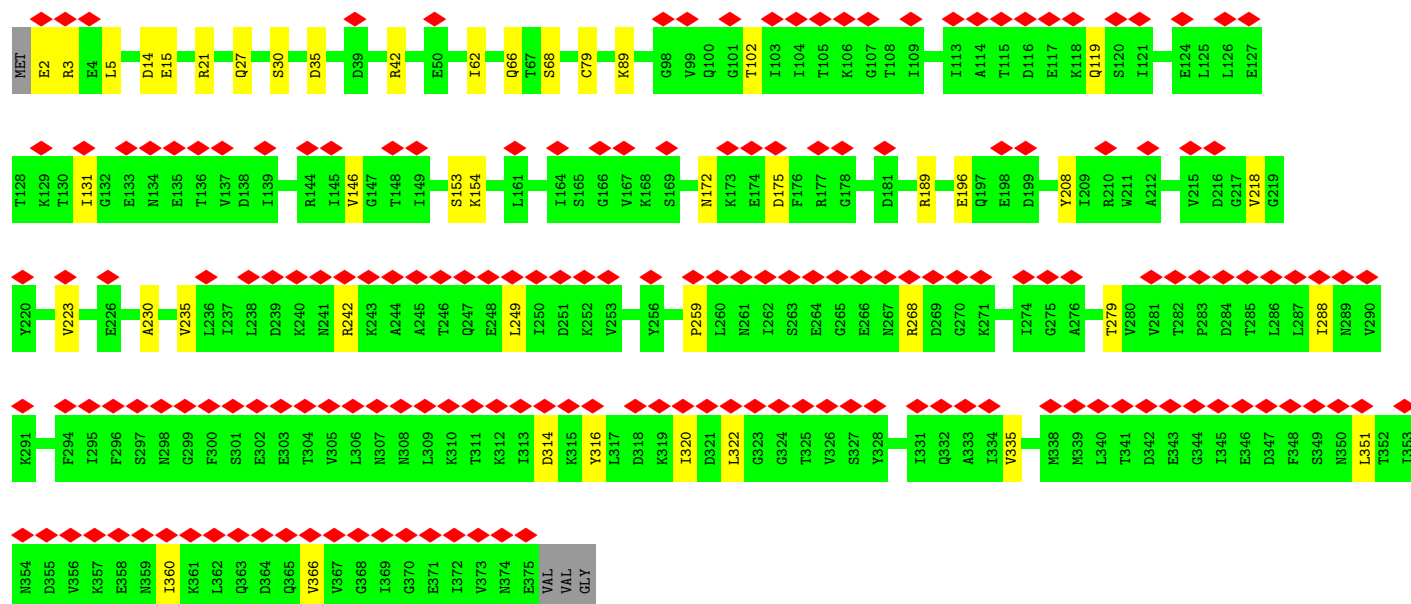






- Molecule 1: gp65 - Triplex 1a protein

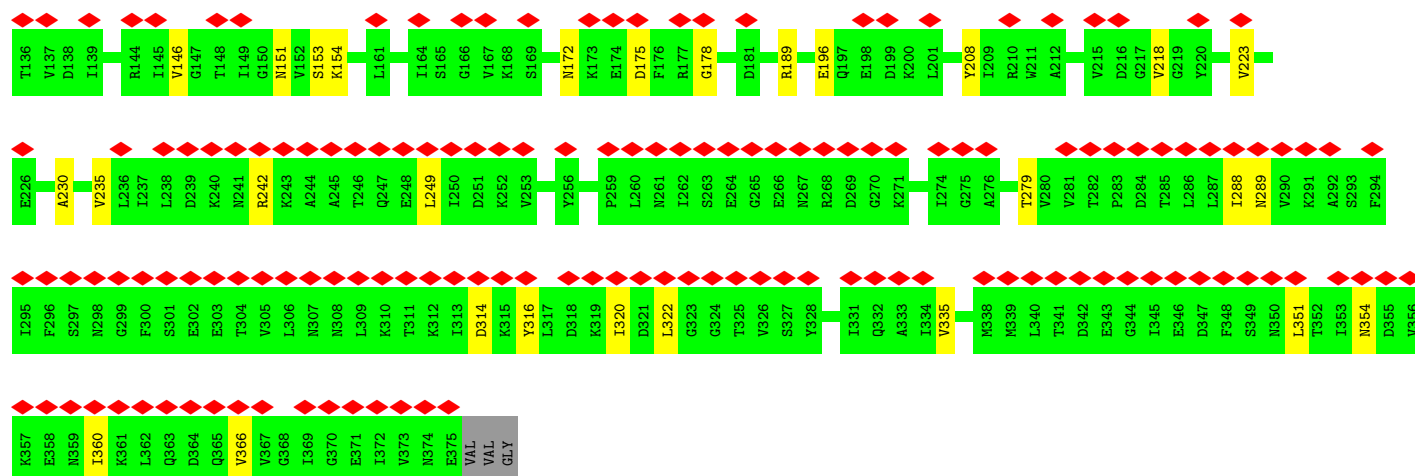
Chain I: 47% 87% 12% .



- Molecule 1: gp65 - Triplex 1a protein

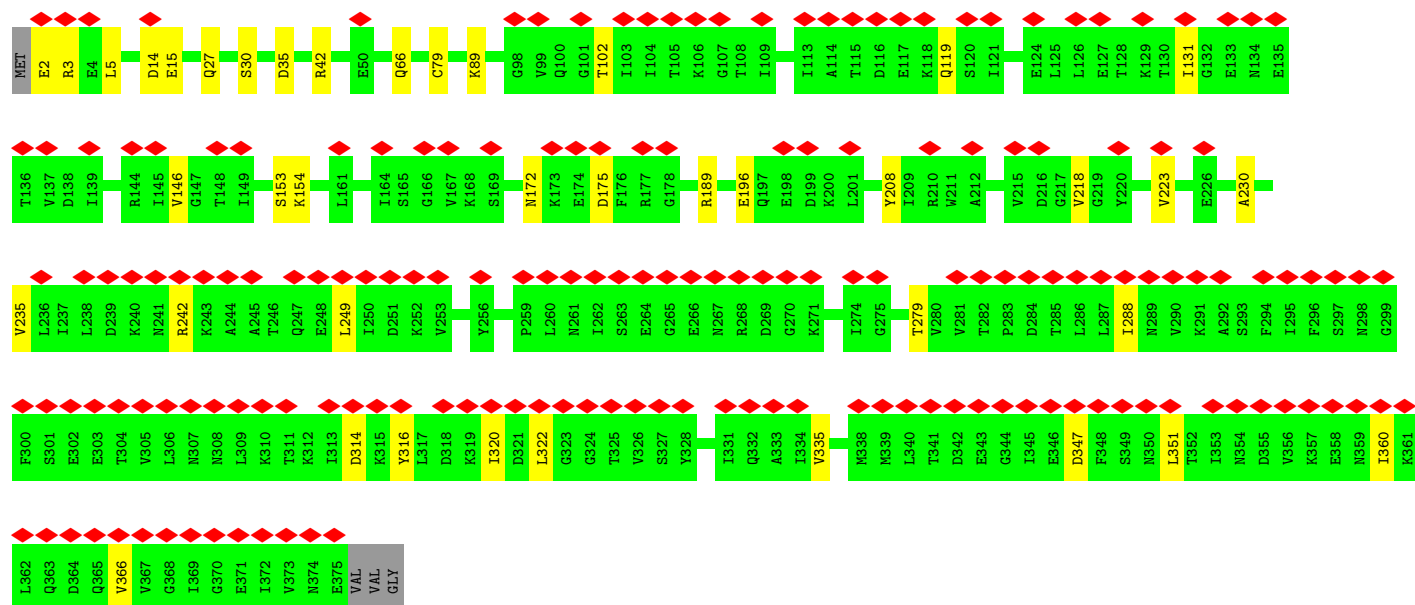
Chain J: 47% 88% 11% .





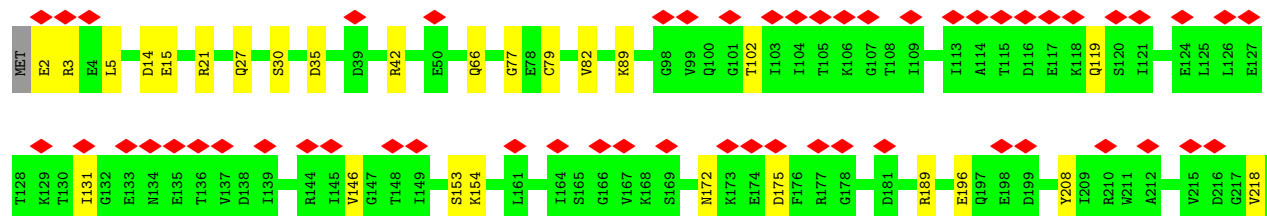
- Molecule 1: gp65 - Triplex 1a protein

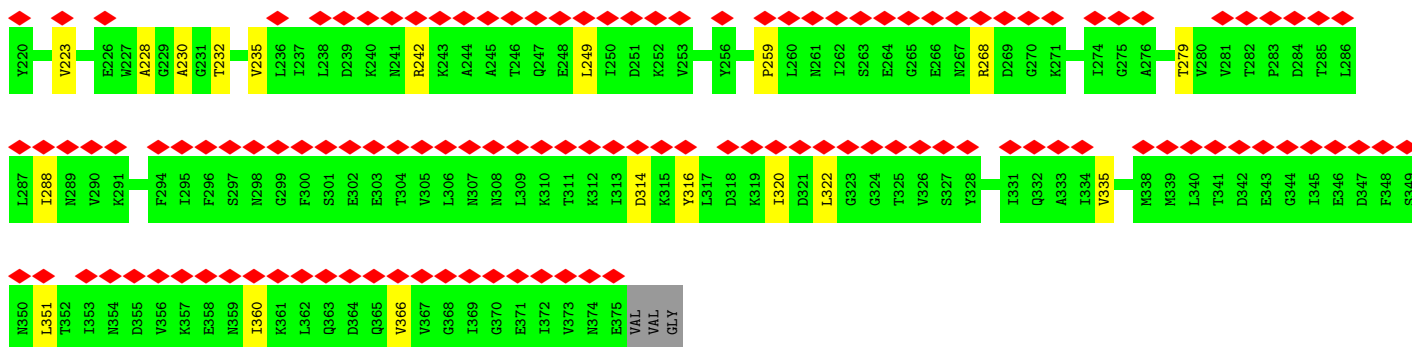
Chain K: 46% 88% 11% .



