



wwPDB EM Validation Summary Report ⓘ

Mar 14, 2026 – 07:15 AM UTC

PDB ID : 6WXG / pdb_00006wxg
EMDB ID : EMD-21957
Title : Cryo-EM reconstruction of VP5*/VP8* assembly from rhesus rotavirus particles - Reversed conformation
Authors : Herrmann, T.; Harrison, S.C.; Jenni, S.
Deposited on : 2020-05-10
Resolution : 3.30 Å (reported)
Based on initial models : 4V7Q, 1SLQ

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

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

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

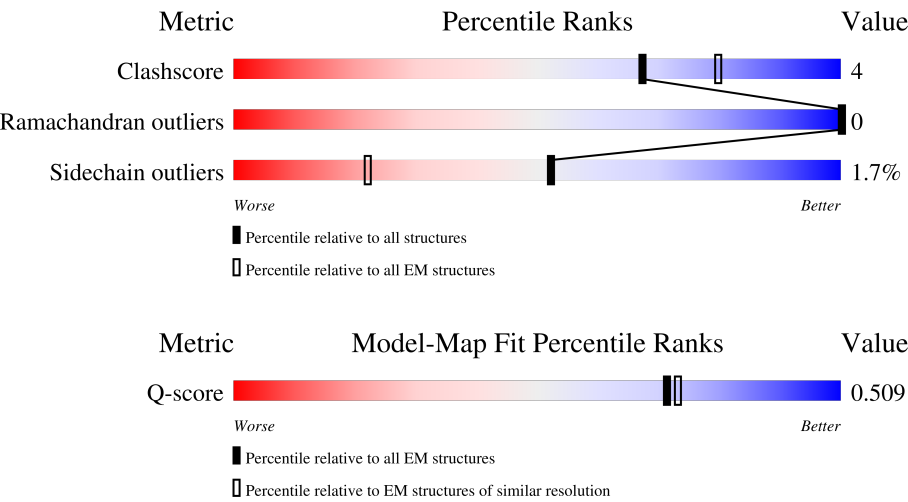
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

1 Overall quality at a glance i

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

The reported resolution of this entry is 3.30 Å.

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



Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	15087 (2.80 - 3.80)

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	1	776	34% 31% 65%
1	2	776	33% 32% 65%
1	3	776	33% 31% 65%
2	A	397	28% 88% 12%

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Mol	Chain	Length	Quality of chain
2	B	397	32% 90% 10%
2	C	397	31% 89% 11%
2	D	397	29% 88% 12%
2	E	397	31% 87% 12%
2	F	397	29% 88% 11%
2	G	397	29% 88% 12%
2	H	397	32% 90% 10%
2	I	397	31% 89% 10%
2	J	397	36% 90% 10%
2	K	397	35% 88% 12%
2	L	397	37% 88% 12%
2	M	397	34% 89% 11%
2	N	397	37% 91% 9%
2	O	397	39% 90% 9%
2	P	397	34% 87% 13%
2	Q	397	32% 90% 10%
2	R	397	38% 90% 10%
3	a	326	51% 69% 11% 21%
3	b	326	48% 71% 13% 17%
3	c	326	44% 71% 9% 20%
3	d	326	43% 72% 10% 19%
3	e	326	40% 67% 12% 20%
3	f	326	44% 71% 10% 19%
3	g	326	54% 74% 9% 17%
3	h	326	50% 71% 8% 21%

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Mol	Chain	Length	Quality of chain
3	i	326	48% 74% 10% 17%
3	j	326	52% 73% 10% 17%
3	k	326	51% 73% 8% 19%
3	l	326	47% 71% 10% 19%
3	m	326	51% 73% 10% 17%
3	n	326	52% 73% 9% 19%
3	o	326	50% 73% 8% 19%
3	p	326	56% 75% 8% 17%
3	q	326	52% 73% 6% 21%
3	r	326	53% 74% 9% 17%

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 201212 atoms, of which 99503 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 Outer capsid protein VP4.

Mol	Chain	Residues	Atoms						AltConf	Trace
1	1	275	Total	C	H	N	O	S	0	0
			4291	1382	2107	368	427	7		
1	2	275	Total	C	H	N	O	S	0	0
			4291	1382	2107	368	427	7		
1	3	275	Total	C	H	N	O	S	0	0
			4291	1382	2107	368	427	7		

- Molecule 2 is a protein called Intermediate capsid protein VP6.

Mol	Chain	Residues	Atoms						AltConf	Trace
2	A	397	Total	C	H	N	O	S	0	0
			6275	2004	3112	551	593	15		
2	B	397	Total	C	H	N	O	S	0	0
			6275	2004	3112	551	593	15		
2	C	397	Total	C	H	N	O	S	0	0
			6275	2004	3112	551	593	15		
2	D	397	Total	C	H	N	O	S	0	0
			6275	2004	3112	551	593	15		
2	E	397	Total	C	H	N	O	S	0	0
			6276	2004	3113	551	593	15		
2	F	397	Total	C	H	N	O	S	0	0
			6276	2004	3113	551	593	15		
2	G	397	Total	C	H	N	O	S	0	0
			6276	2004	3113	551	593	15		
2	H	397	Total	C	H	N	O	S	0	0
			6276	2004	3113	551	593	15		
2	I	397	Total	C	H	N	O	S	0	0
			6275	2004	3112	551	593	15		
2	J	397	Total	C	H	N	O	S	0	0
			6275	2004	3112	551	593	15		
2	K	397	Total	C	H	N	O	S	0	0
			6275	2004	3112	551	593	15		
2	L	397	Total	C	H	N	O	S	0	0
			6276	2004	3113	551	593	15		

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Mol	Chain	Residues	Atoms						AltConf	Trace
2	M	397	Total	C	H	N	O	S	0	0
			6276	2004	3113	551	593	15		
2	N	397	Total	C	H	N	O	S	0	0
			6275	2004	3112	551	593	15		
2	O	397	Total	C	H	N	O	S	0	0
			6276	2004	3113	551	593	15		
2	P	397	Total	C	H	N	O	S	0	0
			6275	2004	3112	551	593	15		
2	Q	397	Total	C	H	N	O	S	0	0
			6276	2004	3113	551	593	15		
2	R	397	Total	C	H	N	O	S	0	0
			6275	2004	3112	551	593	15		

- Molecule 3 is a protein called Outer capsid glycoprotein VP7.

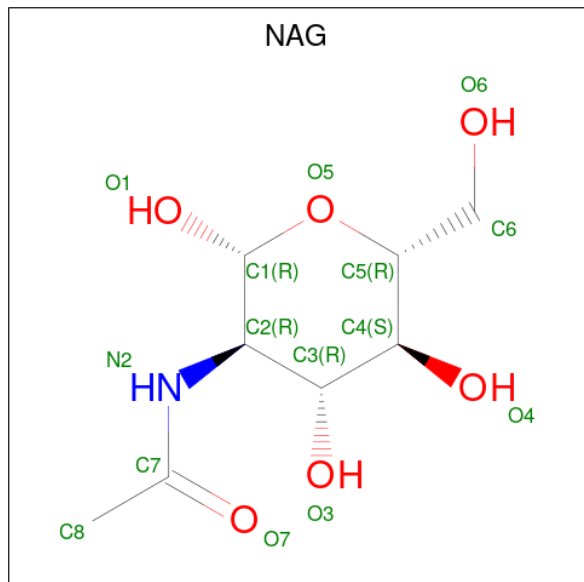
Mol	Chain	Residues	Atoms						AltConf	Trace
3	a	259	Total	C	H	N	O	S	0	0
			4041	1298	1994	324	409	16		
3	b	272	Total	C	H	N	O	S	0	0
			4254	1369	2099	342	428	16		
3	c	261	Total	C	H	N	O	S	0	0
			4074	1308	2011	327	412	16		
3	d	265	Total	C	H	N	O	S	0	0
			4133	1328	2037	333	419	16		
3	e	260	Total	C	H	N	O	S	0	0
			4050	1302	1998	323	411	16		
3	f	265	Total	C	H	N	O	S	0	0
			4133	1328	2037	333	419	16		
3	g	272	Total	C	H	N	O	S	0	0
			4254	1369	2099	342	428	16		
3	h	259	Total	C	H	N	O	S	0	0
			4041	1298	1994	324	409	16		
3	i	272	Total	C	H	N	O	S	0	0
			4254	1369	2099	342	428	16		
3	j	272	Total	C	H	N	O	S	0	0
			4254	1369	2099	342	428	16		
3	k	265	Total	C	H	N	O	S	0	0
			4133	1328	2037	333	419	16		
3	l	265	Total	C	H	N	O	S	0	0
			4133	1328	2037	333	419	16		
3	m	272	Total	C	H	N	O	S	0	0
			4254	1369	2099	342	428	16		

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Mol	Chain	Residues	Atoms						AltConf	Trace
3	n	265	Total	C	H	N	O	S	0	0
			4133	1328	2037	333	419	16		
3	o	265	Total	C	H	N	O	S	0	0
			4133	1328	2037	333	419	16		
3	p	272	Total	C	H	N	O	S	0	0
			4254	1369	2099	342	428	16		
3	q	259	Total	C	H	N	O	S	0	0
			4041	1298	1994	324	409	16		
3	r	272	Total	C	H	N	O	S	0	0
			4254	1369	2099	342	428	16		

- Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms					AltConf
4	a	1	Total	C	H	N	O	0
			28	8	14	1	5	
4	b	1	Total	C	H	N	O	0
			28	8	14	1	5	
4	c	1	Total	C	H	N	O	0
			28	8	14	1	5	
4	d	1	Total	C	H	N	O	0
			28	8	14	1	5	
4	e	1	Total	C	H	N	O	0
			28	8	14	1	5	
4	f	1	Total	C	H	N	O	0
			28	8	14	1	5	

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Mol	Chain	Residues	Atoms					AltConf
4	g	1	Total	C	H	N	O	0
			28	8	14	1	5	
4	h	1	Total	C	H	N	O	0
			28	8	14	1	5	
4	i	1	Total	C	H	N	O	0
			28	8	14	1	5	
4	j	1	Total	C	H	N	O	0
			28	8	14	1	5	
4	k	1	Total	C	H	N	O	0
			28	8	14	1	5	
4	l	1	Total	C	H	N	O	0
			28	8	14	1	5	
4	m	1	Total	C	H	N	O	0
			28	8	14	1	5	
4	n	1	Total	C	H	N	O	0
			28	8	14	1	5	
4	o	1	Total	C	H	N	O	0
			28	8	14	1	5	
4	p	1	Total	C	H	N	O	0
			28	8	14	1	5	
4	q	1	Total	C	H	N	O	0
			28	8	14	1	5	
4	r	1	Total	C	H	N	O	0
			28	8	14	1	5	

- Molecule 5 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		AltConf
5	a	5	Total	Ca	0
			5	5	
5	b	3	Total	Ca	0
			3	3	
5	c	1	Total	Ca	0
			1	1	
5	d	5	Total	Ca	0
			5	5	
5	e	3	Total	Ca	0
			3	3	
5	f	1	Total	Ca	0
			1	1	
5	g	5	Total	Ca	0
			5	5	

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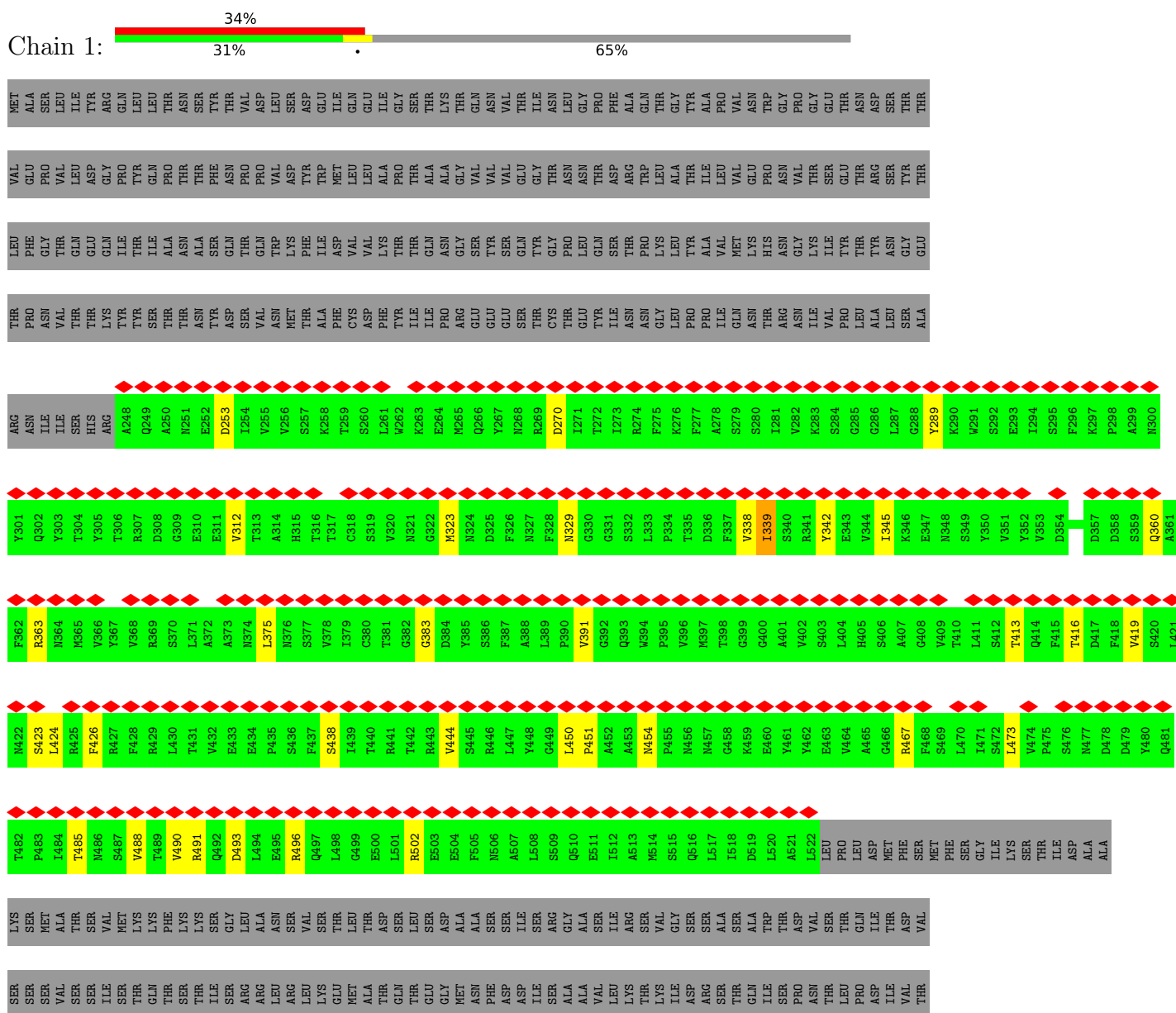
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Mol	Chain	Residues	Atoms		AltConf
5	h	3	Total 3	Ca 3	0
5	i	1	Total 1	Ca 1	0
5	j	5	Total 5	Ca 5	0
5	k	3	Total 3	Ca 3	0
5	l	1	Total 1	Ca 1	0
5	m	5	Total 5	Ca 5	0
5	n	3	Total 3	Ca 3	0
5	o	1	Total 1	Ca 1	0
5	p	5	Total 5	Ca 5	0
5	q	3	Total 3	Ca 3	0
5	r	1	Total 1	Ca 1	0