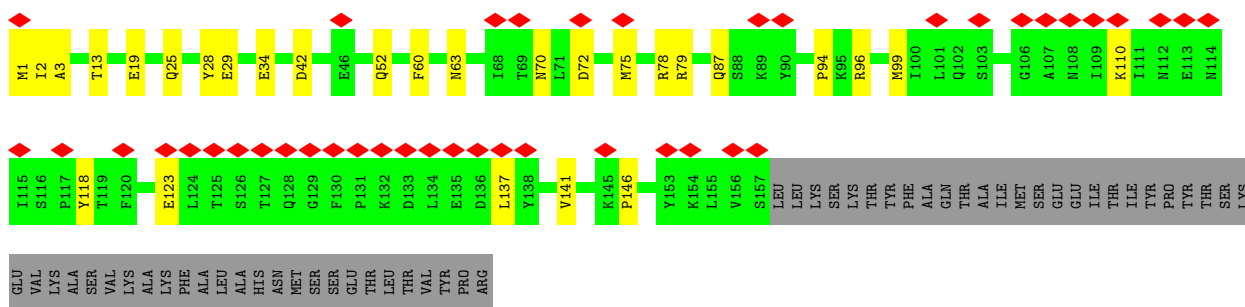
- Molecule 1: gp65 - Triplex 1a protein

Chain L: 47% 87% 12% .

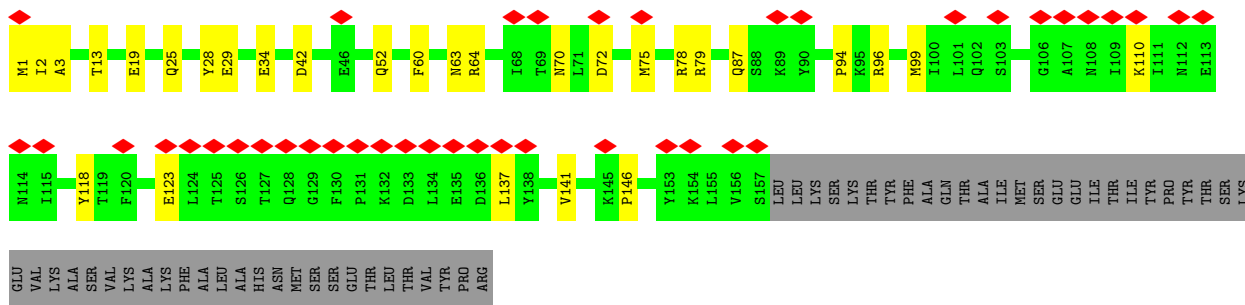




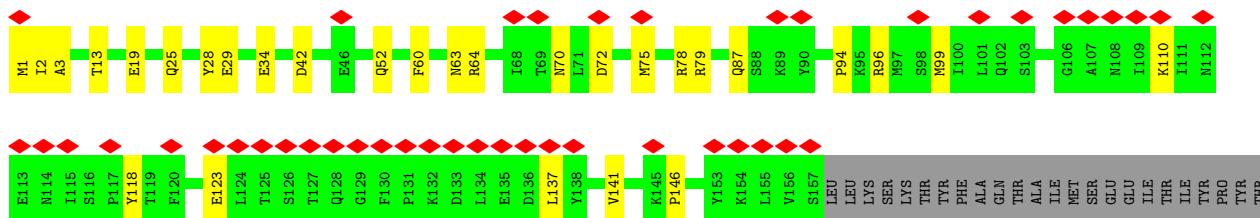
- Molecule 2: gp66 - Triplex 2 protein



- Molecule 2: gp66 - Triplex 2 protein



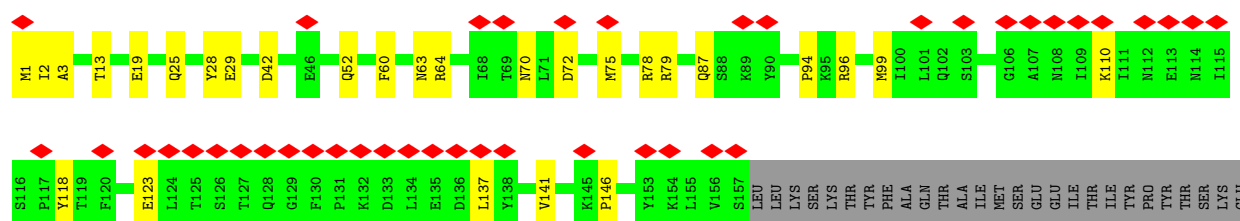
- Molecule 2: gp66 - Triplex 2 protein



SER LYS GLU VAL LYS ALA SER VAL LYS ALA LYS PHE LEU ALA HIS ASN MET SER SER SER THR THR VAL TYR PRO ARG

• Molecule 2: gp66 - Triplex 2 protein

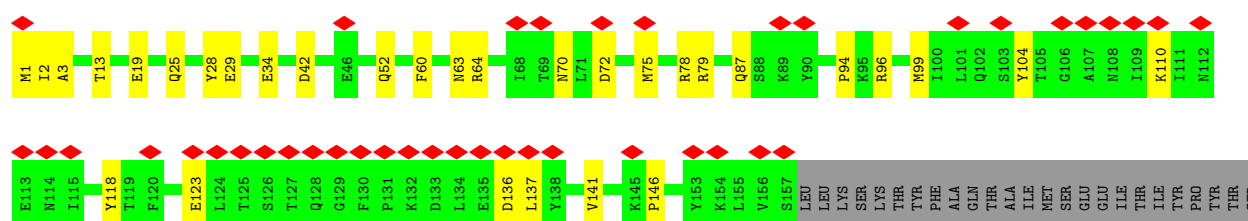
Chain P: 20% 62% 13% 25%



VAL LYS ALA VAL LYS ALA LYS PHE LEU ALA HIS ASN MET SER SER SER THR THR VAL TYR PRO ARG

• Molecule 2: gp66 - Triplex 2 protein

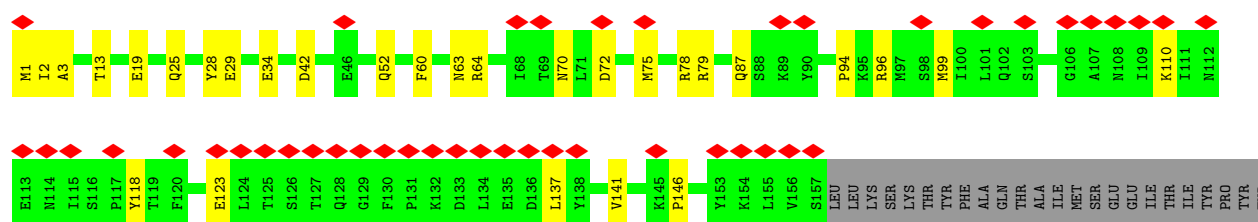
Chain Q: 20% 60% 15% 25%



LYS GLU VAL LYS ALA SER VAL LYS ALA LYS PHE LEU ALA HIS ASN MET SER SER SER THR THR VAL TYR PRO ARG

• Molecule 2: gp66 - Triplex 2 protein

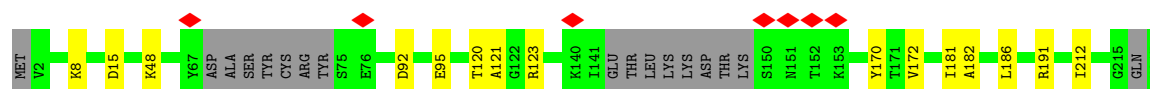
Chain R: 21% 61% 14% 25%



SER LYS GLU VAL LYS ALA SER VAL LYS ALA LYS PHE LEU ALA HIS ASN MET SER SER SER THR THR VAL TYR PRO ARG

• Molecule 3: gp61 - Tail tube initiator

Chain S: 86% 7% 8%





- Molecule 3: gp61 - Tail tube initiator

Chain T: 87% 6% 8%



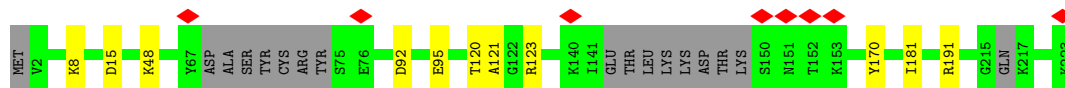
- Molecule 3: gp61 - Tail tube initiator

Chain U: 87% 5% 8%



- Molecule 3: gp61 - Tail tube initiator

Chain V: 87% 5% 8%



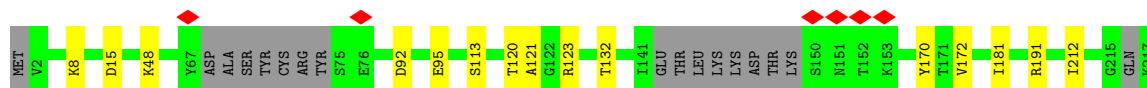
- Molecule 3: gp61 - Tail tube initiator

Chain W: 86% 7% 8%



- Molecule 3: gp61 - Tail tube initiator

Chain X: 86% 7% 8%



- Molecule 4: gp56 - Tail tube protein

Chain Y:  83% 12% 5%



- Molecule 4: gp56 - Tail tube protein

Chain Z:  86% 9% 5%



- Molecule 4: gp56 - Tail tube protein

Chain a:  84% 11% 5%



- Molecule 4: gp56 - Tail tube protein

Chain b:  85% 9% 5%



- Molecule 4: gp56 - Tail tube protein

Chain c:  87% 8% 5%



- Molecule 4: gp56 - Tail tube protein

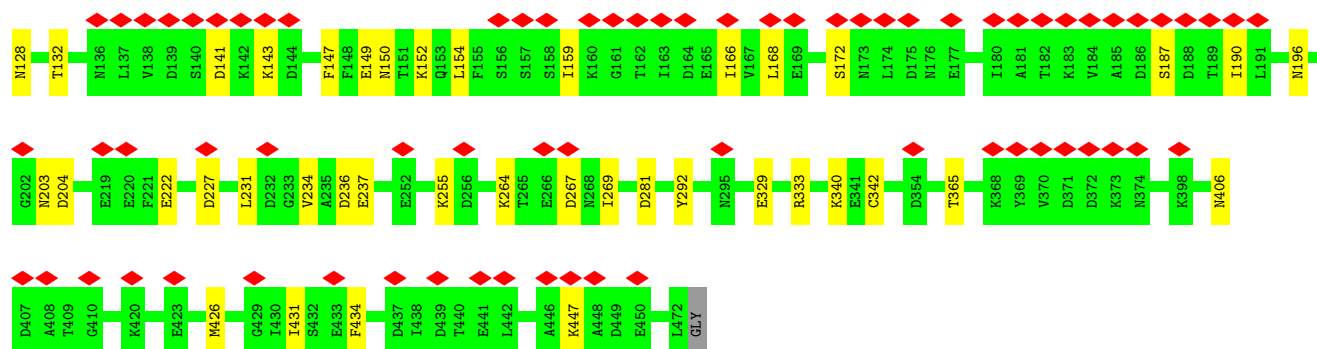
Chain d:  84% 11% 5%



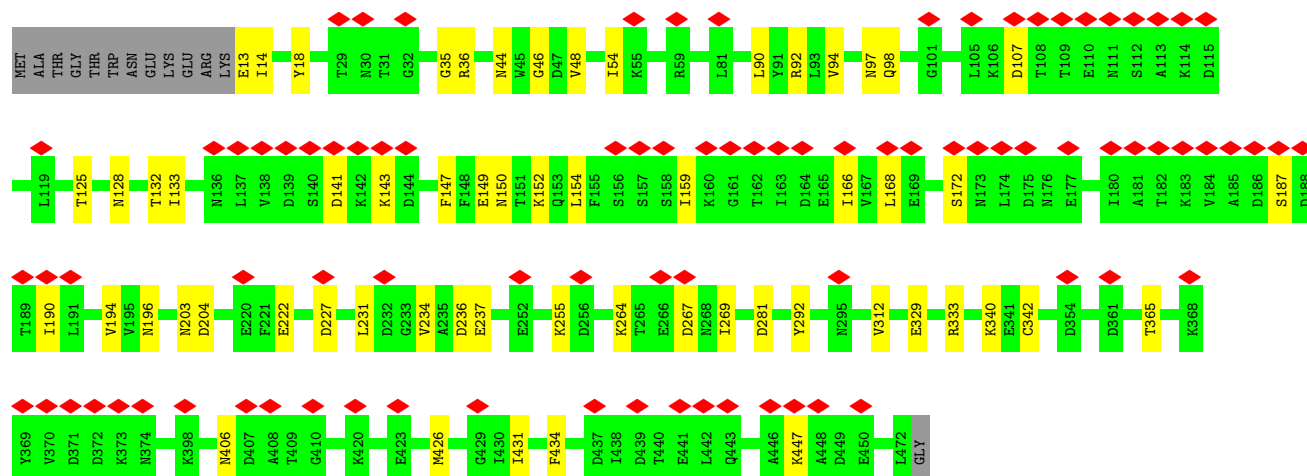
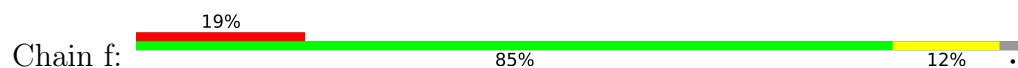
- Molecule 5: gp55 - Tail sheath protein

Chain e:  19% 86% 11% 5%

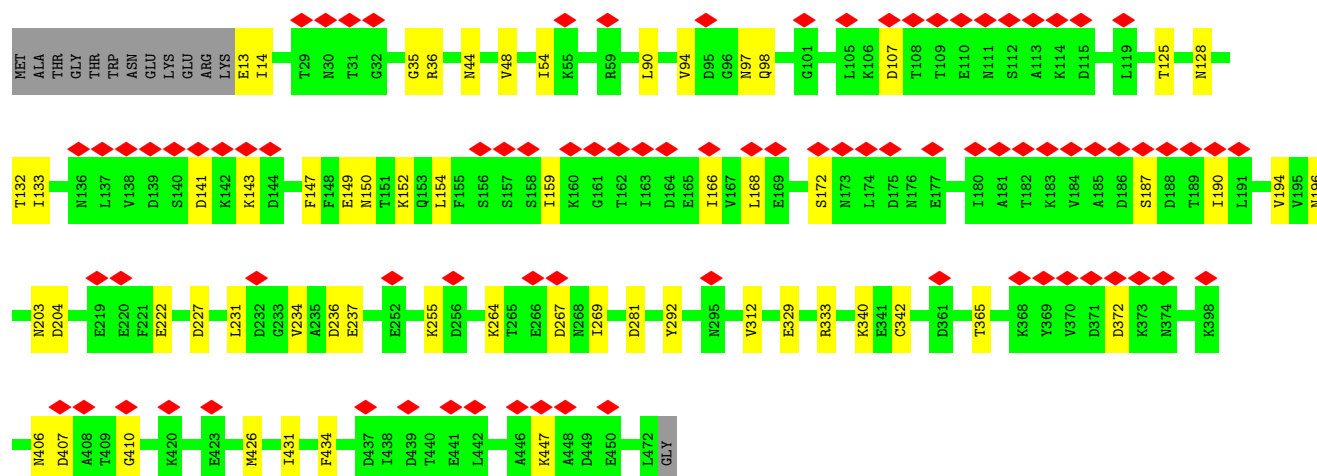
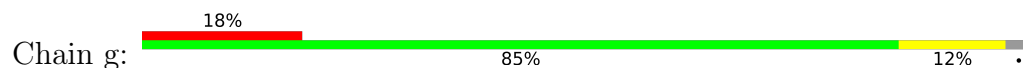




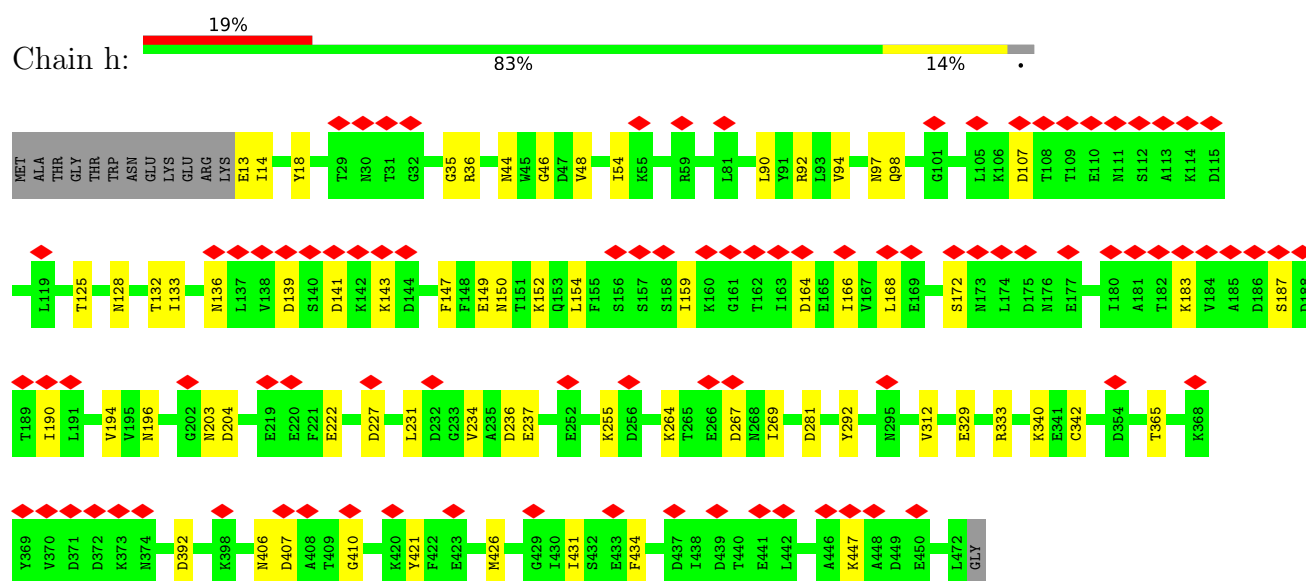
• Molecule 5: gp55 - Tail sheath protein



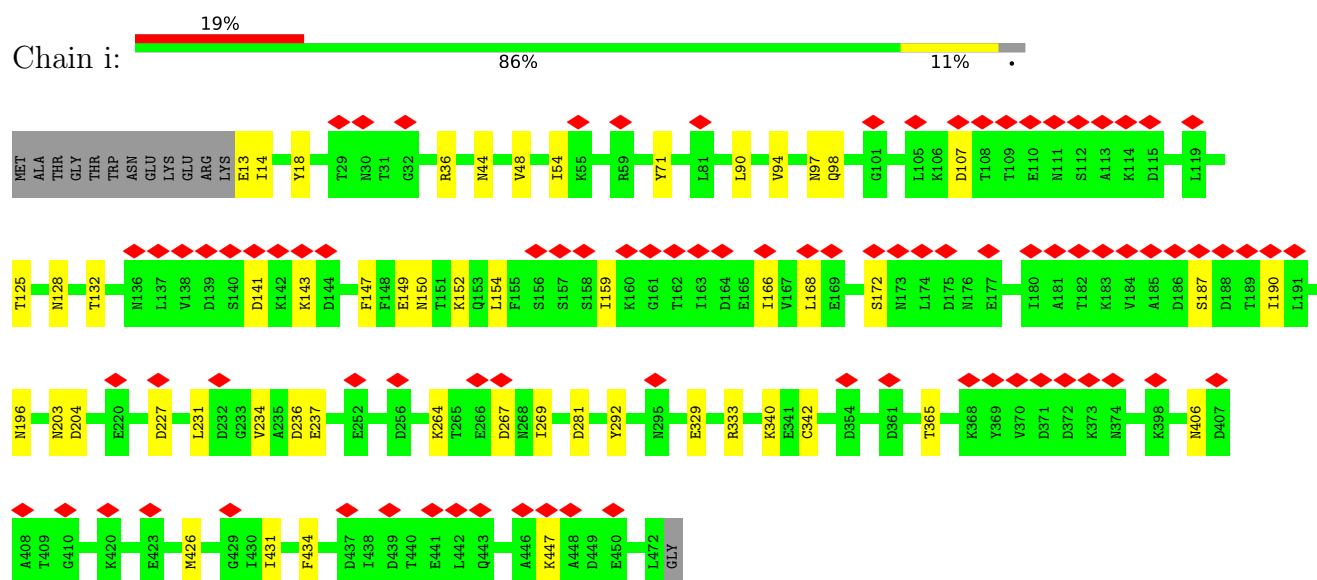
• Molecule 5: gp55 - Tail sheath protein