3 Residue-property plots

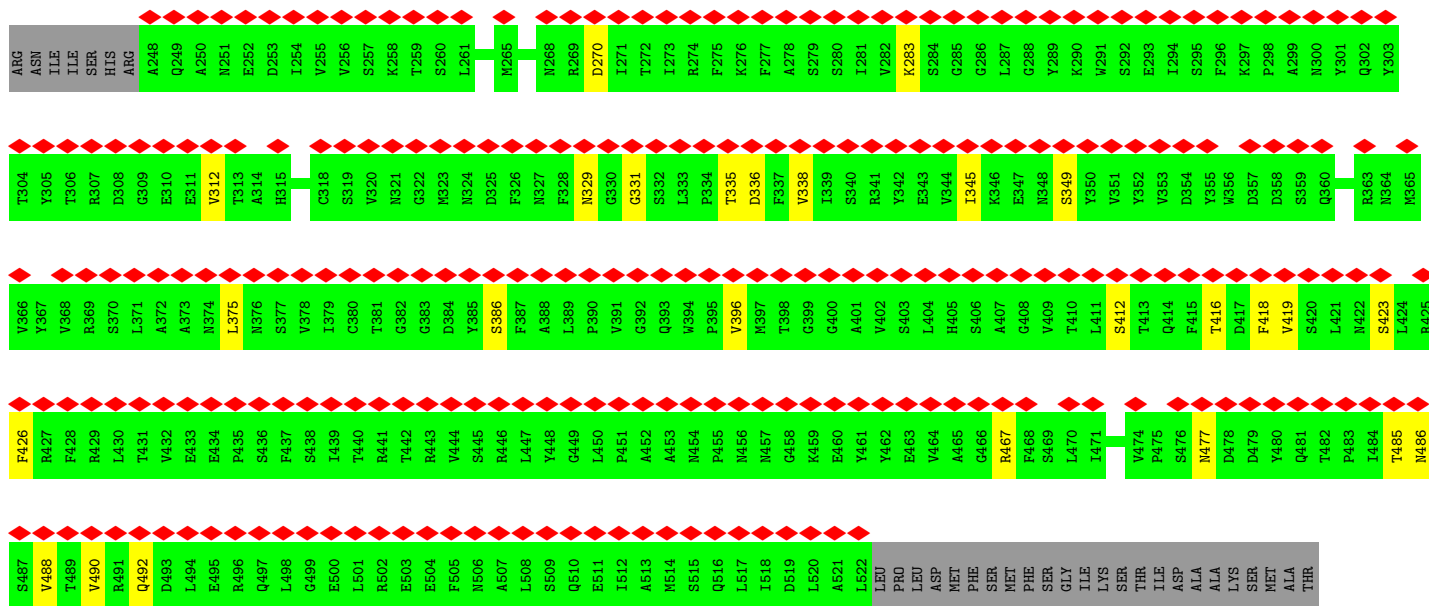
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion in map density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red diamond above a residue indicates a poor fit to the EM map for this residue (all-atom inclusion < 40%). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: Outer capsid protein VP4



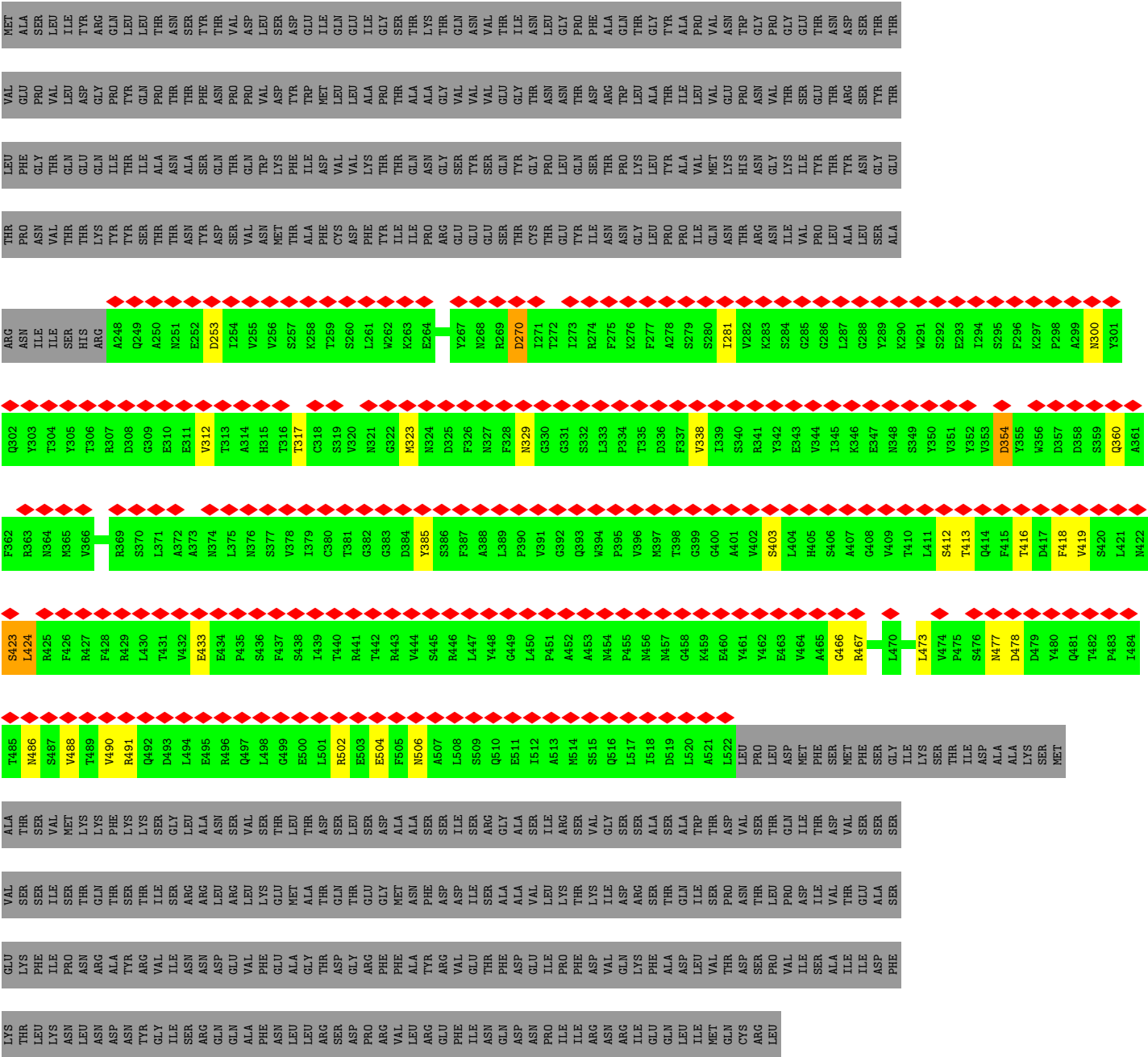
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- Molecule 1: Outer capsid protein VP4

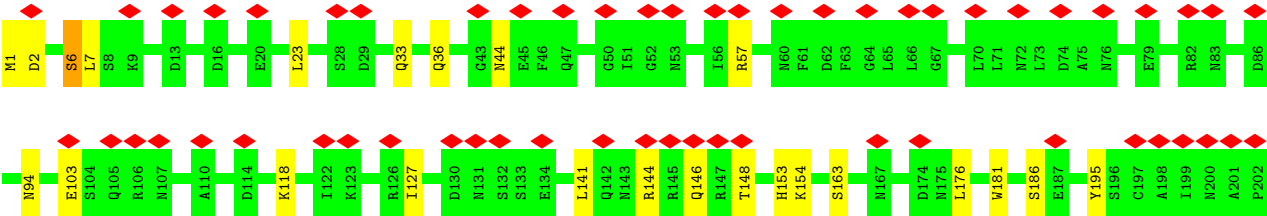
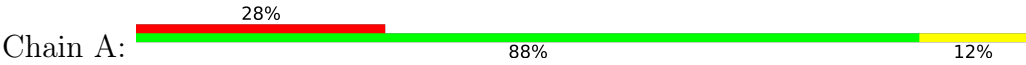
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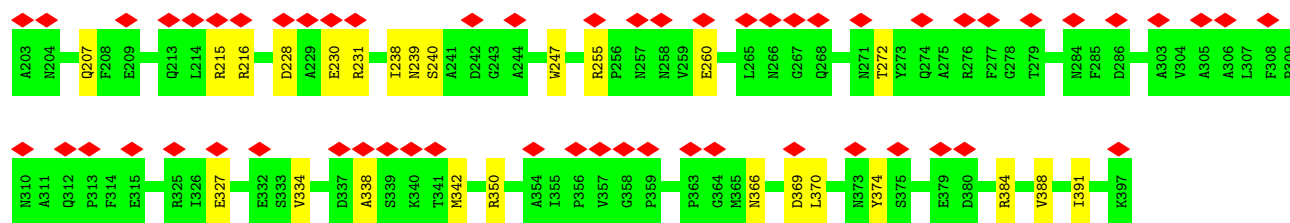
- Molecule 1: Outer capsid protein VP4



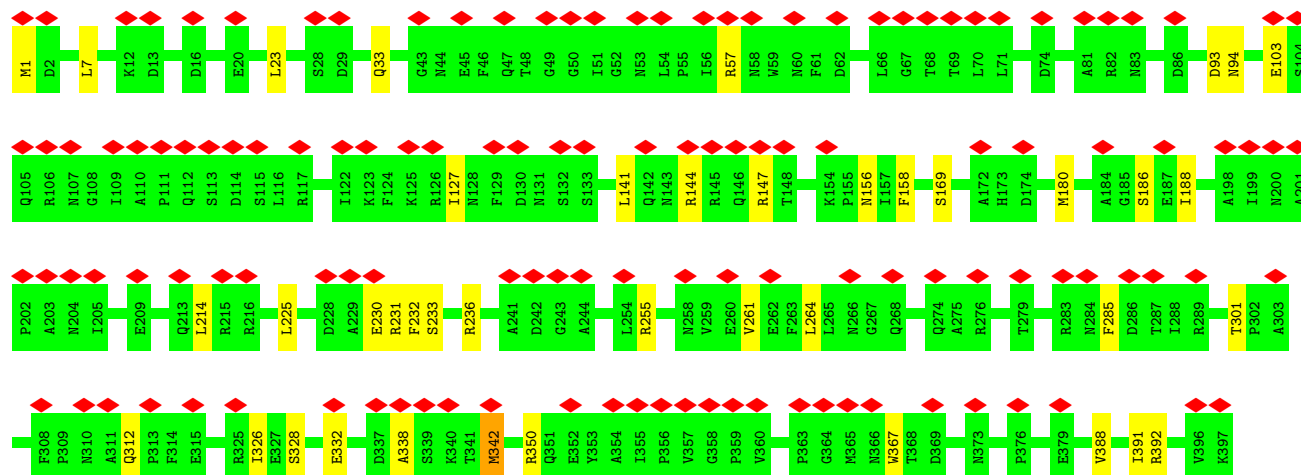
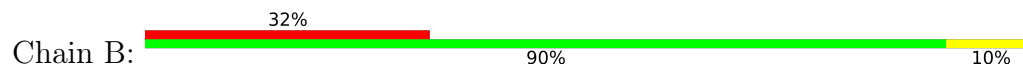


● Molecule 2: Intermediate capsid protein VP6

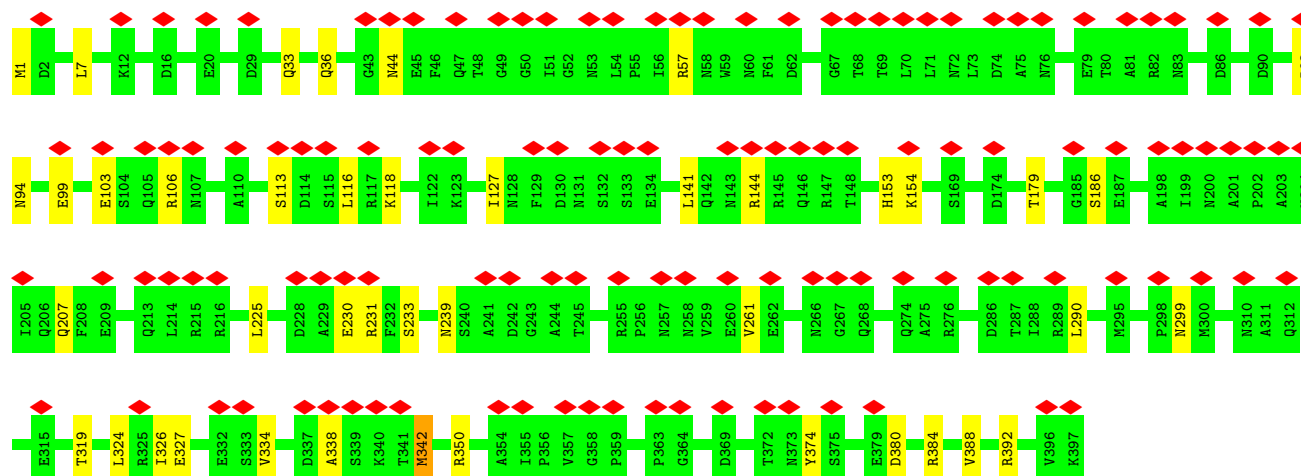
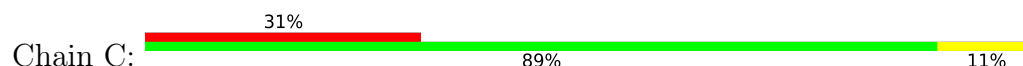




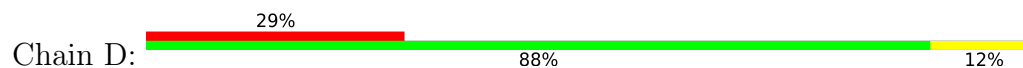
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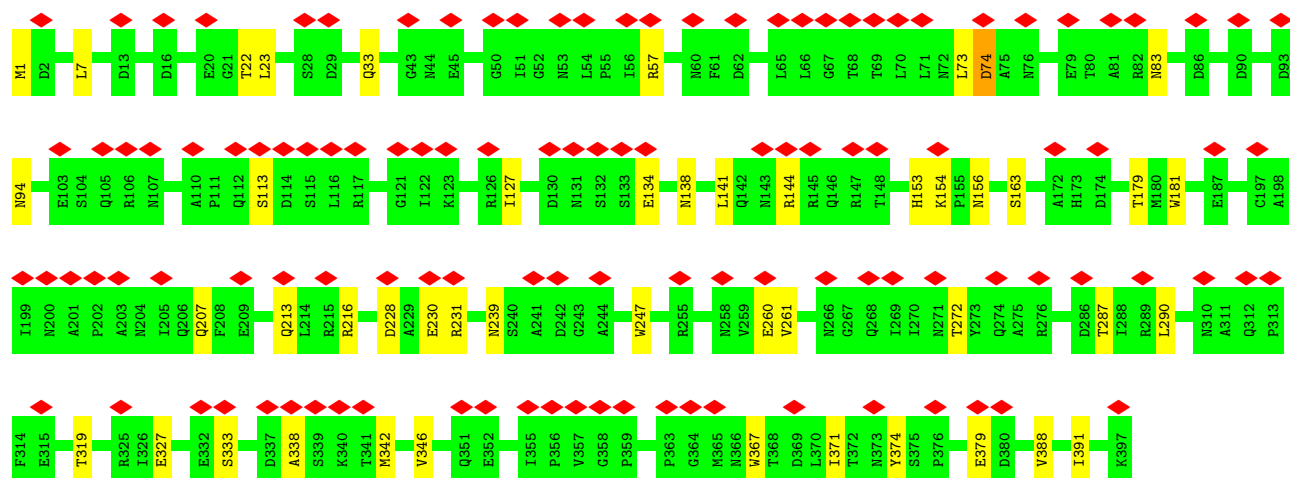


• Molecule 2: Intermediate capsid protein VP6

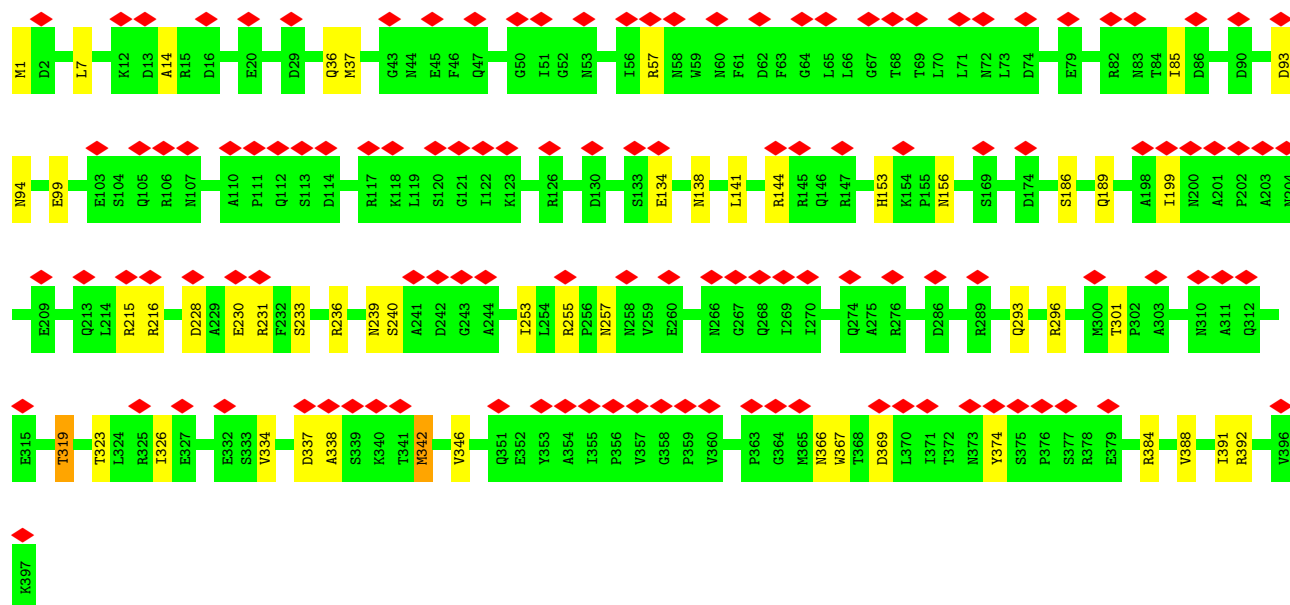
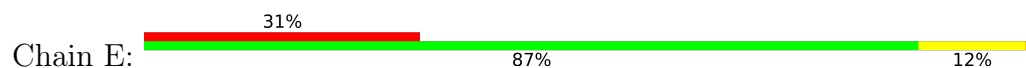


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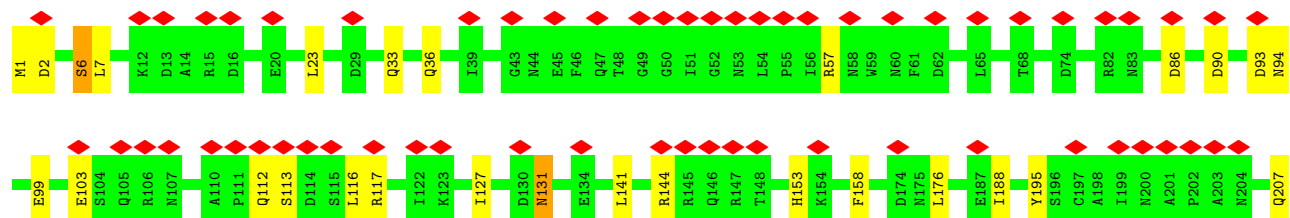
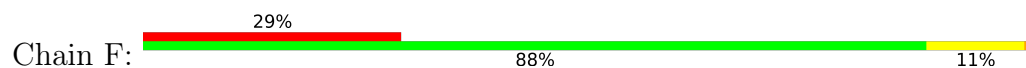


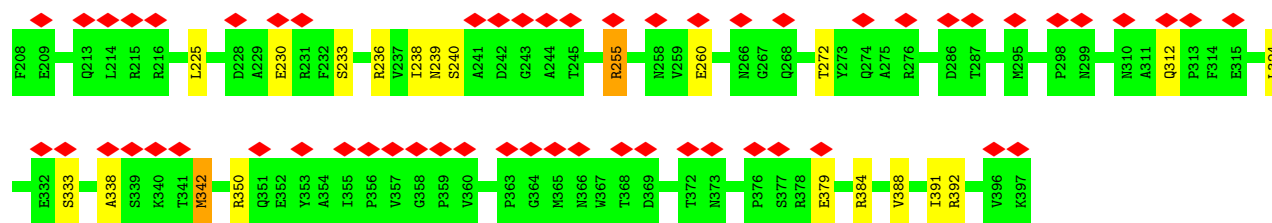


• Molecule 2: Intermediate capsid protein VP6

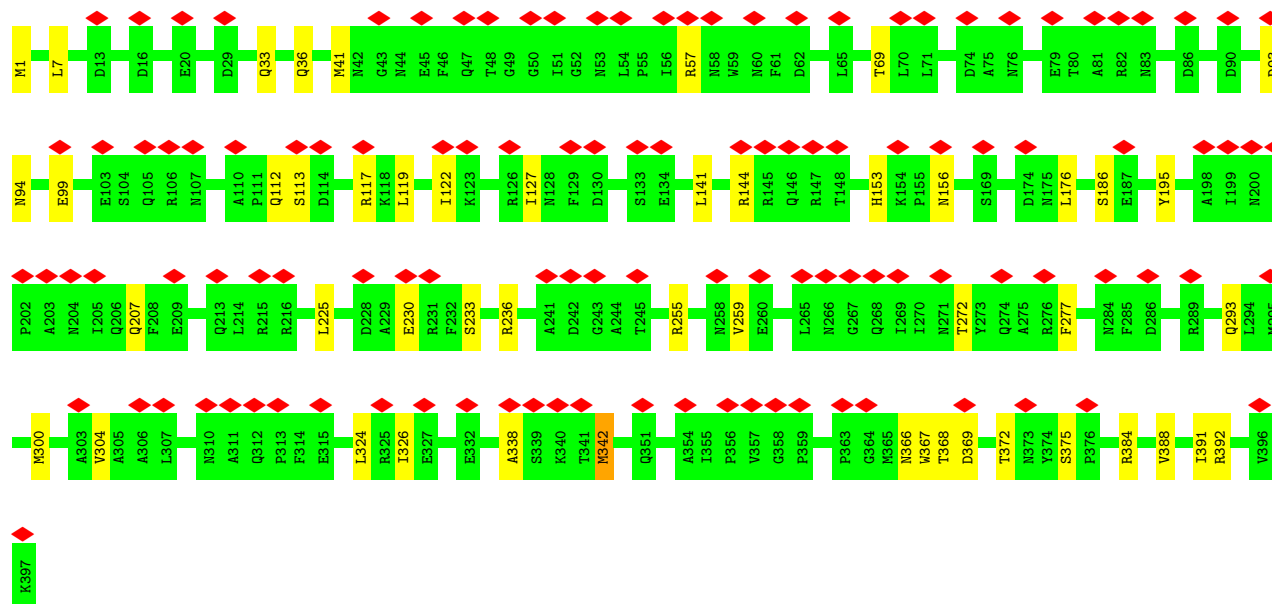
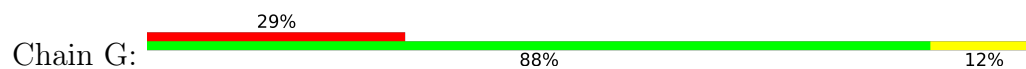


• Molecule 2: Intermediate capsid protein VP6

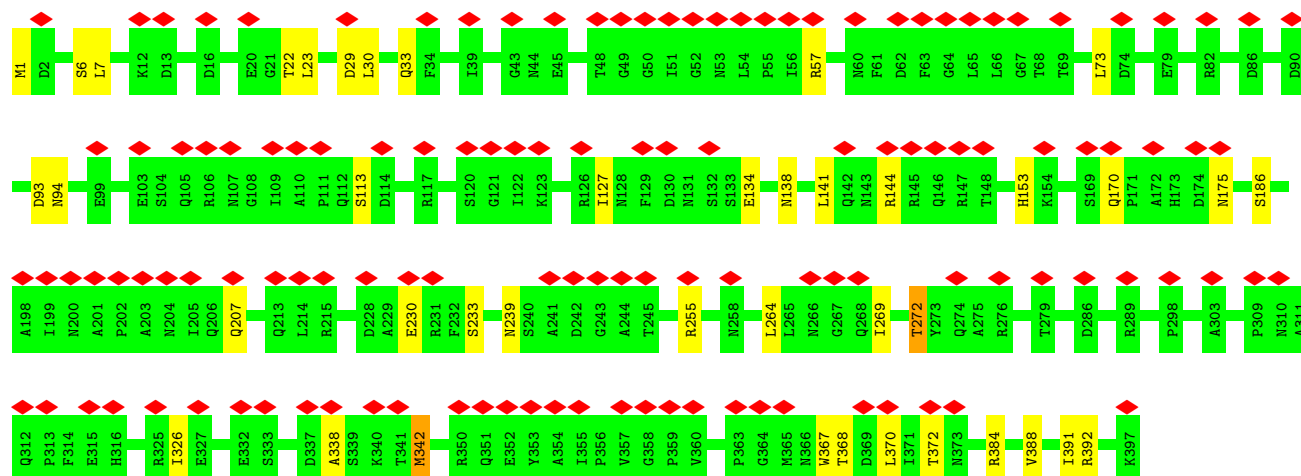
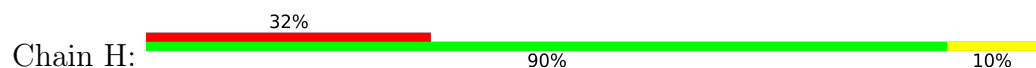




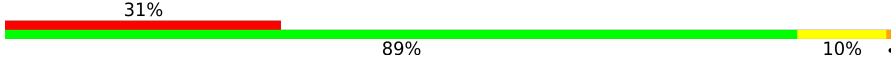
• Molecule 2: Intermediate capsid protein VP6