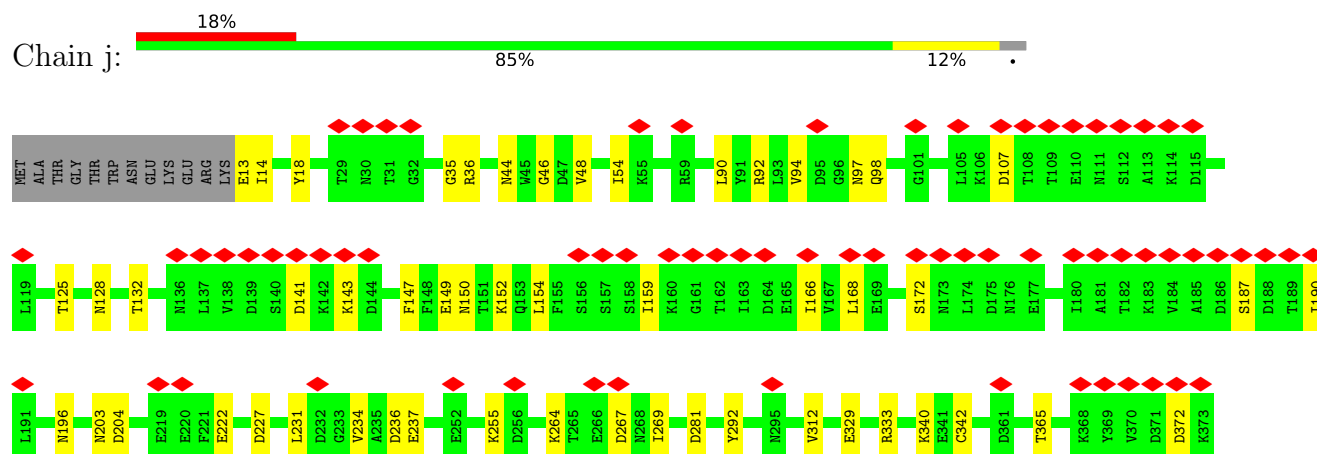
• Molecule 5: gp55 - Tail sheath protein



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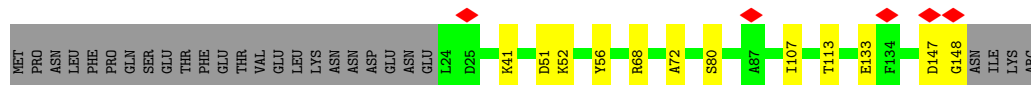


• Molecule 5: gp55 - Tail sheath protein





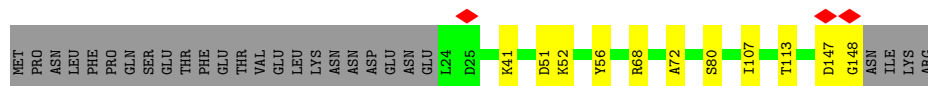
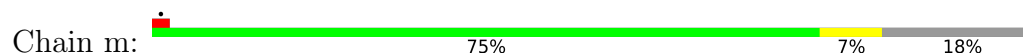
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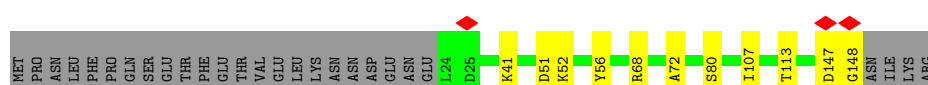
- Molecule 6: gp64 - Sheath initiator



- Molecule 6: gp64 - Sheath initiator



- Molecule 6: gp64 - Sheath initiator



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	19276	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	42	Depositor
Minimum defocus (nm)	1500	Depositor
Maximum defocus (nm)	3000	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	2.189	Depositor
Minimum map value	-1.057	Depositor
Average map value	0.006	Depositor
Map value standard deviation	0.089	Depositor
Recommended contour level	0.469	Depositor
Map size (Å)	530.0, 530.0, 530.0	wwPDB
Map dimensions	500, 500, 500	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.06, 1.06, 1.06	Depositor

5 Model quality

5.1 Standard geometry

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	0.14	0/2948	0.29	0/3993
1	B	0.15	0/2948	0.29	0/3993
1	C	0.15	0/2948	0.29	0/3993
1	D	0.14	0/2948	0.29	0/3993
1	E	0.14	0/2948	0.29	0/3993
1	F	0.14	0/2948	0.29	0/3993
1	G	0.14	0/2930	0.29	0/3968
1	H	0.14	0/2930	0.29	0/3968
1	I	0.14	0/2930	0.29	0/3968
1	J	0.14	0/2930	0.29	0/3968
1	K	0.14	0/2930	0.29	0/3968
1	L	0.14	0/2930	0.29	0/3968
2	M	0.15	0/1286	0.31	0/1734
2	N	0.15	0/1286	0.30	0/1734
2	O	0.15	0/1286	0.30	0/1734
2	P	0.15	0/1286	0.31	0/1734
2	Q	0.15	0/1286	0.30	0/1734
2	R	0.15	0/1286	0.30	0/1734
3	S	0.20	0/1702	0.34	0/2288
3	T	0.20	0/1702	0.34	0/2288
3	U	0.20	0/1702	0.34	0/2288
3	V	0.20	0/1702	0.34	0/2288
3	W	0.20	0/1702	0.34	0/2288
3	X	0.20	0/1702	0.34	0/2288
4	Y	0.18	0/1041	0.31	0/1394
4	Z	0.18	0/1041	0.32	0/1394
4	a	0.18	0/1041	0.31	0/1394
4	b	0.18	0/1041	0.31	0/1394
4	c	0.18	0/1041	0.32	0/1394
4	d	0.18	0/1041	0.30	0/1394
5	e	0.15	0/3667	0.31	0/4946
5	f	0.15	0/3667	0.31	0/4946
5	g	0.15	0/3667	0.31	0/4946
5	h	0.15	0/3667	0.31	0/4946

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
5	i	0.15	0/3667	0.31	0/4946
5	j	0.15	0/3667	0.31	0/4946
6	k	0.19	0/1038	0.32	0/1399
6	l	0.19	0/1038	0.32	0/1399
6	m	0.19	0/1038	0.32	0/1399
6	n	0.19	0/1038	0.32	0/1399
6	o	0.19	0/1038	0.32	0/1399
6	p	0.19	0/1038	0.32	0/1399
All	All	0.16	0/87672	0.31	0/118332

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

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	2910	2899	2909	33	0
1	B	2910	2899	2909	30	0
1	C	2910	2899	2909	33	0
1	D	2910	2899	2909	33	0
1	E	2910	2899	2909	33	0
1	F	2910	2899	2909	34	0
1	G	2892	2885	2888	32	0
1	H	2892	2885	2888	31	0
1	I	2892	2885	2888	30	0
1	J	2892	2885	2888	29	0
1	K	2892	2885	2888	28	0
1	L	2892	2885	2888	31	0
2	M	1264	1295	1295	22	0
2	N	1264	1295	1295	23	0
2	O	1264	1295	1295	23	0
2	P	1264	1295	1295	22	0
2	Q	1264	1295	1295	24	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	R	1264	1295	1295	23	0
3	S	1675	1736	1728	11	0
3	T	1675	1736	1728	10	0
3	U	1675	1736	1728	9	0
3	V	1675	1736	1728	9	0
3	W	1675	1736	1728	11	0
3	X	1675	1736	1728	11	0
4	Y	1029	1027	1030	10	0
4	Z	1029	1027	1030	8	0
4	a	1029	1027	1030	9	0
4	b	1029	1027	1030	8	0
4	c	1029	1027	1030	7	0
4	d	1029	1027	1030	9	0
5	e	3615	3618	3612	38	0
5	f	3615	3618	3612	40	0
5	g	3615	3618	3612	39	0
5	h	3615	3618	3612	44	0
5	i	3615	3618	3612	37	0
5	j	3615	3618	3612	39	0
6	k	1019	1018	1017	9	0
6	l	1019	1018	1017	9	0
6	m	1019	1018	1017	8	0
6	n	1019	1018	1017	9	0
6	o	1019	1018	1017	9	0
6	p	1019	1018	1017	8	0
All	All	86424	86868	86874	782	0

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

The worst 5 of 782 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:E:242:ARG:NH2	1:E:317:LEU:O	2.20	0.75
1:D:242:ARG:NH2	1:D:317:LEU:O	2.20	0.75
1:A:242:ARG:NH2	1:A:317:LEU:O	2.20	0.75
1:F:242:ARG:NH2	1:F:317:LEU:O	2.20	0.75
1:C:242:ARG:NH2	1:C:317:LEU:O	2.20	0.75

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	375/378 (99%)	358 (96%)	17 (4%)	0	100	100
1	B	375/378 (99%)	358 (96%)	17 (4%)	0	100	100
1	C	375/378 (99%)	358 (96%)	17 (4%)	0	100	100
1	D	375/378 (99%)	358 (96%)	17 (4%)	0	100	100
1	E	375/378 (99%)	358 (96%)	17 (4%)	0	100	100
1	F	375/378 (99%)	358 (96%)	17 (4%)	0	100	100
1	G	372/378 (98%)	358 (96%)	14 (4%)	0	100	100
1	H	372/378 (98%)	357 (96%)	15 (4%)	0	100	100
1	I	372/378 (98%)	357 (96%)	15 (4%)	0	100	100
1	J	372/378 (98%)	358 (96%)	14 (4%)	0	100	100
1	K	372/378 (98%)	357 (96%)	15 (4%)	0	100	100
1	L	372/378 (98%)	357 (96%)	15 (4%)	0	100	100
2	M	155/209 (74%)	148 (96%)	7 (4%)	0	100	100
2	N	155/209 (74%)	149 (96%)	6 (4%)	0	100	100
2	O	155/209 (74%)	149 (96%)	6 (4%)	0	100	100
2	P	155/209 (74%)	148 (96%)	7 (4%)	0	100	100
2	Q	155/209 (74%)	149 (96%)	6 (4%)	0	100	100
2	R	155/209 (74%)	149 (96%)	6 (4%)	0	100	100
3	S	198/223 (89%)	191 (96%)	7 (4%)	0	100	100
3	T	198/223 (89%)	191 (96%)	7 (4%)	0	100	100
3	U	198/223 (89%)	191 (96%)	7 (4%)	0	100	100
3	V	198/223 (89%)	191 (96%)	7 (4%)	0	100	100
3	W	198/223 (89%)	191 (96%)	7 (4%)	0	100	100
3	X	198/223 (89%)	191 (96%)	7 (4%)	0	100	100
4	Y	128/137 (93%)	121 (94%)	7 (6%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
4	Z	128/137 (93%)	120 (94%)	8 (6%)	0	100	100
4	a	128/137 (93%)	119 (93%)	9 (7%)	0	100	100
4	b	128/137 (93%)	120 (94%)	8 (6%)	0	100	100
4	c	128/137 (93%)	120 (94%)	8 (6%)	0	100	100
4	d	128/137 (93%)	119 (93%)	9 (7%)	0	100	100
5	e	458/473 (97%)	435 (95%)	23 (5%)	0	100	100
5	f	458/473 (97%)	436 (95%)	22 (5%)	0	100	100
5	g	458/473 (97%)	435 (95%)	22 (5%)	1 (0%)	43	73
5	h	458/473 (97%)	435 (95%)	23 (5%)	0	100	100
5	i	458/473 (97%)	436 (95%)	22 (5%)	0	100	100
5	j	458/473 (97%)	435 (95%)	22 (5%)	1 (0%)	43	73
6	k	123/152 (81%)	118 (96%)	5 (4%)	0	100	100
6	l	123/152 (81%)	119 (97%)	4 (3%)	0	100	100
6	m	123/152 (81%)	119 (97%)	4 (3%)	0	100	100
6	n	123/152 (81%)	118 (96%)	5 (4%)	0	100	100
6	o	123/152 (81%)	119 (97%)	4 (3%)	0	100	100
6	p	123/152 (81%)	119 (97%)	4 (3%)	0	100	100
All	All	10854/11700 (93%)	10373 (96%)	479 (4%)	2 (0%)	100	100

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
5	g	372	ASP
5	j	372	ASP

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	324/325 (100%)	324 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	324/325 (100%)	324 (100%)	0	100	100
1	C	324/325 (100%)	324 (100%)	0	100	100
1	D	324/325 (100%)	324 (100%)	0	100	100
1	E	324/325 (100%)	324 (100%)	0	100	100
1	F	324/325 (100%)	324 (100%)	0	100	100
1	G	322/325 (99%)	322 (100%)	0	100	100
1	H	322/325 (99%)	322 (100%)	0	100	100
1	I	322/325 (99%)	322 (100%)	0	100	100
1	J	322/325 (99%)	322 (100%)	0	100	100
1	K	322/325 (99%)	322 (100%)	0	100	100
1	L	322/325 (99%)	322 (100%)	0	100	100
2	M	143/189 (76%)	143 (100%)	0	100	100
2	N	143/189 (76%)	143 (100%)	0	100	100
2	O	143/189 (76%)	143 (100%)	0	100	100
2	P	143/189 (76%)	143 (100%)	0	100	100
2	Q	143/189 (76%)	143 (100%)	0	100	100
2	R	143/189 (76%)	143 (100%)	0	100	100
3	S	189/207 (91%)	189 (100%)	0	100	100
3	T	189/207 (91%)	189 (100%)	0	100	100
3	U	189/207 (91%)	189 (100%)	0	100	100
3	V	189/207 (91%)	189 (100%)	0	100	100
3	W	189/207 (91%)	189 (100%)	0	100	100
3	X	189/207 (91%)	189 (100%)	0	100	100
4	Y	115/122 (94%)	115 (100%)	0	100	100
4	Z	115/122 (94%)	115 (100%)	0	100	100
4	a	115/122 (94%)	115 (100%)	0	100	100
4	b	115/122 (94%)	115 (100%)	0	100	100
4	c	115/122 (94%)	115 (100%)	0	100	100
4	d	115/122 (94%)	115 (100%)	0	100	100
5	e	399/413 (97%)	399 (100%)	0	100	100
5	f	399/413 (97%)	399 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
5	g	399/413 (97%)	399 (100%)	0	100	100
5	h	399/413 (97%)	399 (100%)	0	100	100
5	i	399/413 (97%)	399 (100%)	0	100	100
5	j	399/413 (97%)	399 (100%)	0	100	100
6	k	112/139 (81%)	112 (100%)	0	100	100
6	l	112/139 (81%)	112 (100%)	0	100	100
6	m	112/139 (81%)	112 (100%)	0	100	100
6	n	112/139 (81%)	112 (100%)	0	100	100
6	o	112/139 (81%)	112 (100%)	0	100	100
6	p	112/139 (81%)	112 (100%)	0	100	100
All	All	9624/10320 (93%)	9624 (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
5	e	283	ASN
5	h	176	ASN
6	p	42	ASN
5	f	176	ASN
5	g	176	ASN

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 [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](#)

There are no chain breaks in this entry.

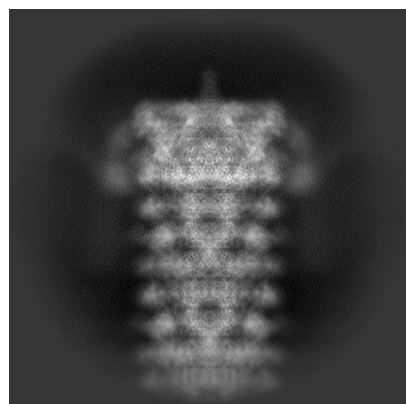
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-51195. These allow visual inspection of the internal detail of the map and identification of artifacts.

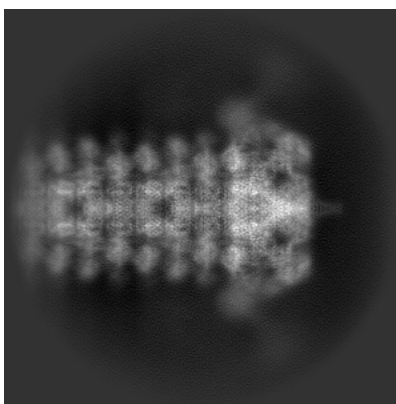
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

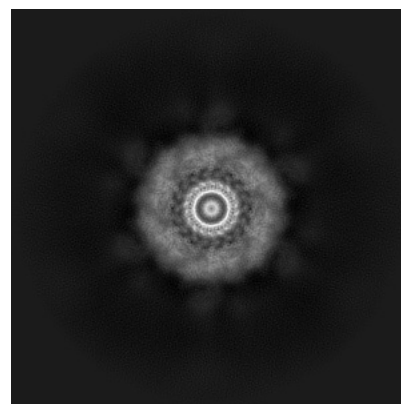
6.1.1 Primary map



X

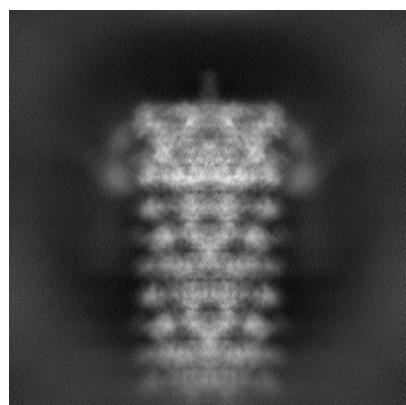


Y

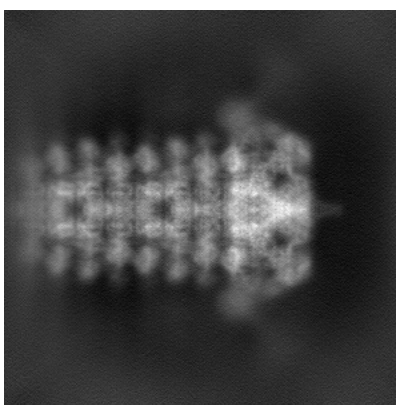


Z

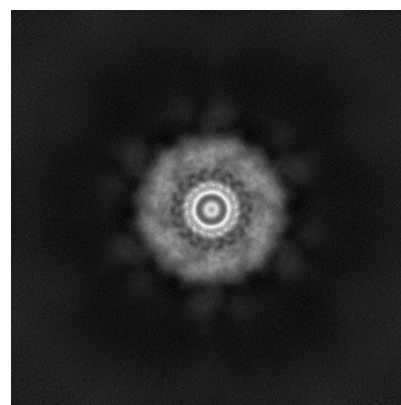
6.1.2 Raw map



X



Y

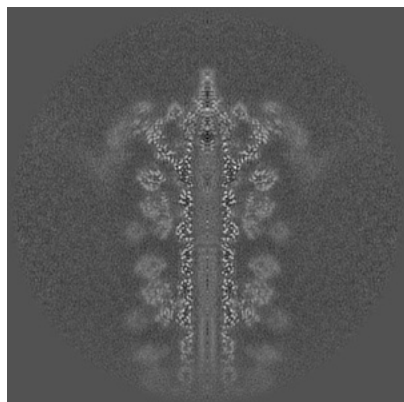


Z

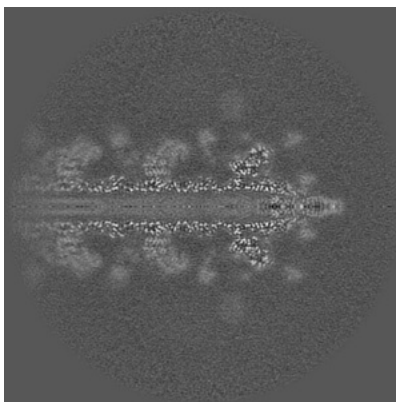
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

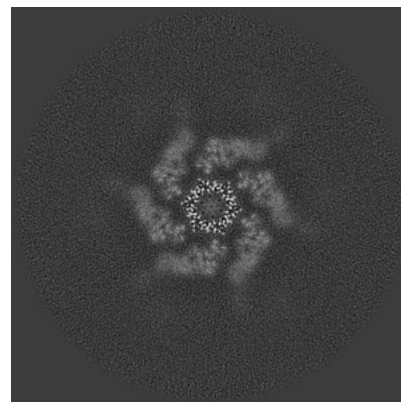
6.2.1 Primary map



X Index: 250

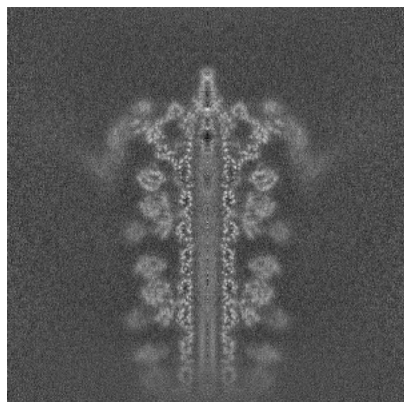


Y Index: 250

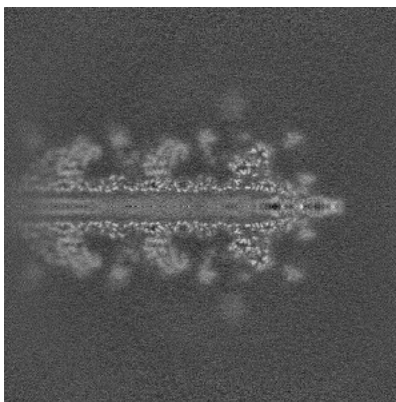


Z Index: 250

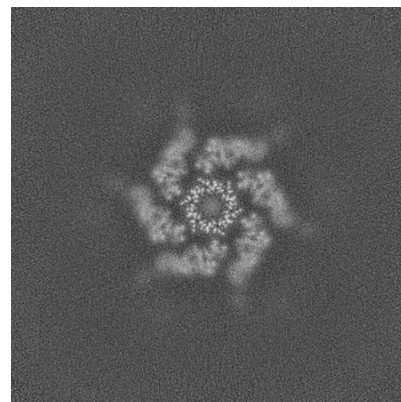
6.2.2 Raw map



X Index: 250



Y Index: 250

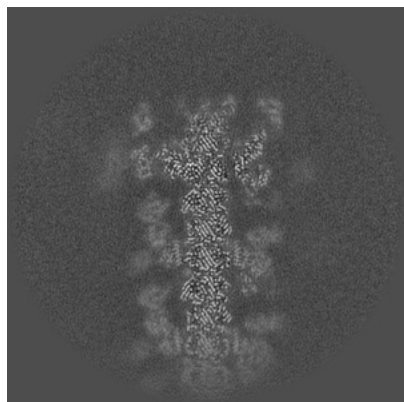


Z Index: 250

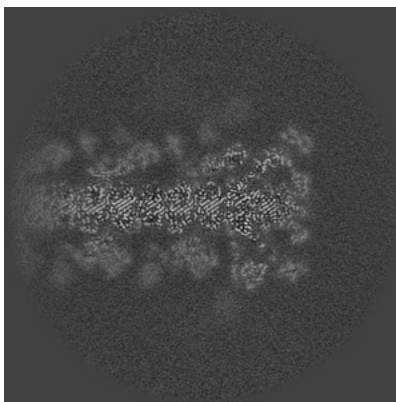
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