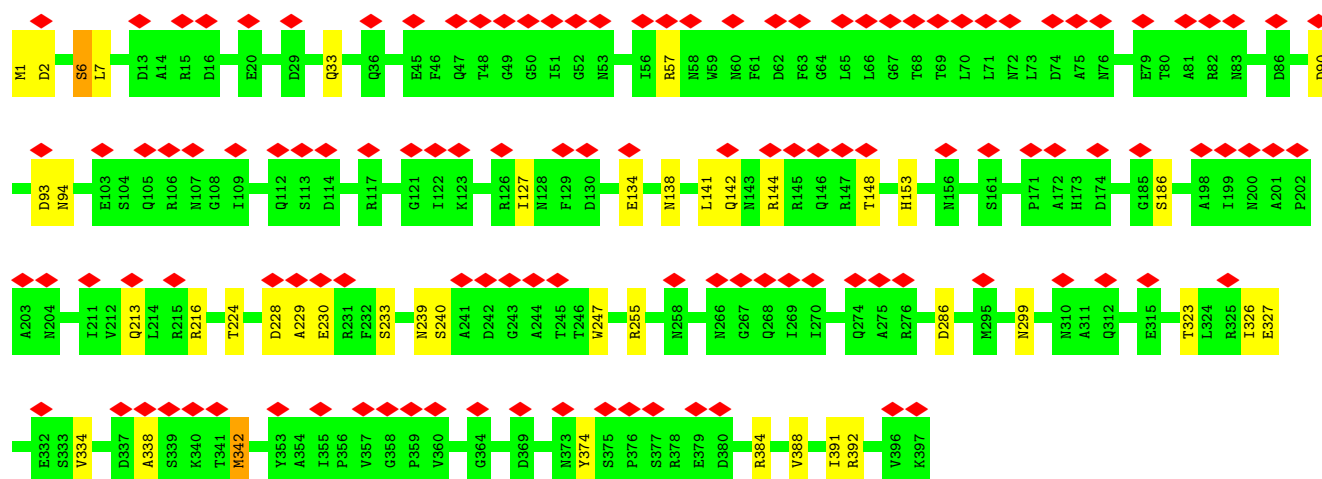


• Molecule 2: Intermediate capsid protein VP6

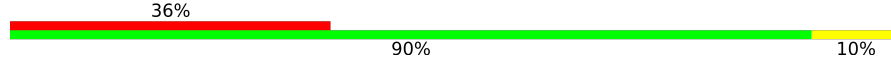


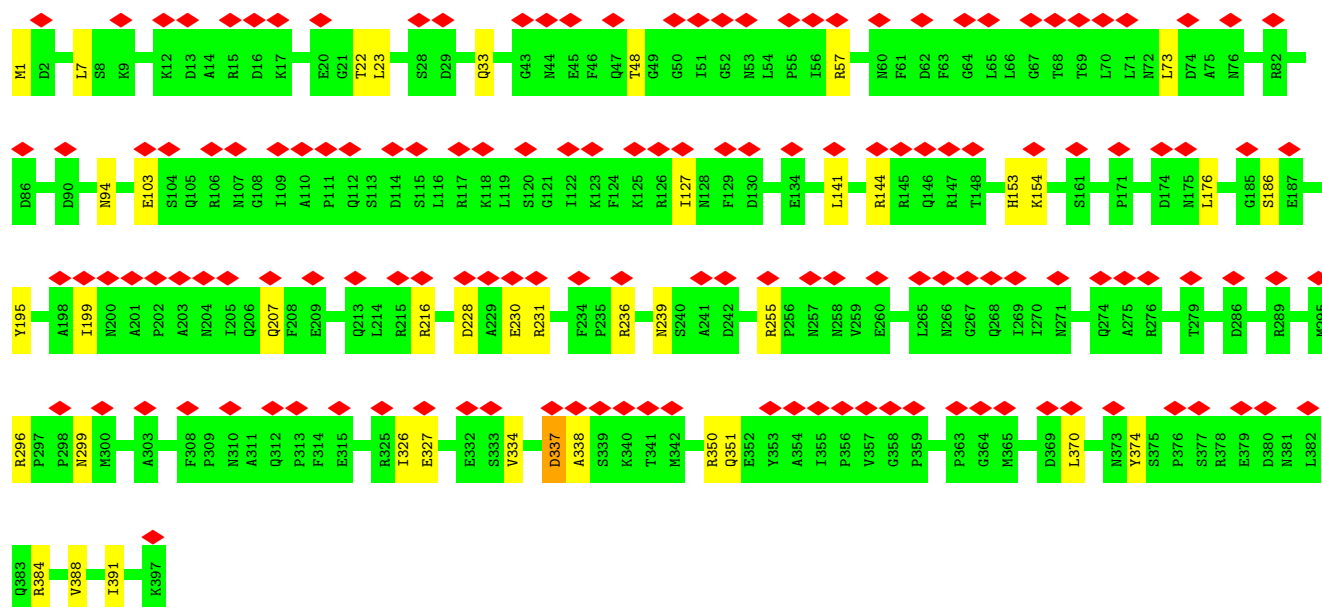
• Molecule 2: Intermediate capsid protein VP6

Chain I: 

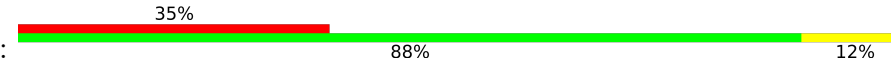


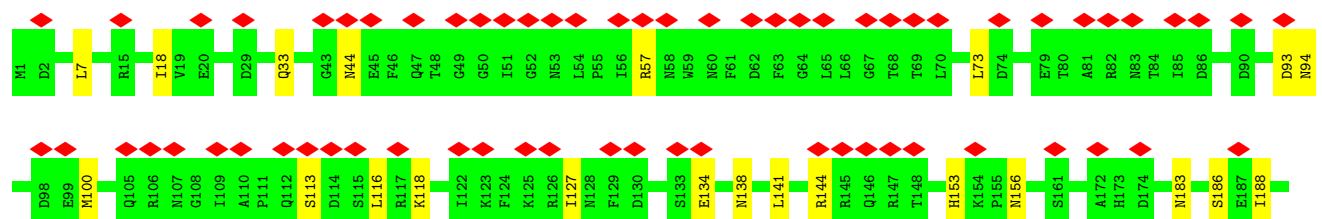
• Molecule 2: Intermediate capsid protein VP6

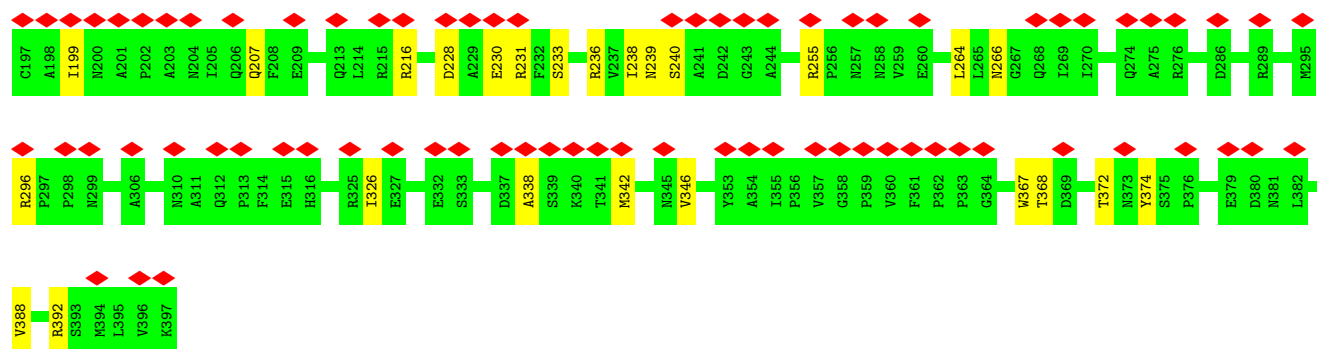
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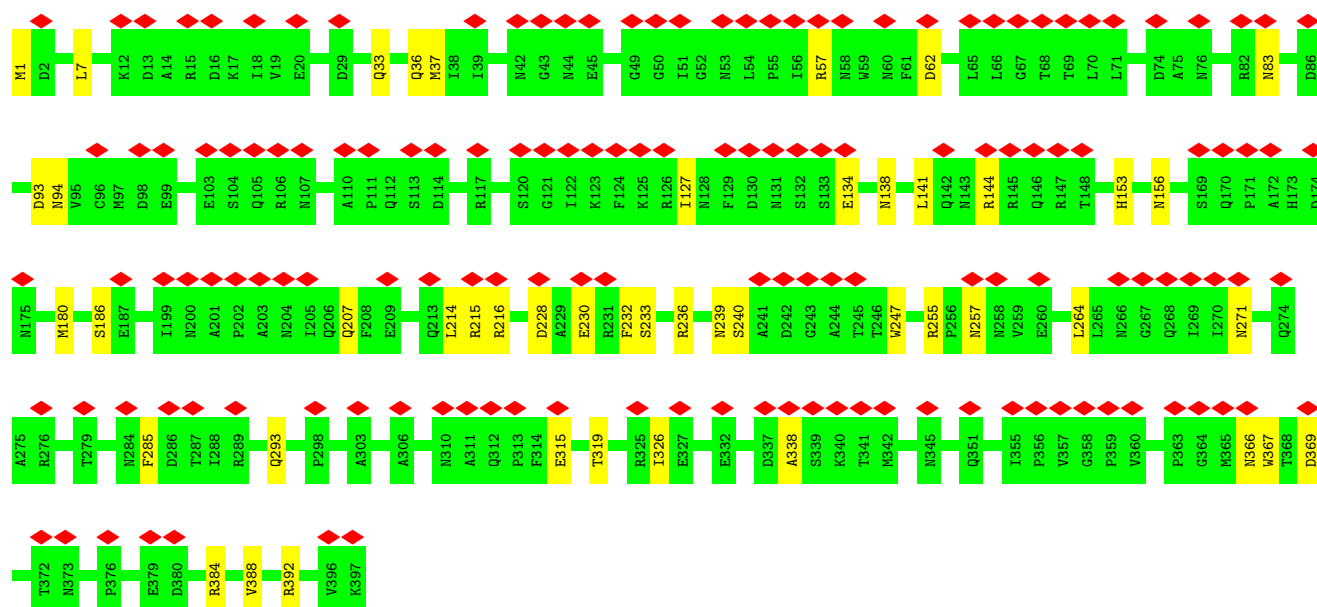
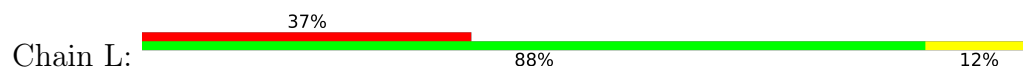
• Molecule 2: Intermediate capsid protein VP6

Chain K: 

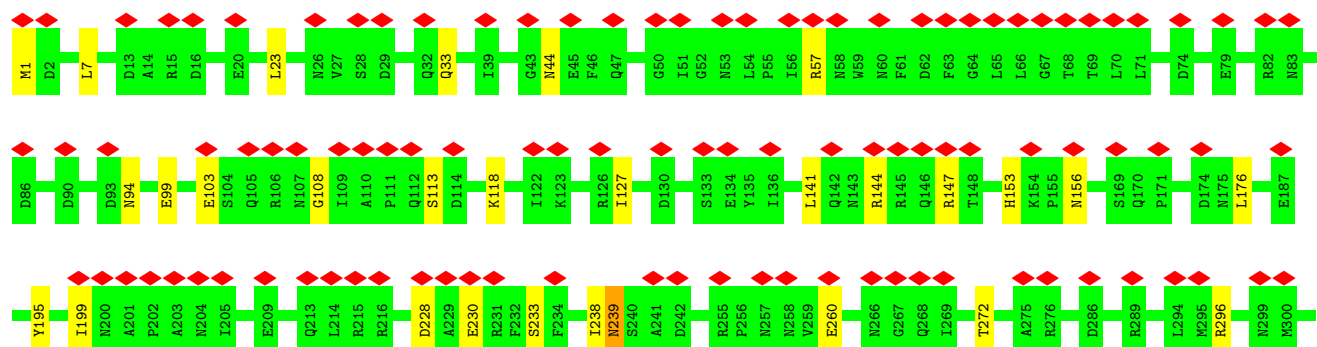
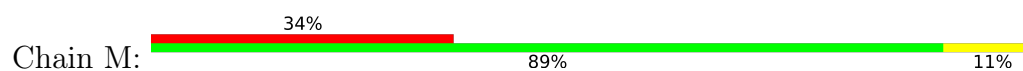


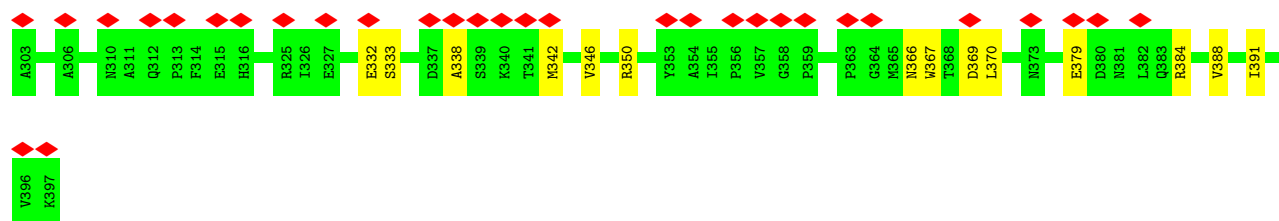


• Molecule 2: Intermediate capsid protein VP6

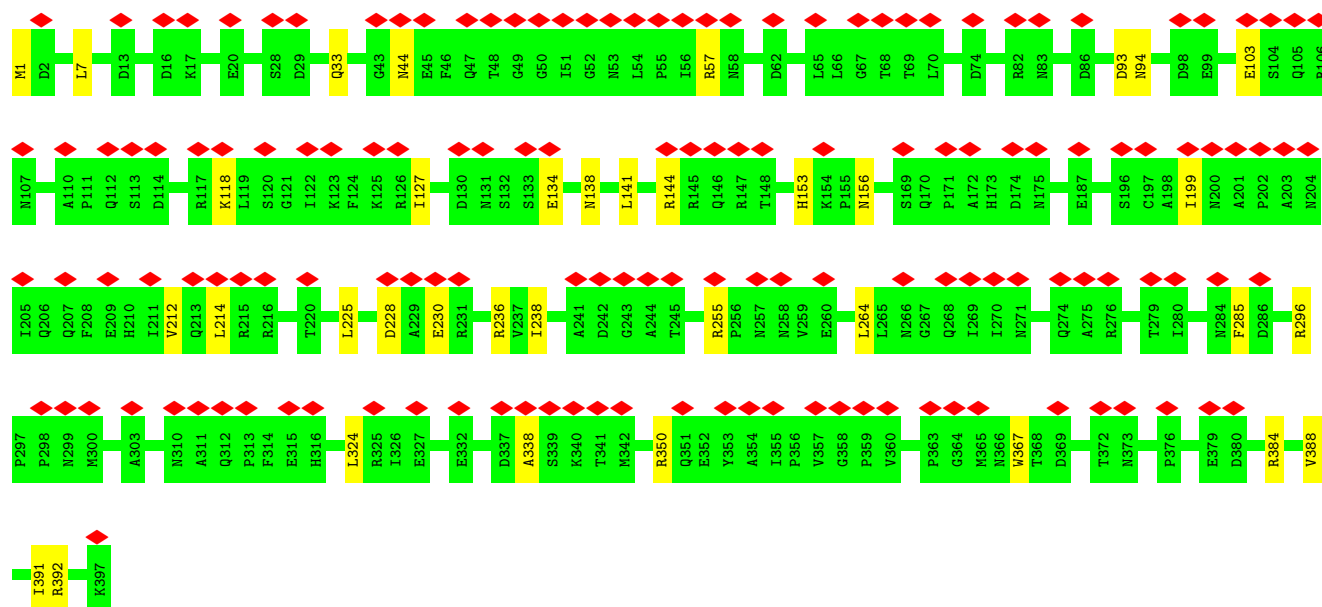
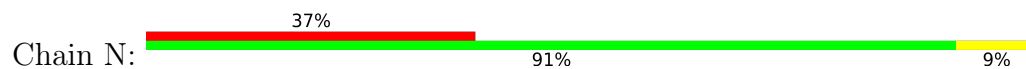