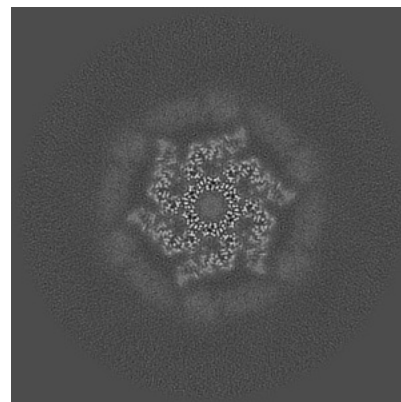
6.3.1 Primary map



X Index: 229

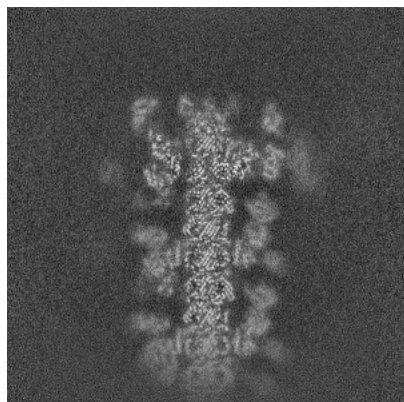


Y Index: 229

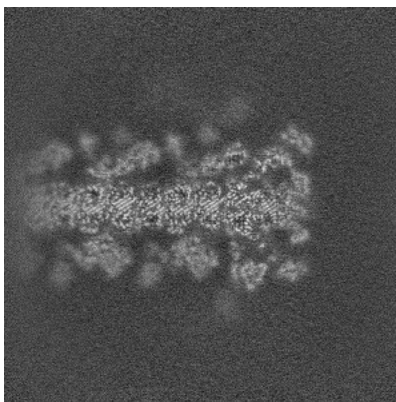


Z Index: 292

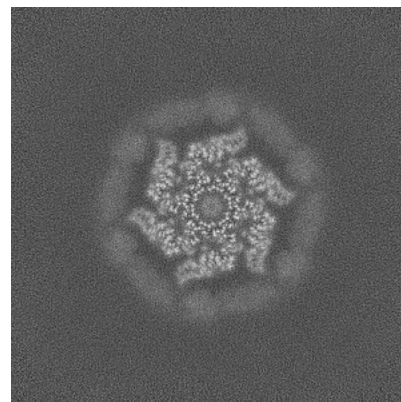
6.3.2 Raw map



X Index: 271



Y Index: 229

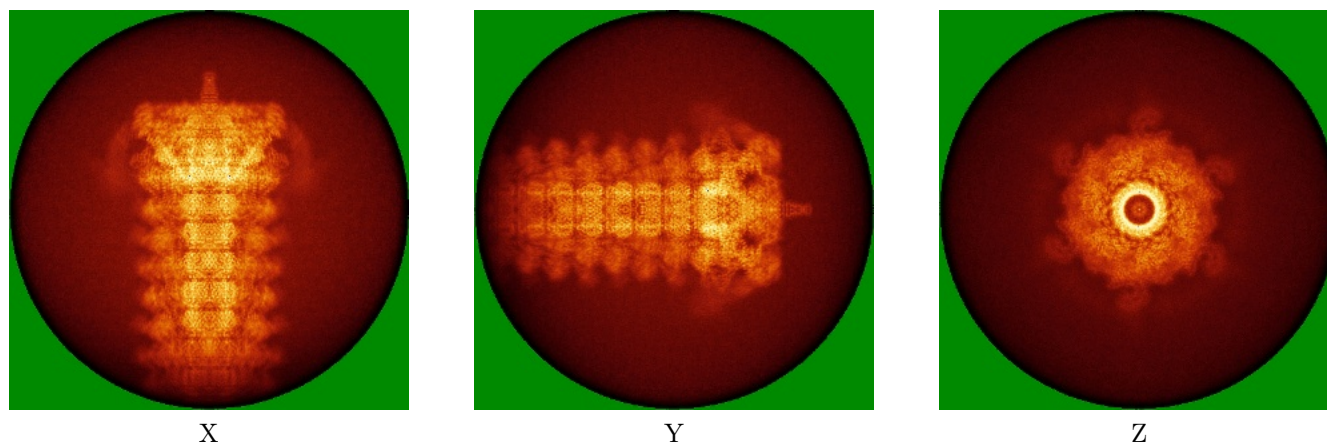


Z Index: 290

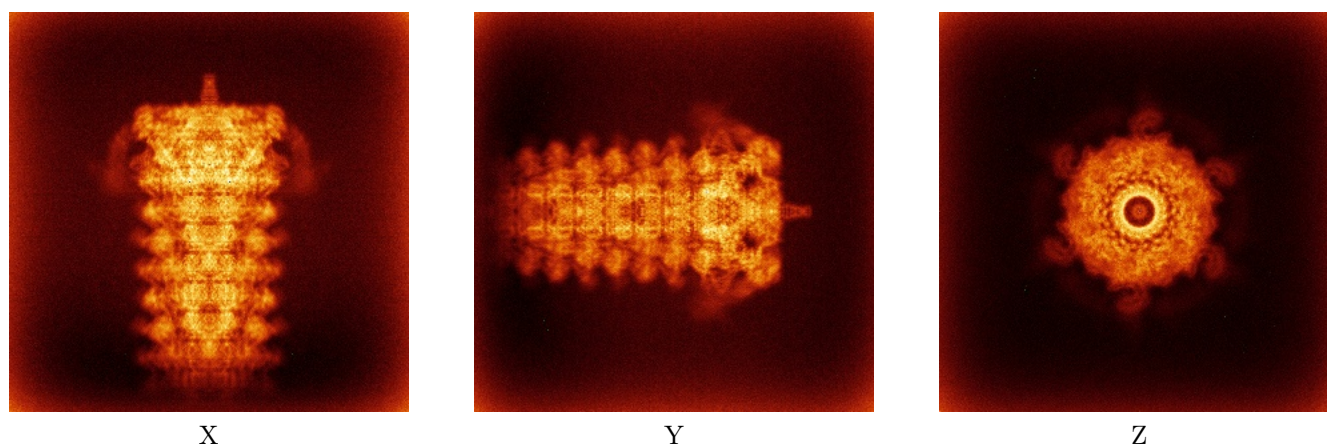
The images above show the largest variance slices of the map in three orthogonal directions.

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

6.4.1 Primary map



6.4.2 Raw map



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](#)

This section was not generated.

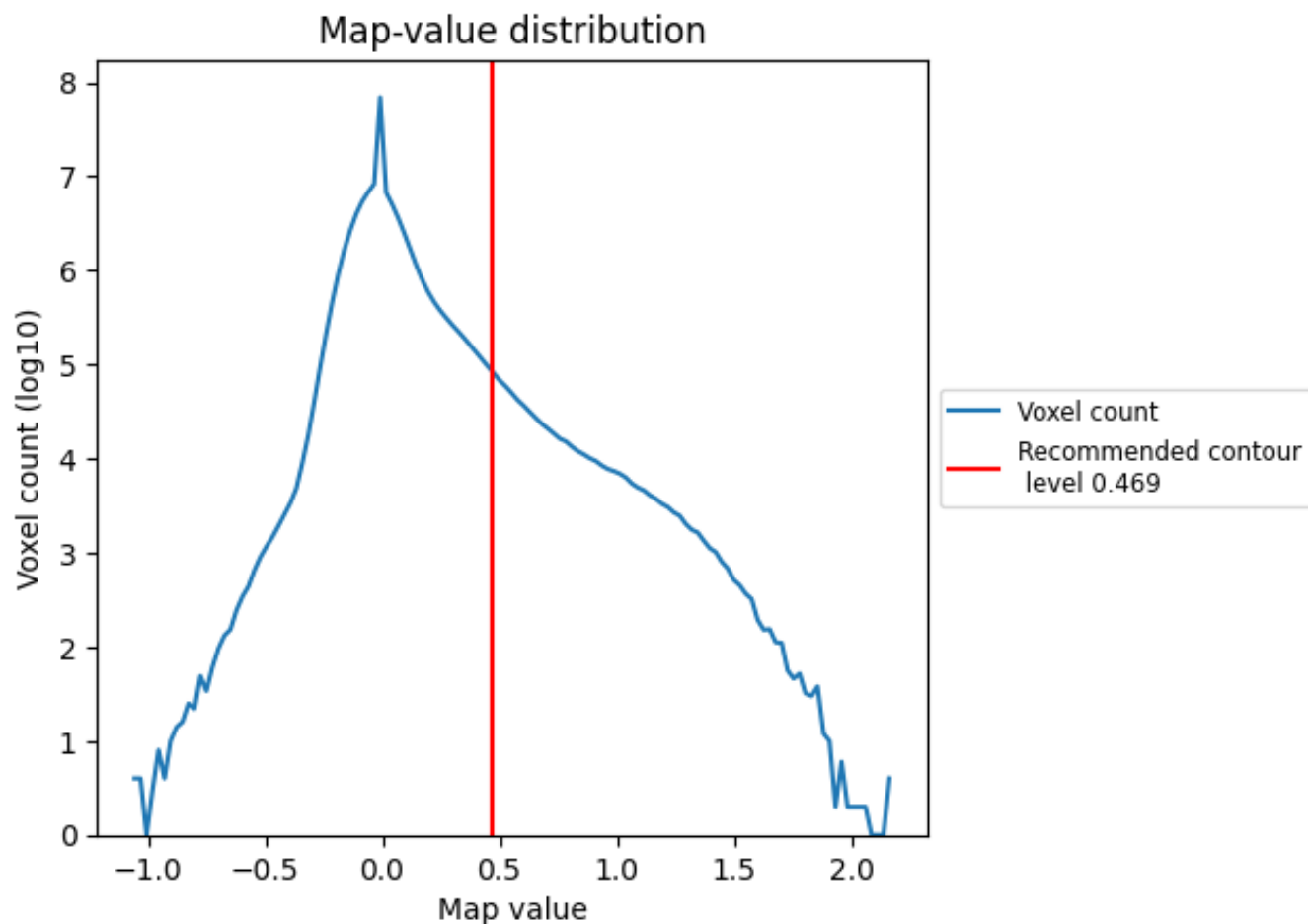
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