• Molecule 2: Intermediate capsid protein VP6

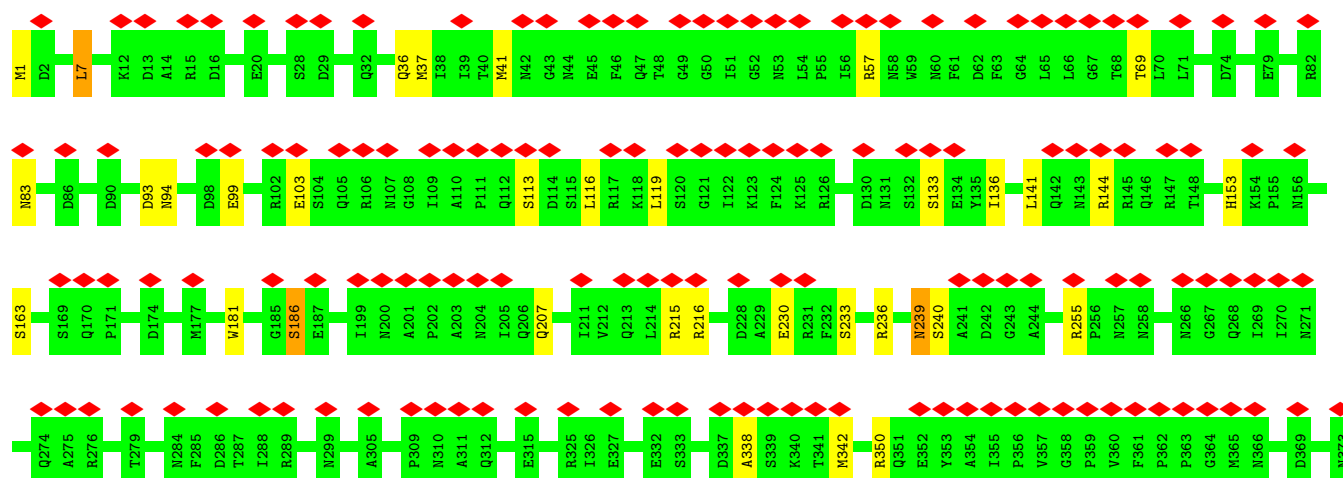
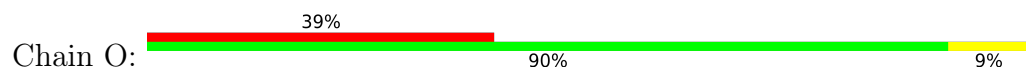


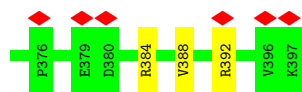


• Molecule 2: Intermediate capsid protein VP6

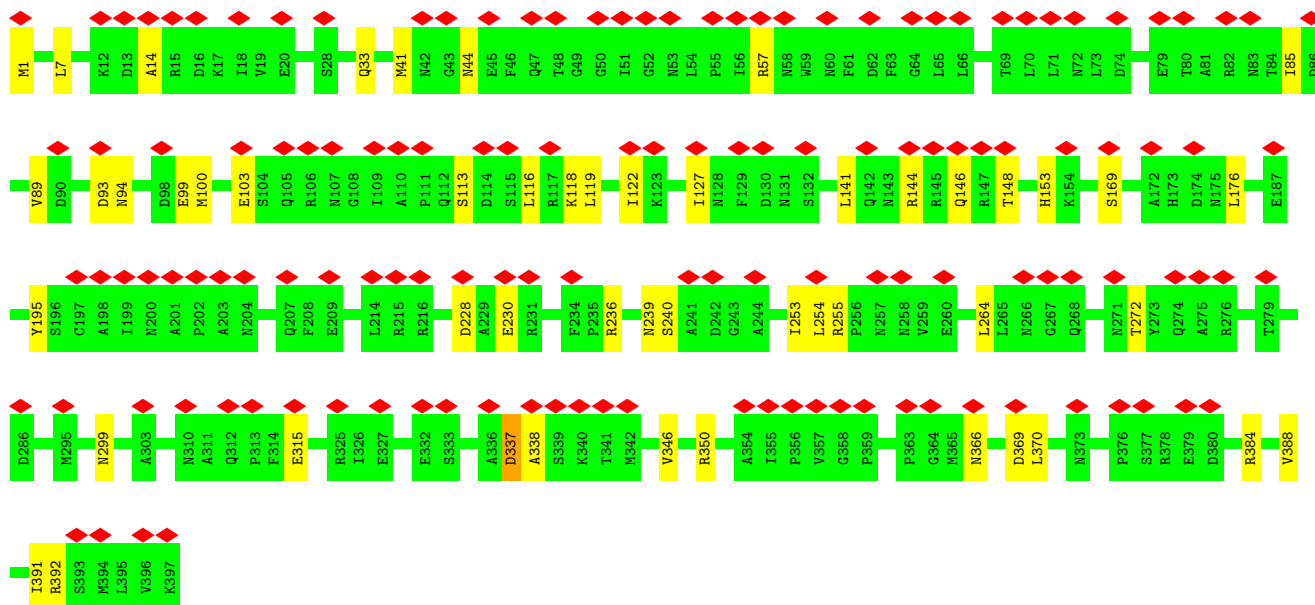
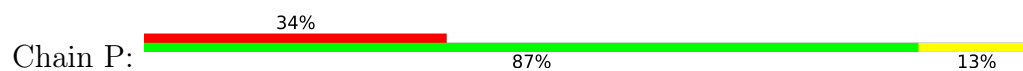


• Molecule 2: Intermediate capsid protein VP6

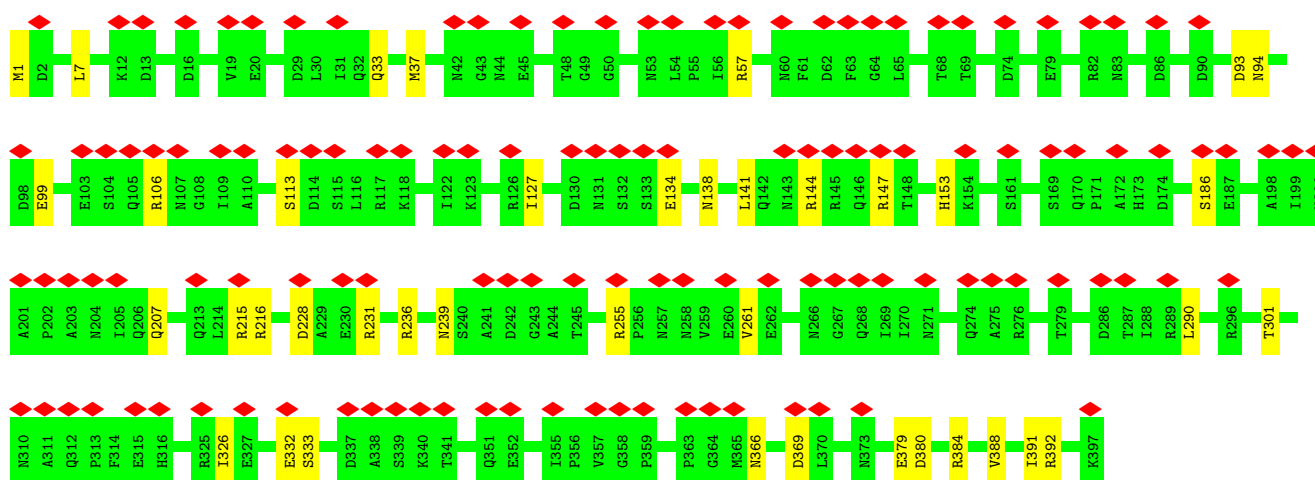
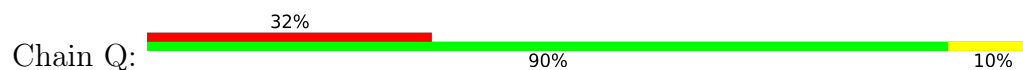




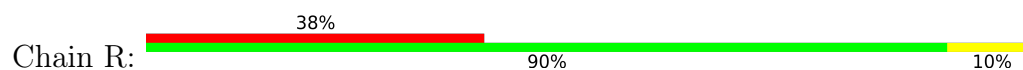
• Molecule 2: Intermediate capsid protein VP6

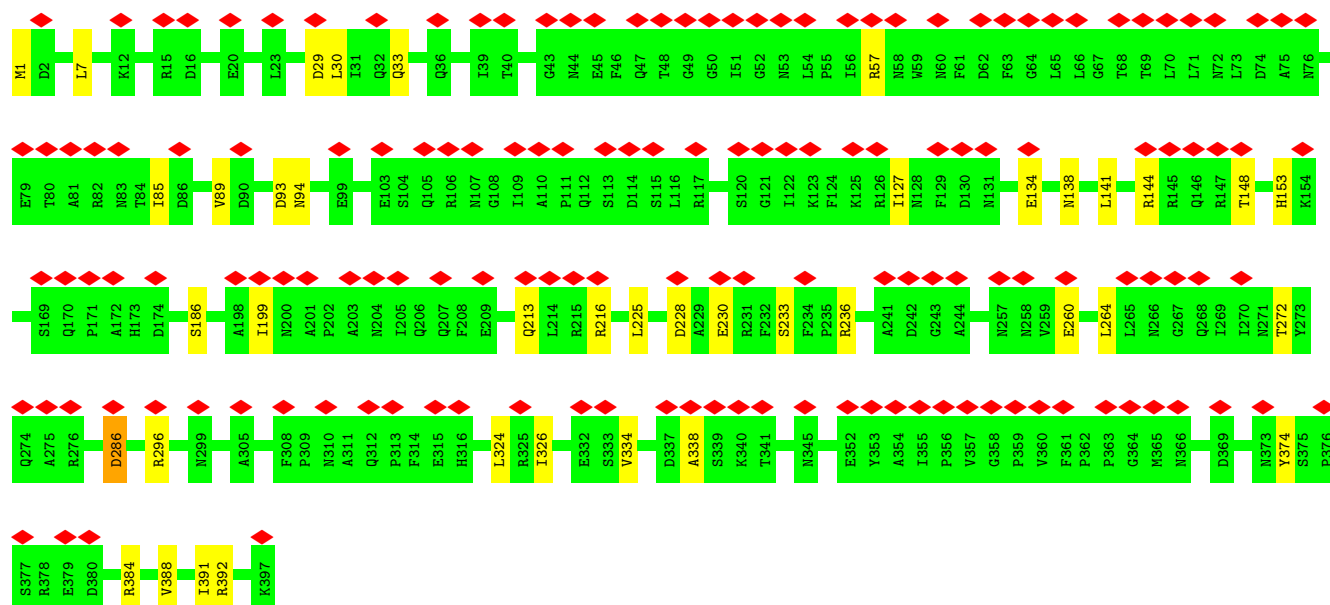


• Molecule 2: Intermediate capsid protein VP6

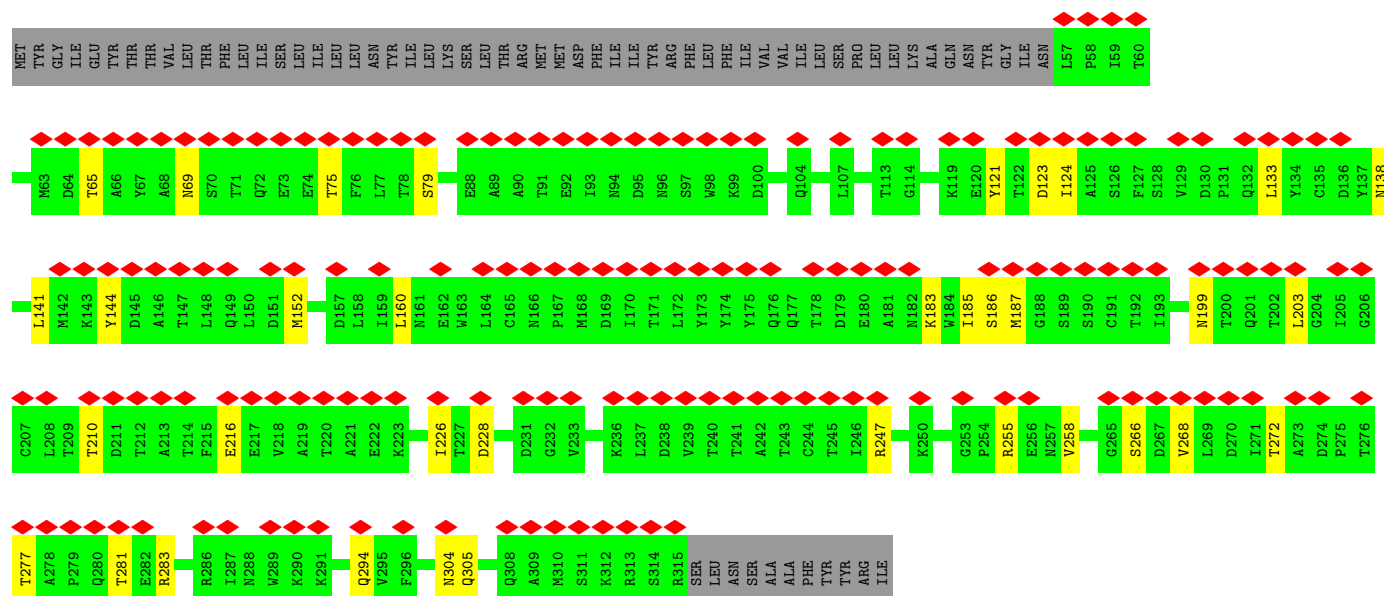


• Molecule 2: Intermediate capsid protein VP6

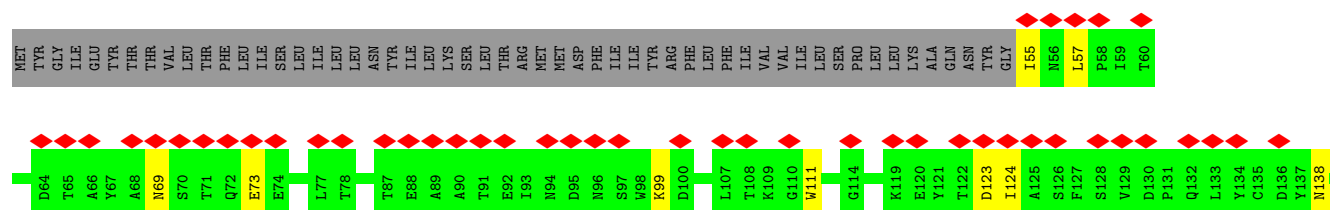


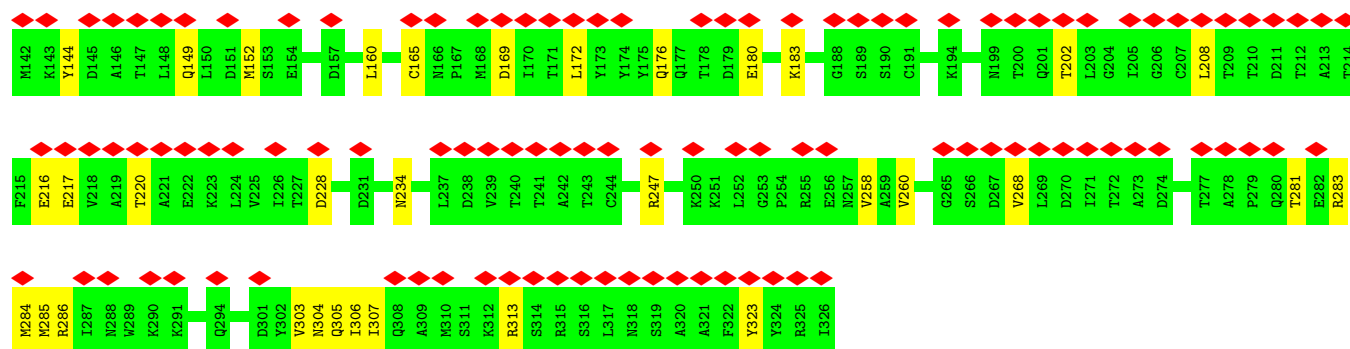


• Molecule 3: Outer capsid glycoprotein VP7

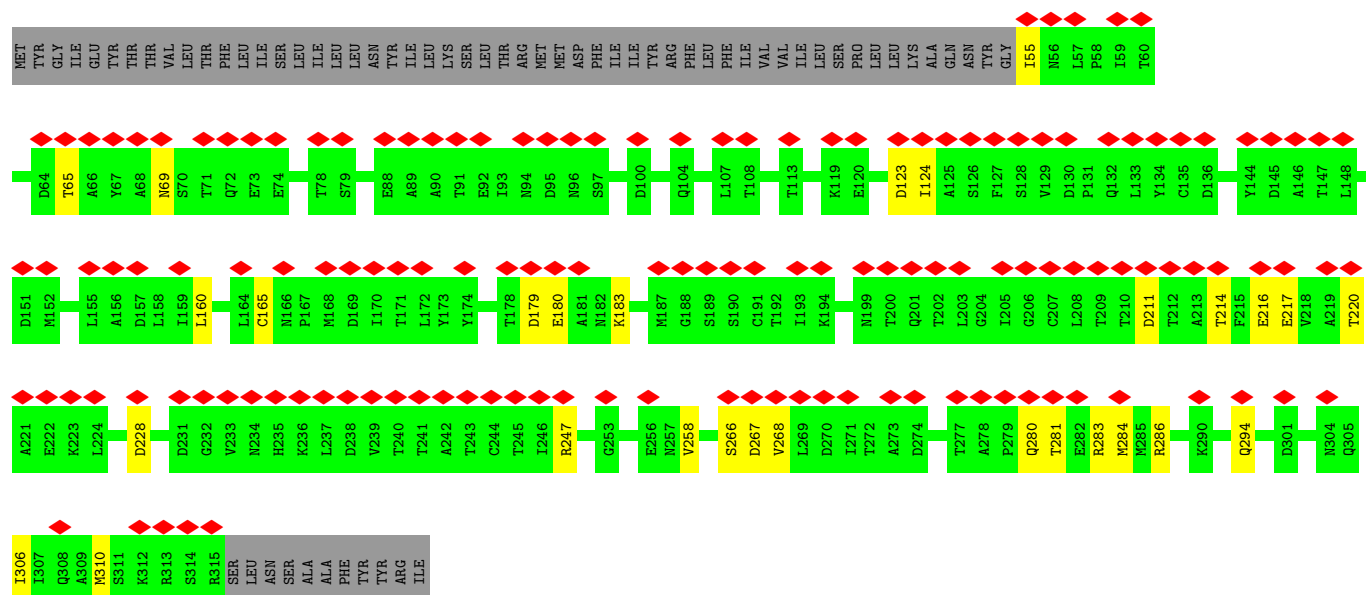
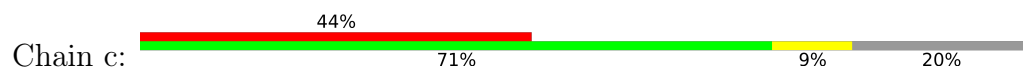


• Molecule 3: Outer capsid glycoprotein VP7

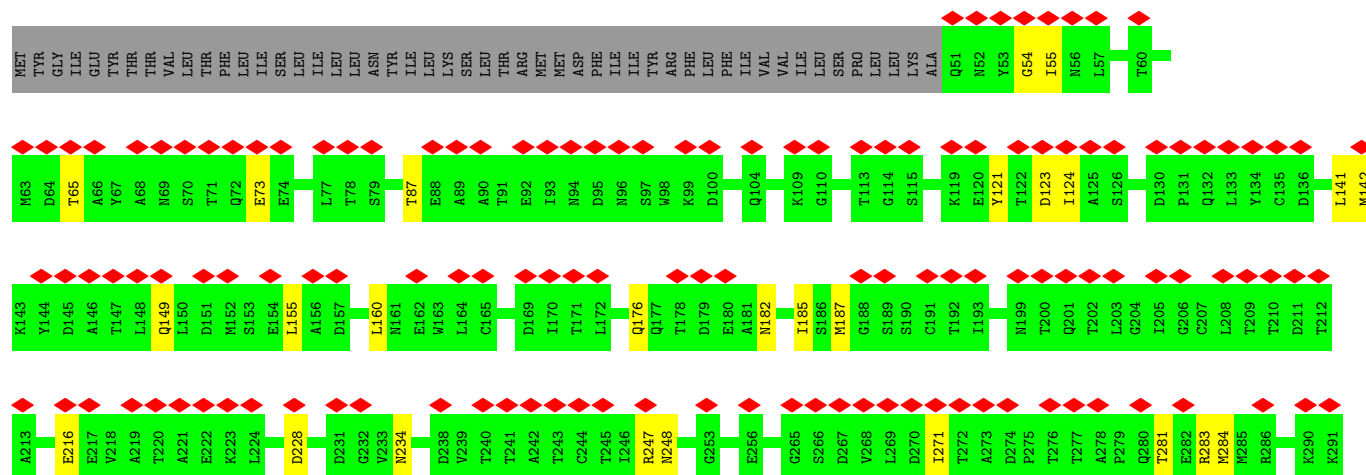
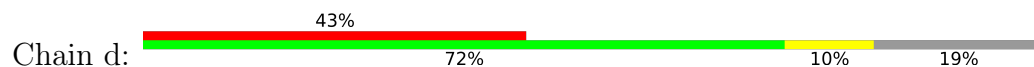


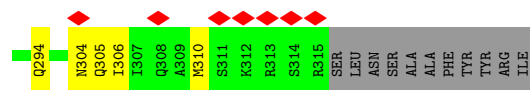


• Molecule 3: Outer capsid glycoprotein VP7

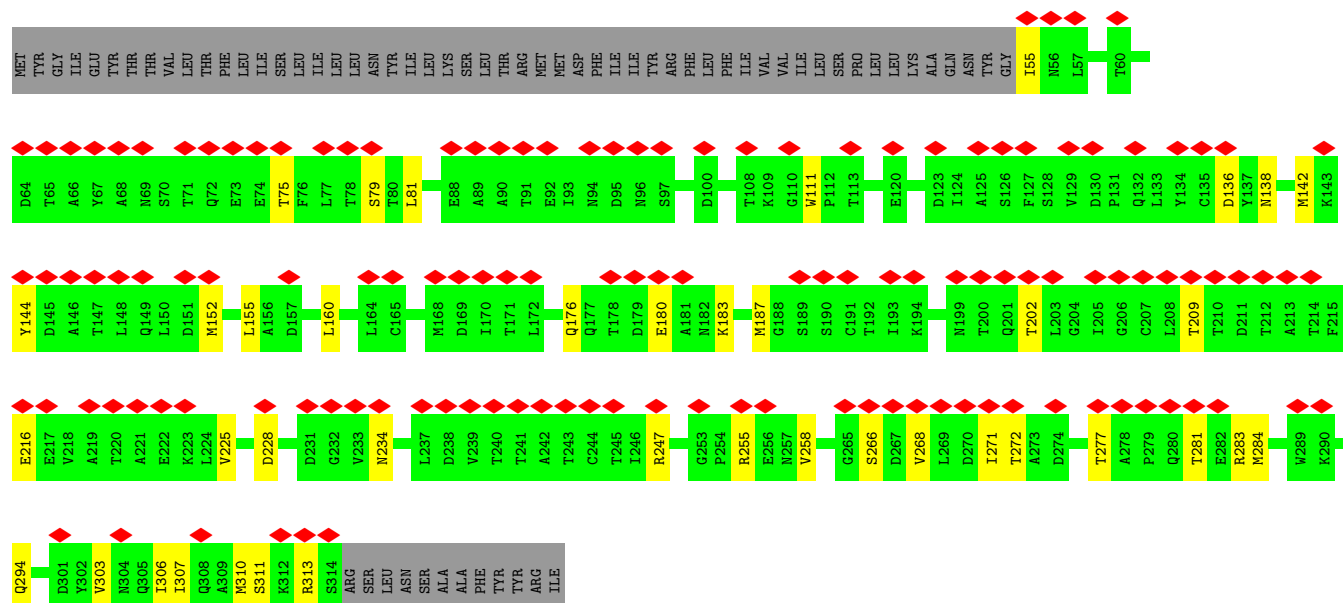
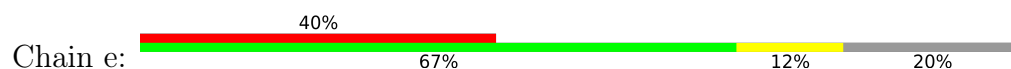


• Molecule 3: Outer capsid glycoprotein VP7

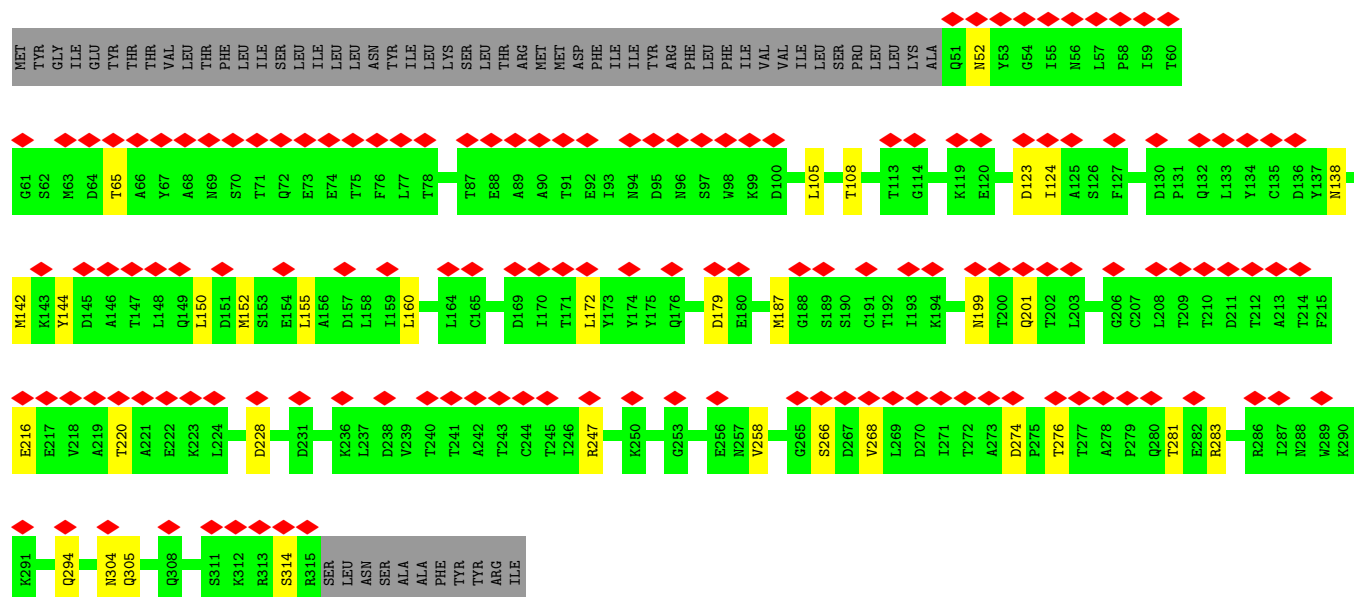
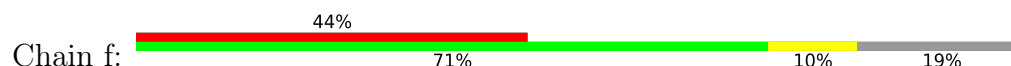