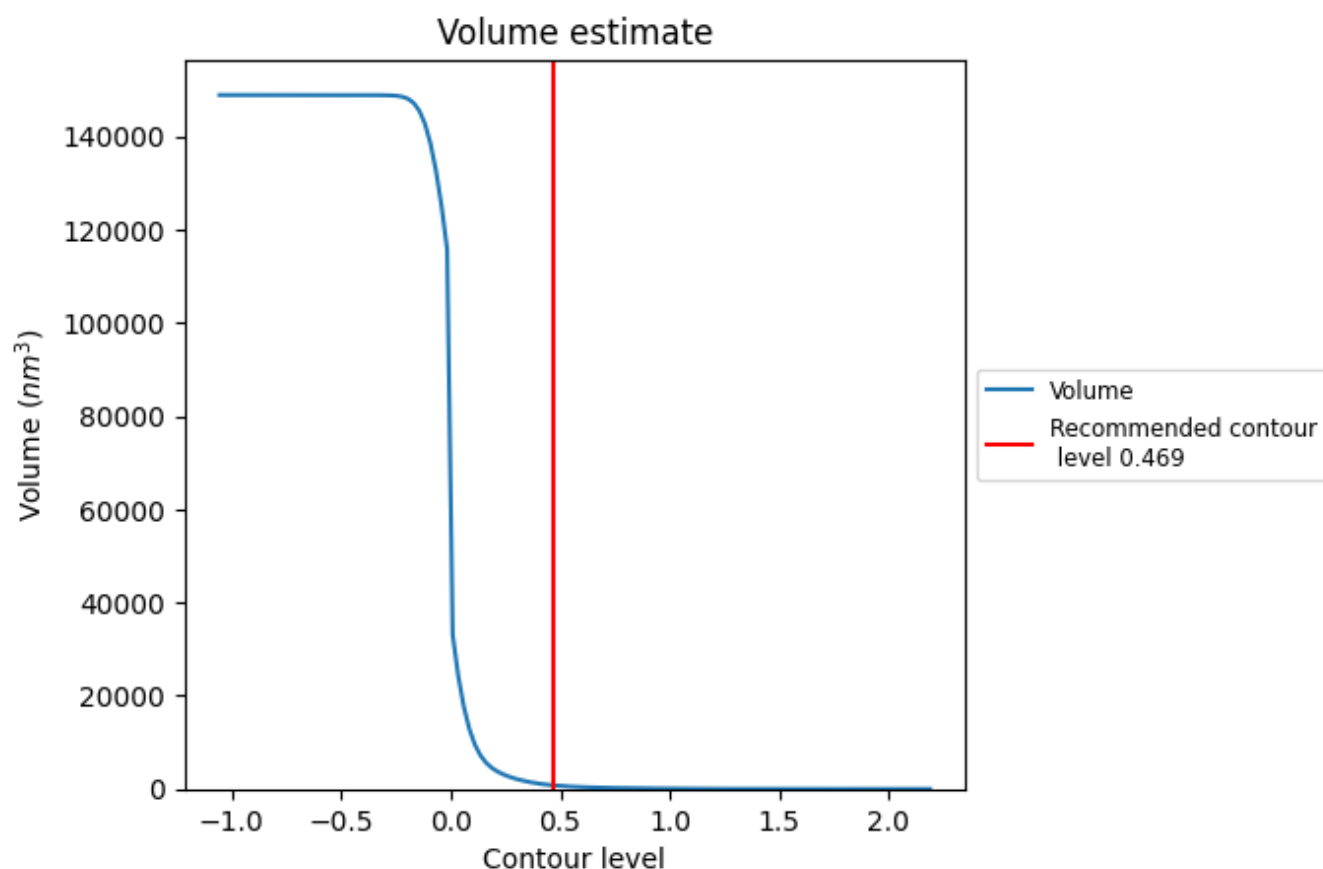
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



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

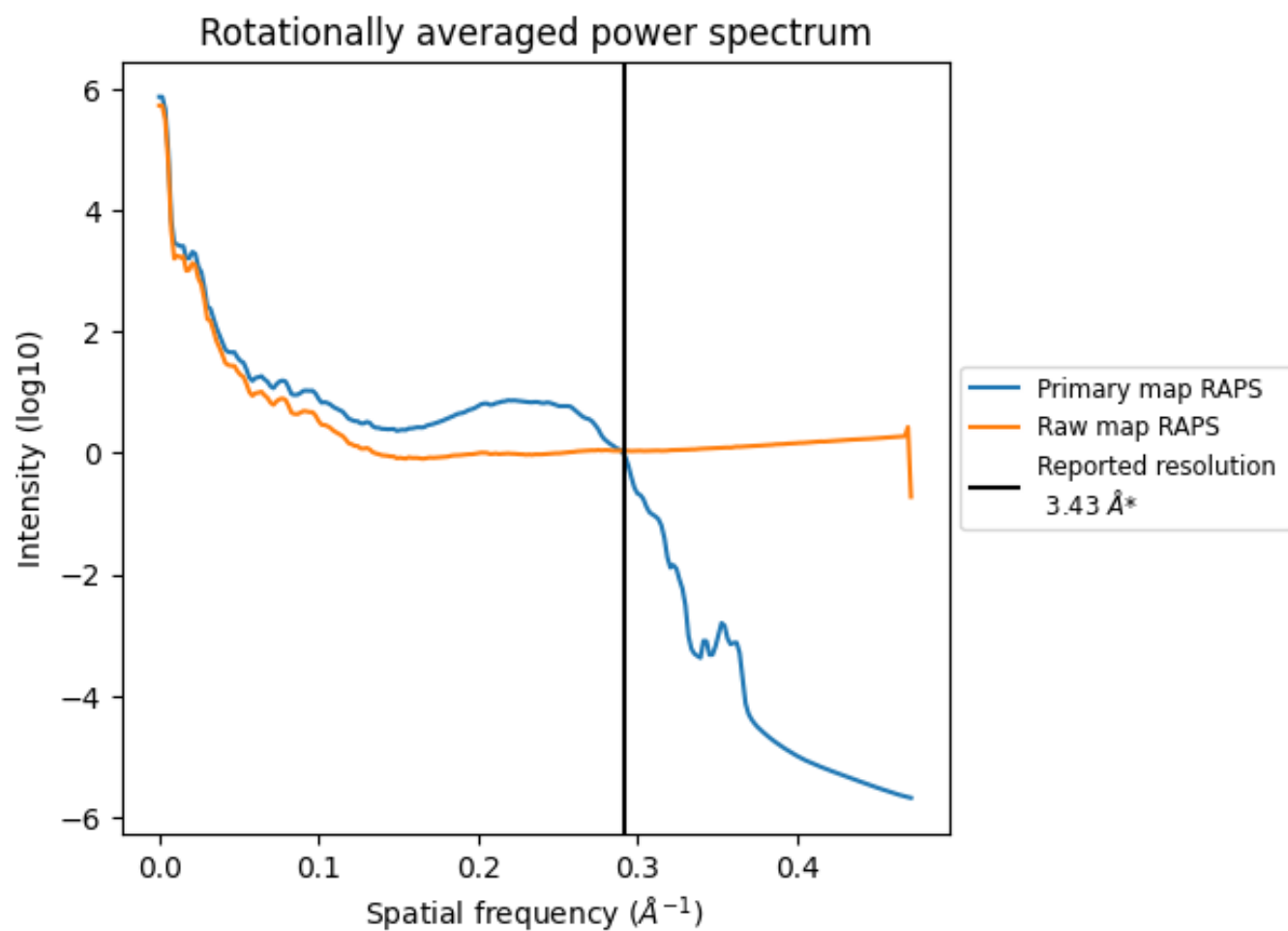
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 780 nm^3 ; this corresponds to an approximate mass of 704 kDa.

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

7.3 Rotationally averaged power spectrum ⓘ

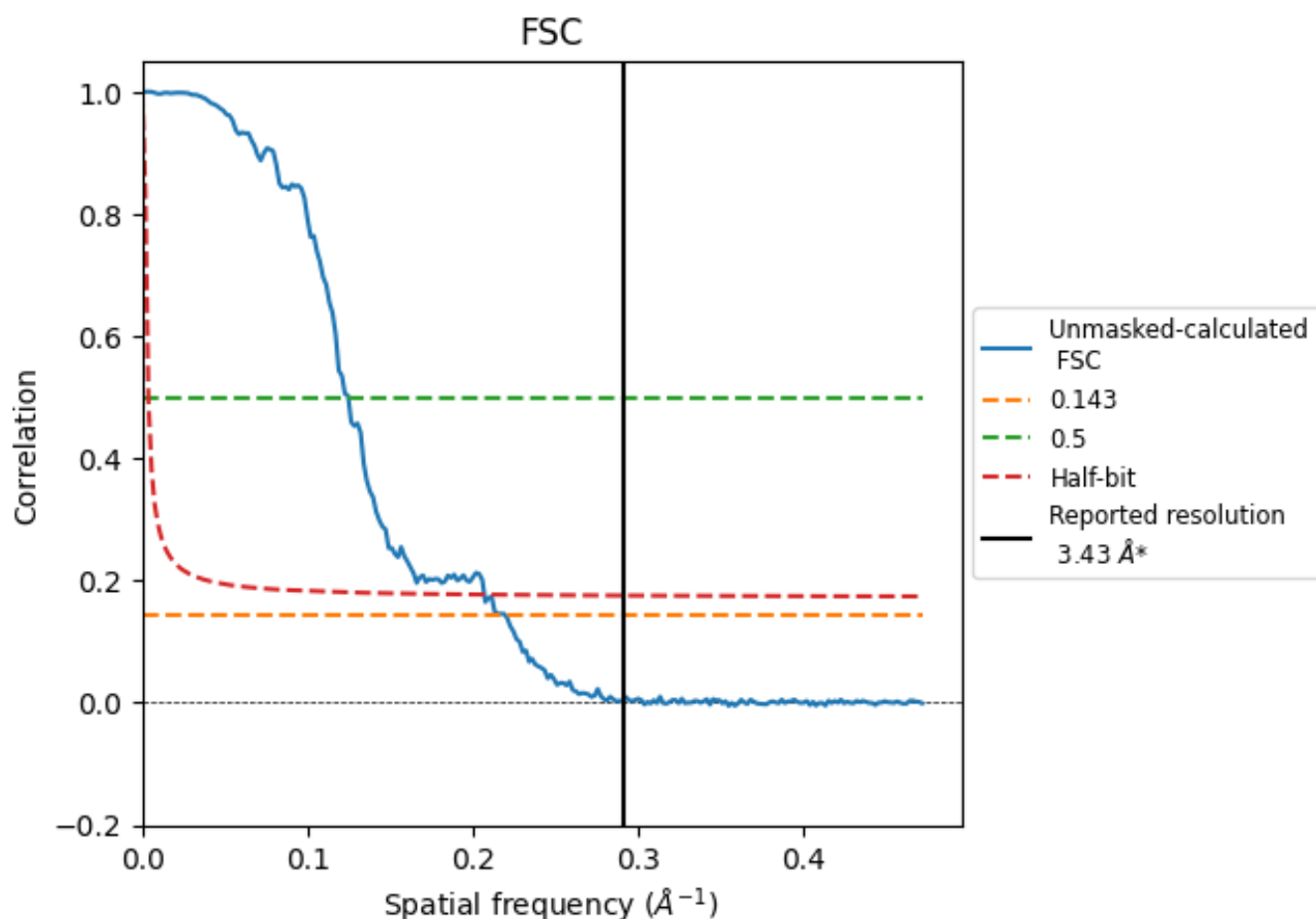


*Reported resolution corresponds to spatial frequency of $0.292 \text{ \AA}^{-1}$