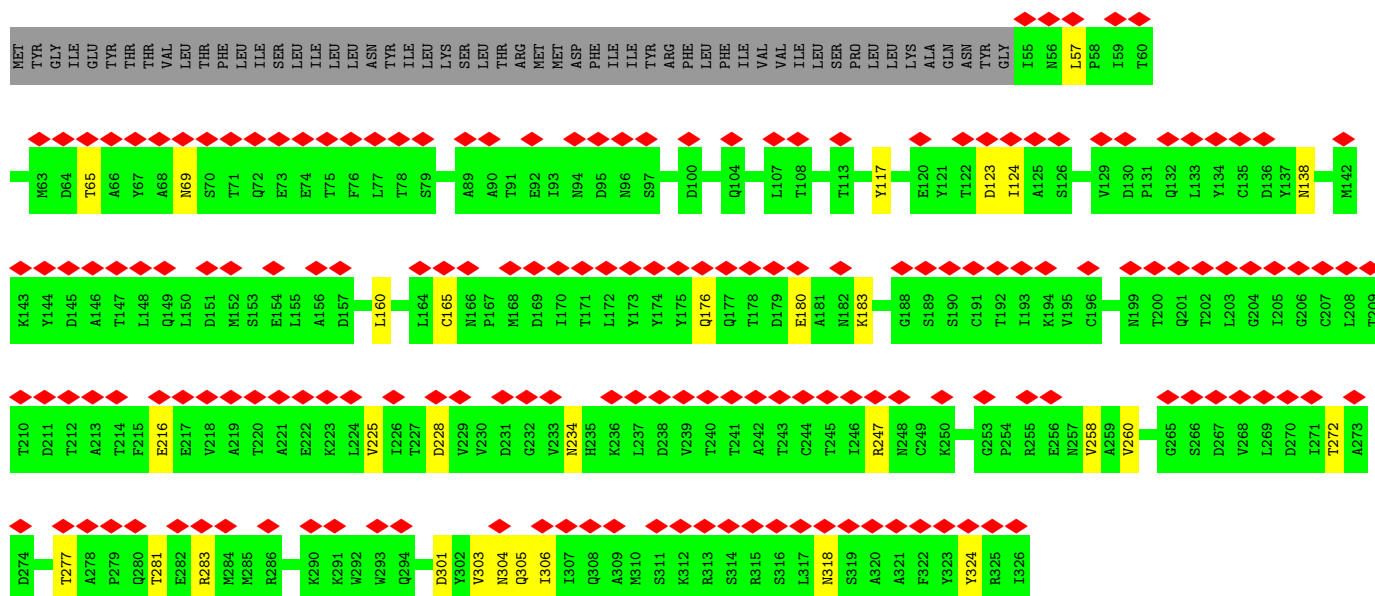
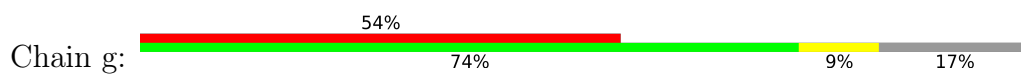
• Molecule 3: Outer capsid glycoprotein VP7



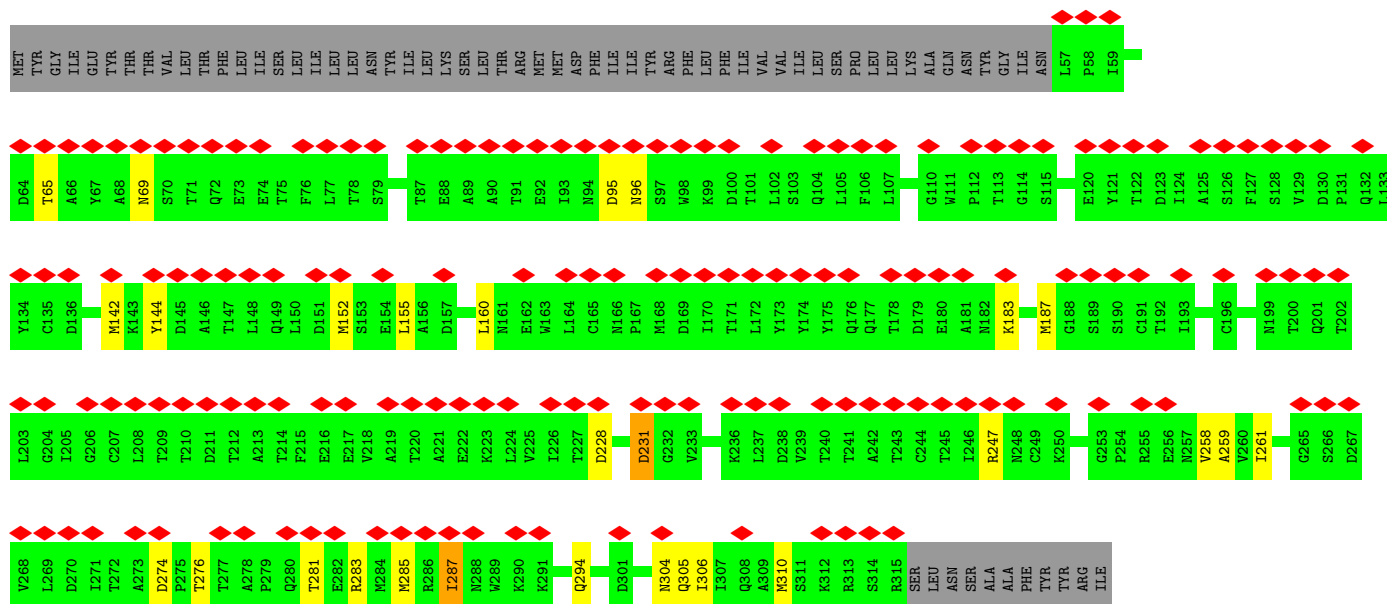
• Molecule 3: Outer capsid glycoprotein VP7



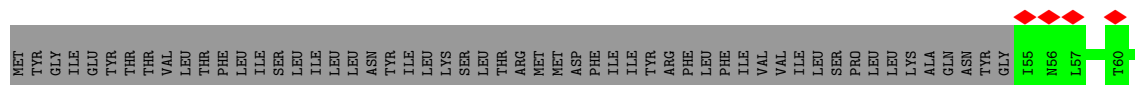
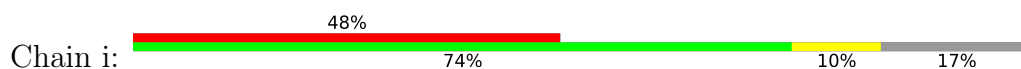
• Molecule 3: Outer capsid glycoprotein VP7

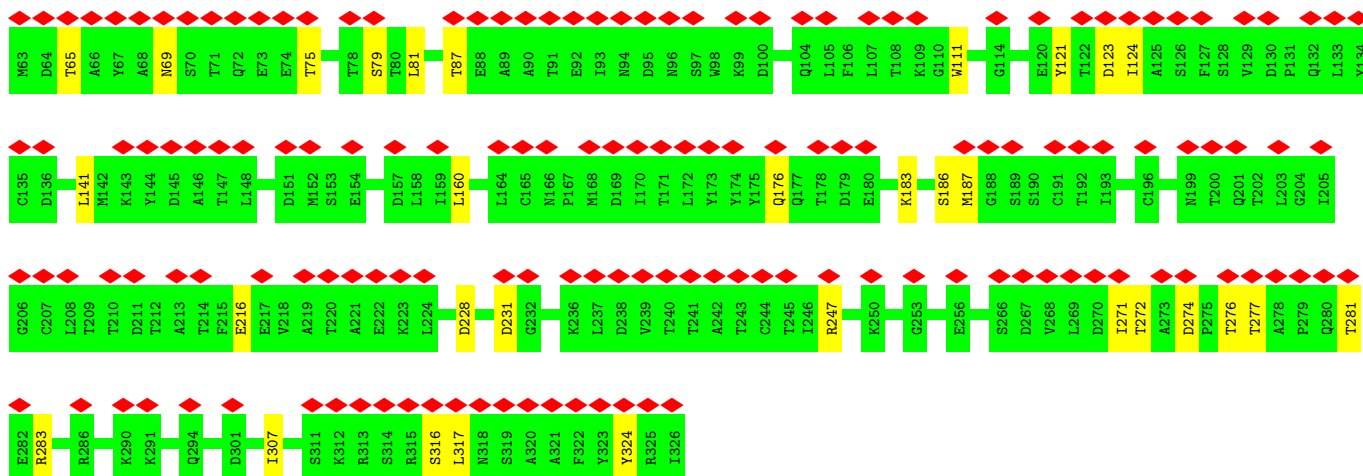


• Molecule 3: Outer capsid glycoprotein VP7

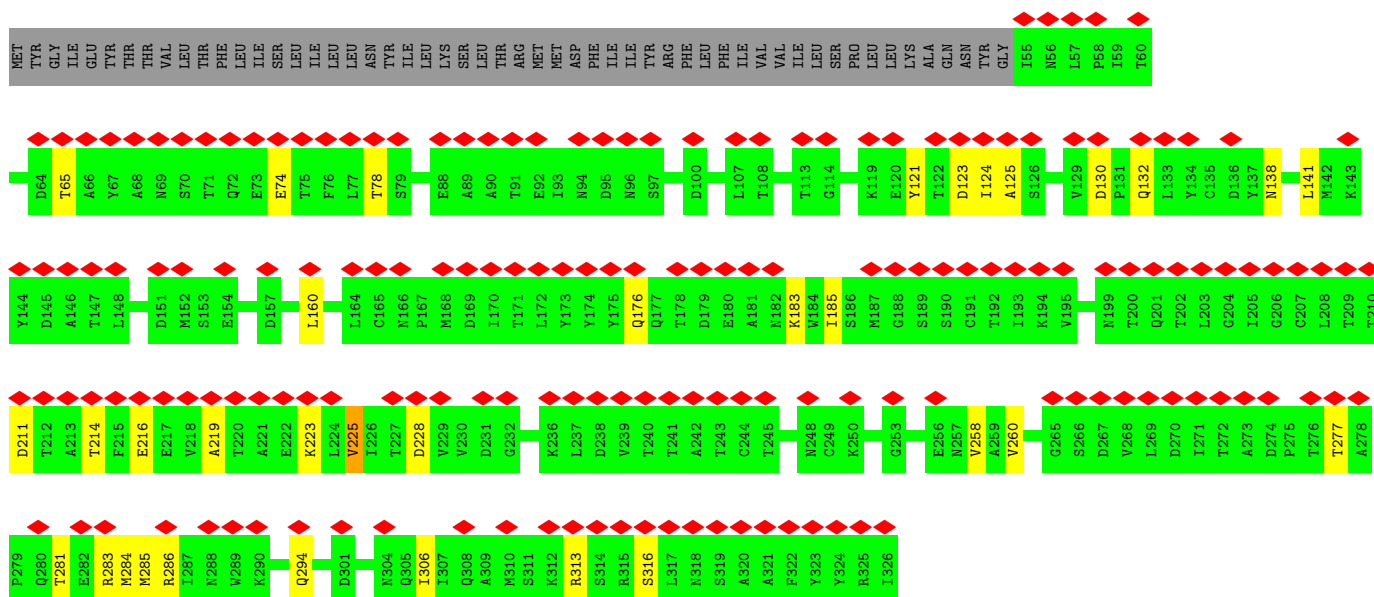
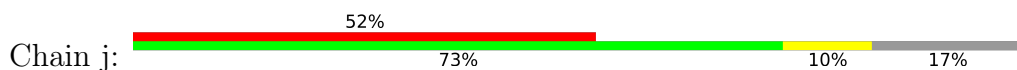


• Molecule 3: Outer capsid glycoprotein VP7

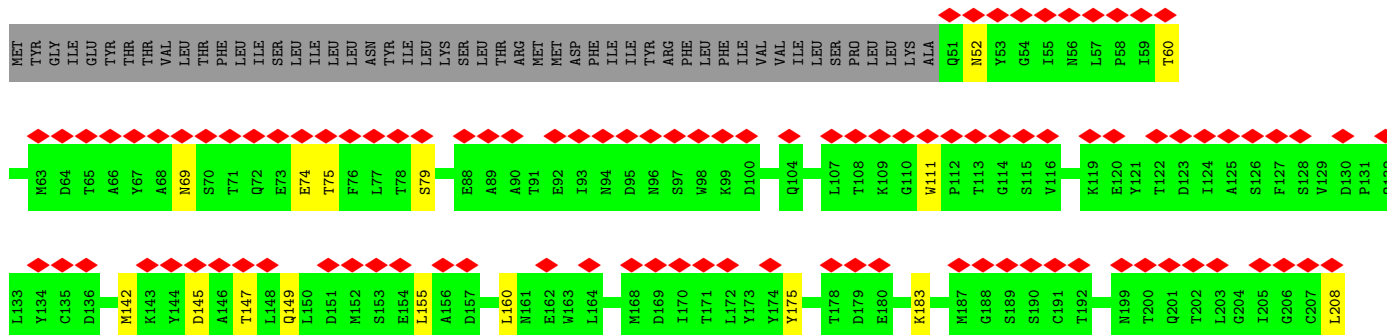
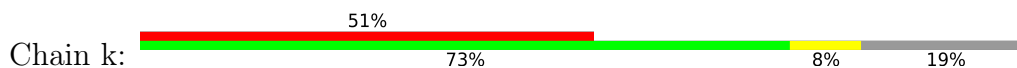