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 0.292 Å⁻¹

8.2 Resolution estimates [i](#)

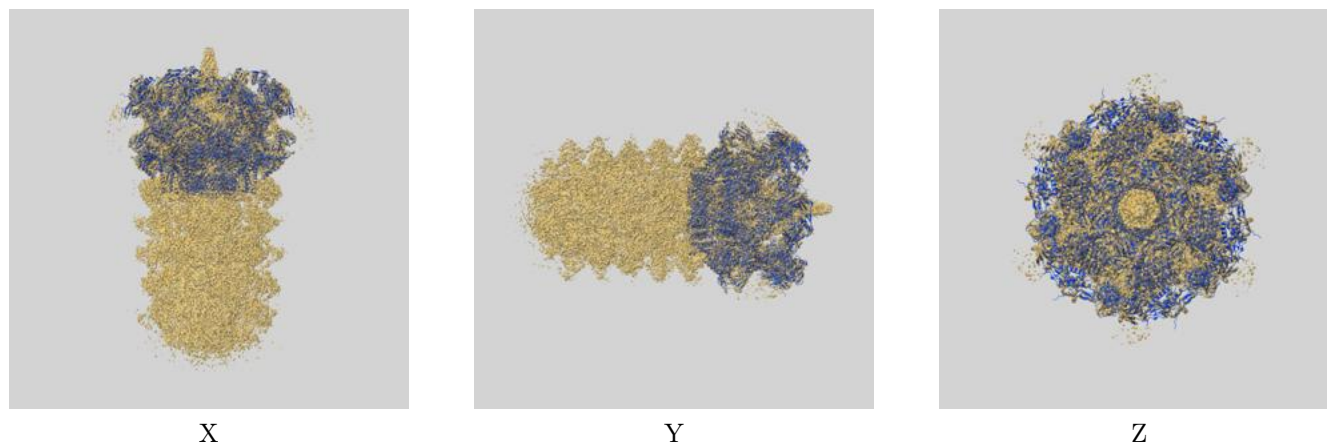
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.43	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	4.56	8.03	4.83

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 4.56 differs from the reported value 3.43 by more than 10 %

9 Map-model fit [i](#)

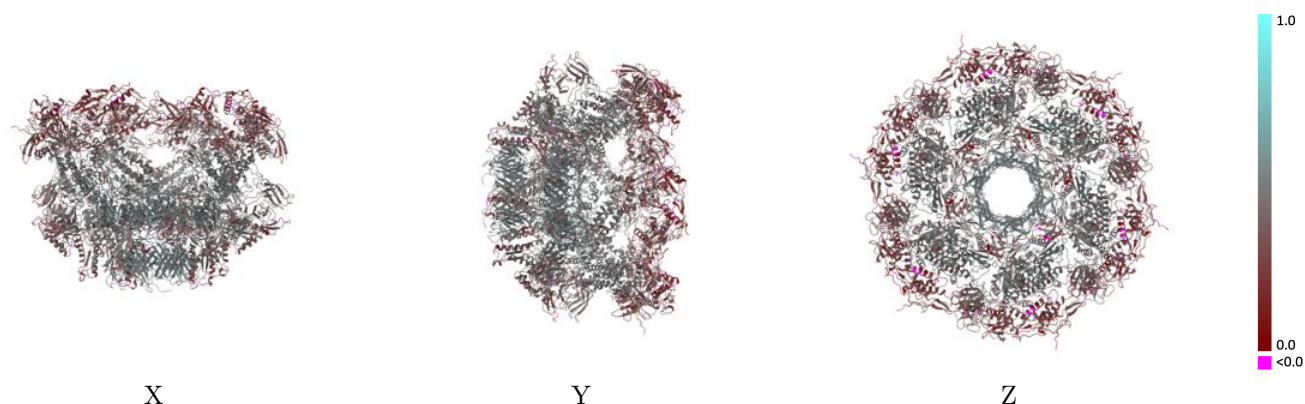
This section contains information regarding the fit between EMDB map EMD-51195 and PDB model 9GB2. Per-residue inclusion information can be found in section [3](#) on page [8](#).

9.1 Map-model overlay [i](#)



The images above show the 3D surface view of the map at the recommended contour level 0.469 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](#)

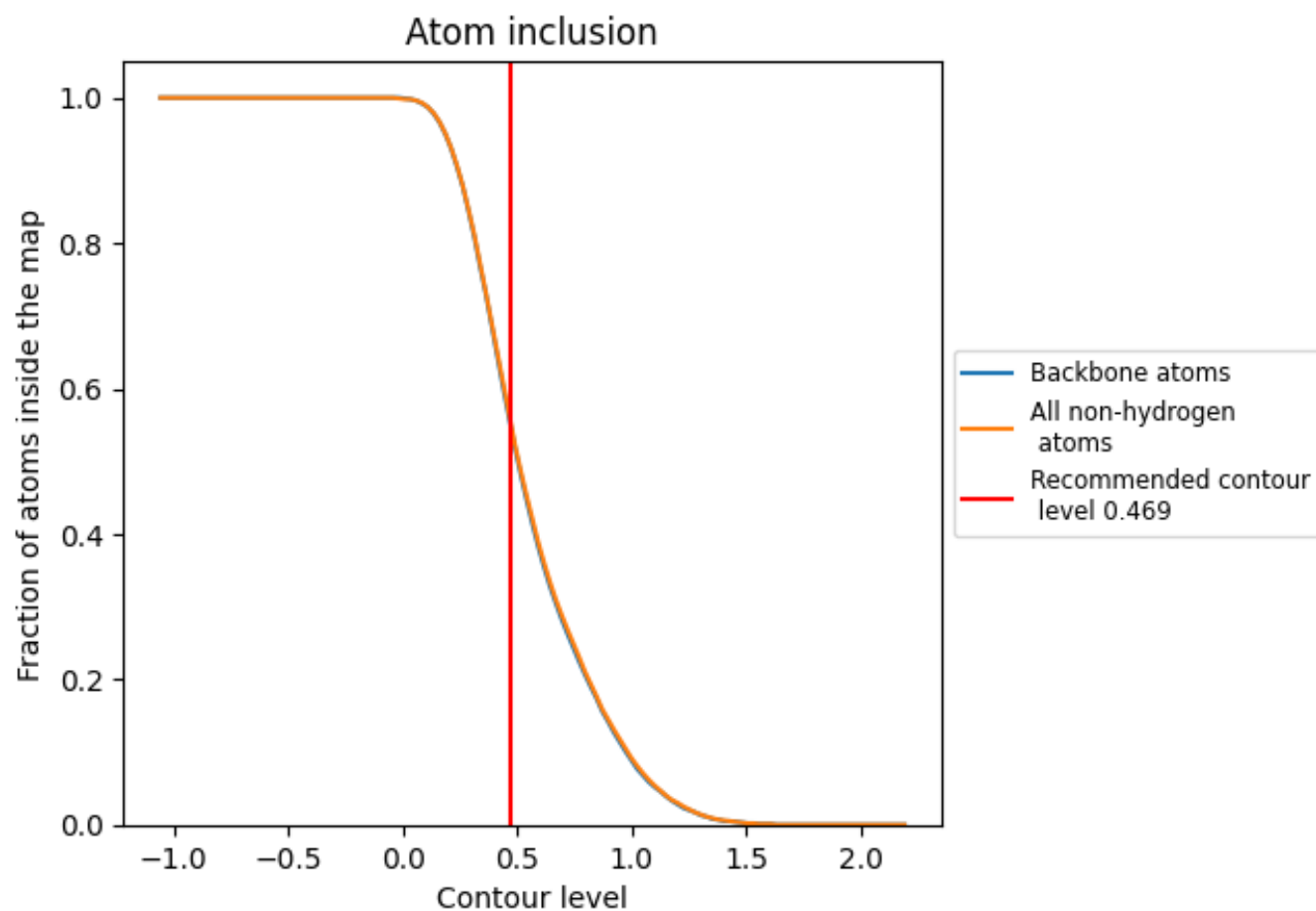


The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

This section was not generated.

9.4 Atom inclusion [i](#)



At the recommended contour level, 55% of all backbone atoms, 56% 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 (0.469) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.5550	 0.4090
A	 0.4910	 0.3810
B	 0.4860	 0.3810
C	 0.4860	 0.3810
D	 0.4910	 0.3830
E	 0.4860	 0.3840
F	 0.4860	 0.3820
G	 0.4030	 0.3180
H	 0.4050	 0.3170
I	 0.4040	 0.3160
J	 0.4030	 0.3170
K	 0.4050	 0.3180
L	 0.4040	 0.3190
M	 0.5500	 0.3980
N	 0.5470	 0.3970
O	 0.5450	 0.4010
P	 0.5500	 0.4030
Q	 0.5470	 0.4000
R	 0.5450	 0.4010
S	 0.7420	 0.5040
T	 0.7330	 0.5030
U	 0.7340	 0.5030
V	 0.7420	 0.5050
W	 0.7330	 0.5030
X	 0.7340	 0.5020
Y	 0.7200	 0.5020
Z	 0.7120	 0.5020
a	 0.7160	 0.5020
b	 0.7200	 0.5020
c	 0.7120	 0.5030
d	 0.7160	 0.5010
e	 0.5780	 0.4120
f	 0.5800	 0.4130
g	 0.5780	 0.4120
h	 0.5780	 0.4110



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Chain	Atom inclusion	Q-score
i	 0.5810	 0.4110
j	 0.5780	 0.4110
k	 0.7190	 0.4930
l	 0.7270	 0.4970
m	 0.7200	 0.4950
n	 0.7190	 0.4950
o	 0.7260	 0.4960
p	 0.7220	 0.4960