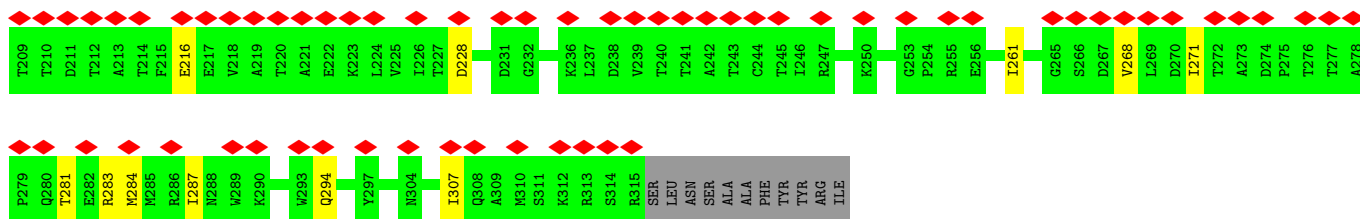


• Molecule 3: Outer capsid glycoprotein VP7

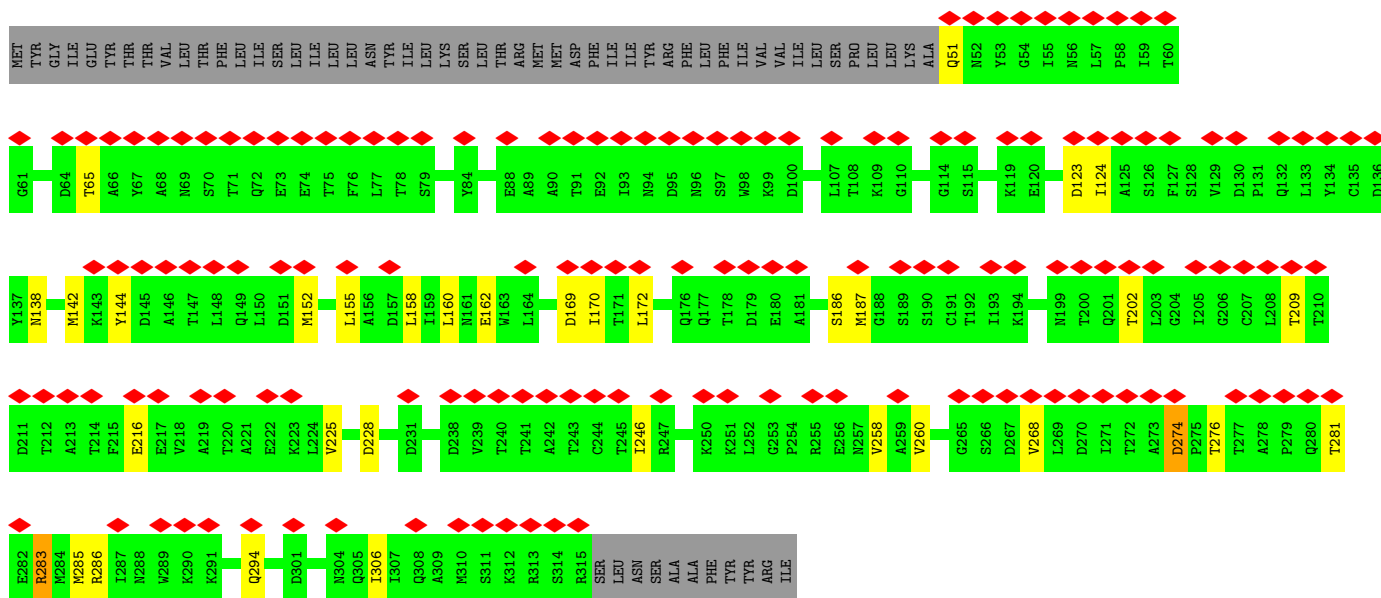


• Molecule 3: Outer capsid glycoprotein VP7

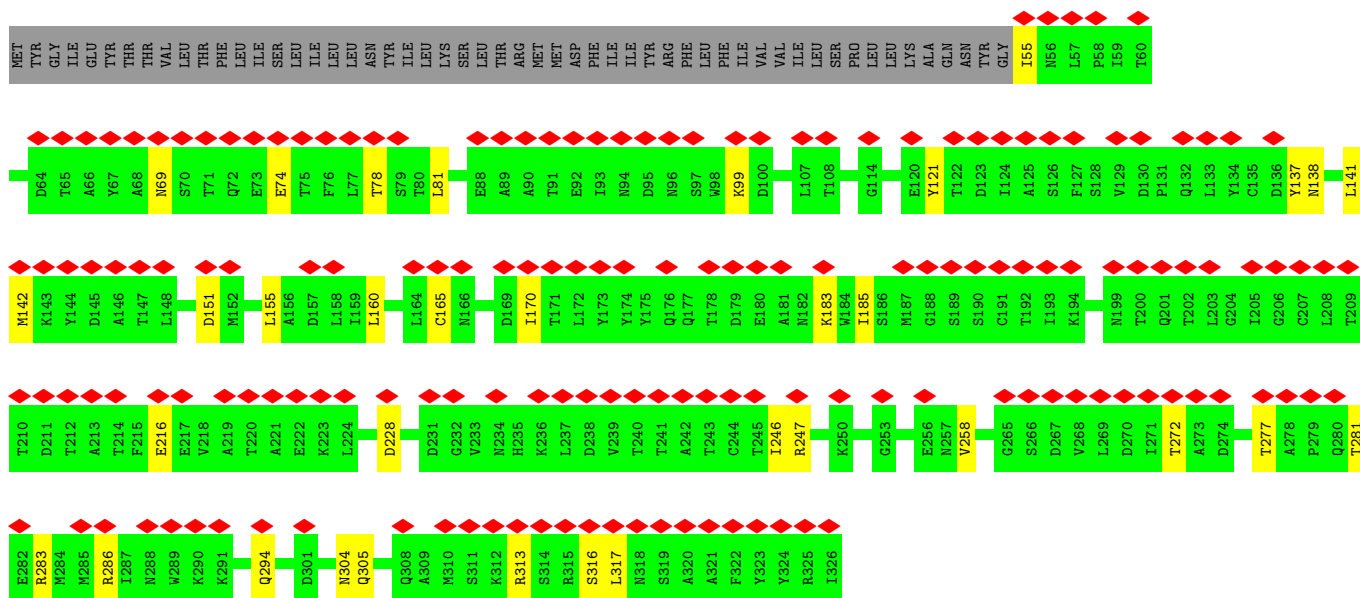
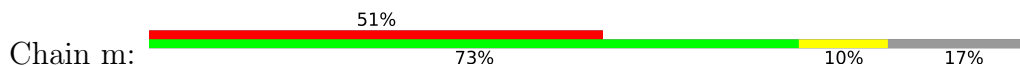




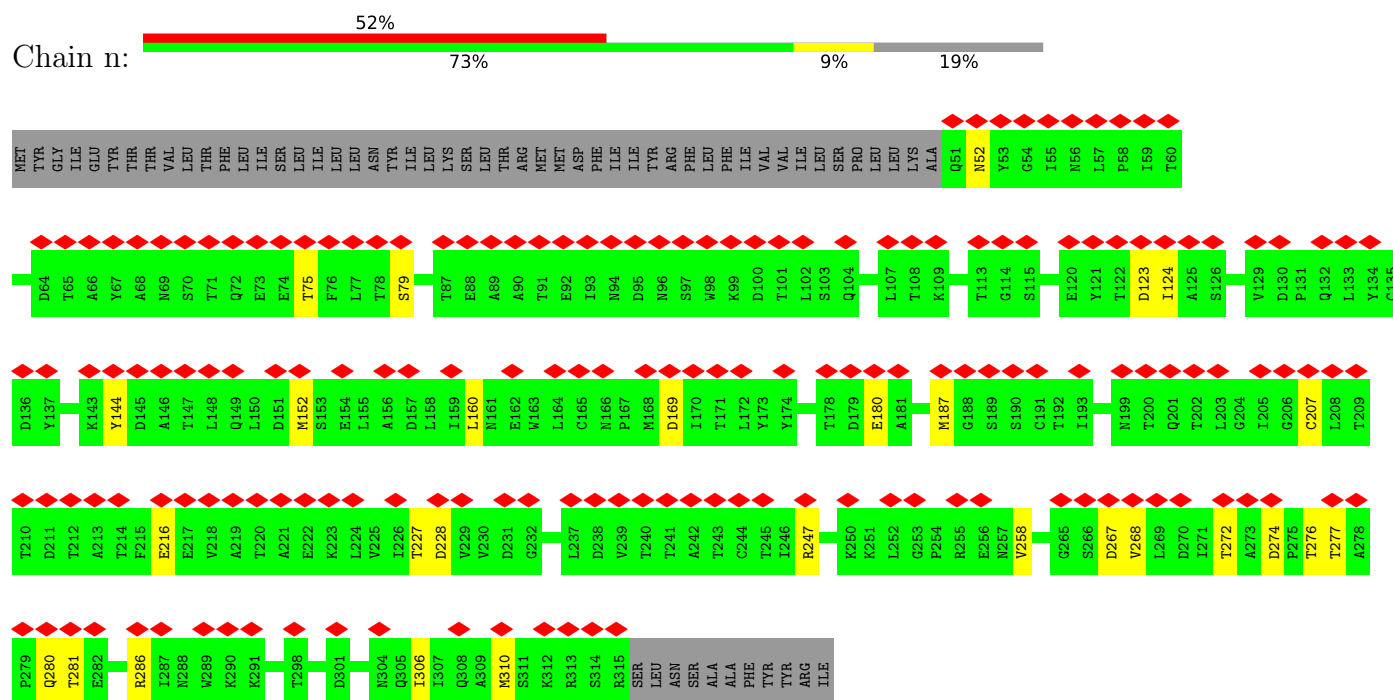
• Molecule 3: Outer capsid glycoprotein VP7



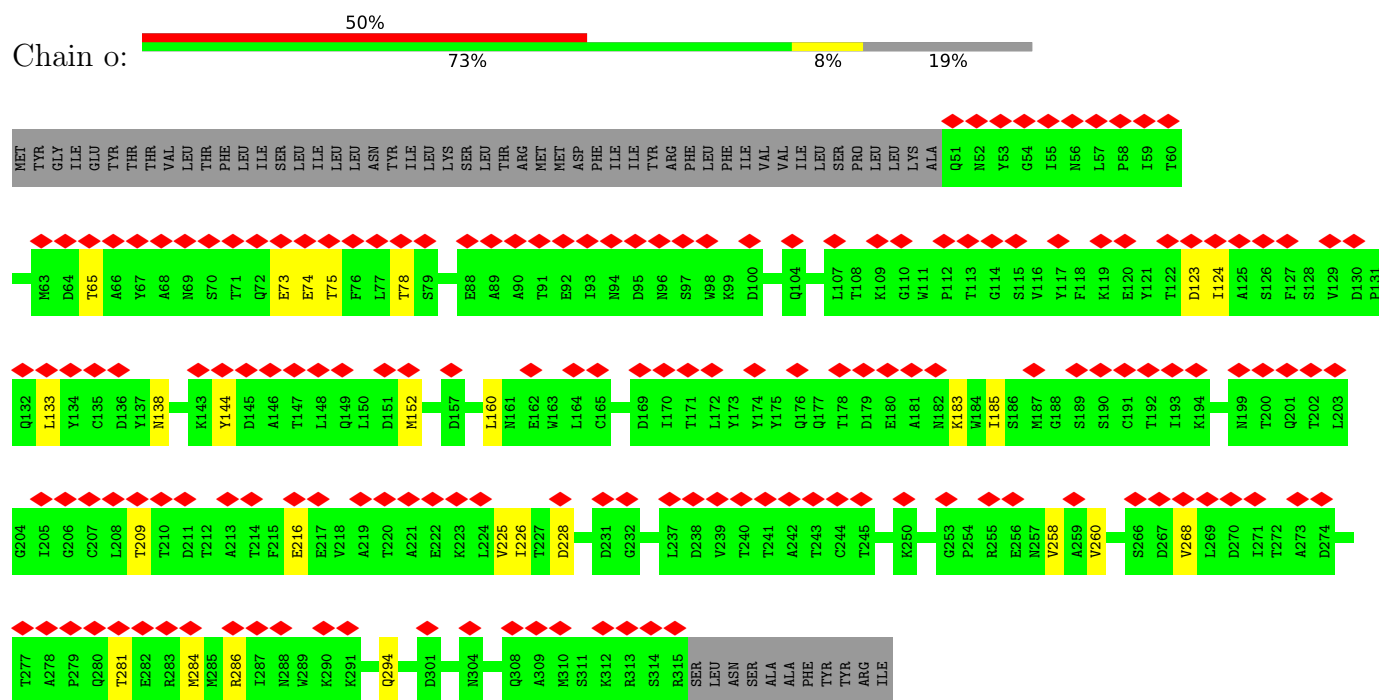
• Molecule 3: Outer capsid glycoprotein VP7



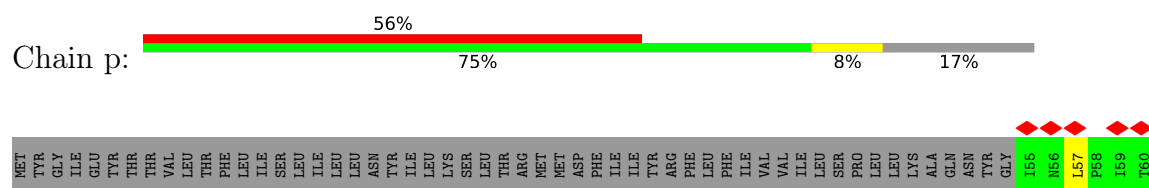
- Molecule 3: Outer capsid glycoprotein VP7

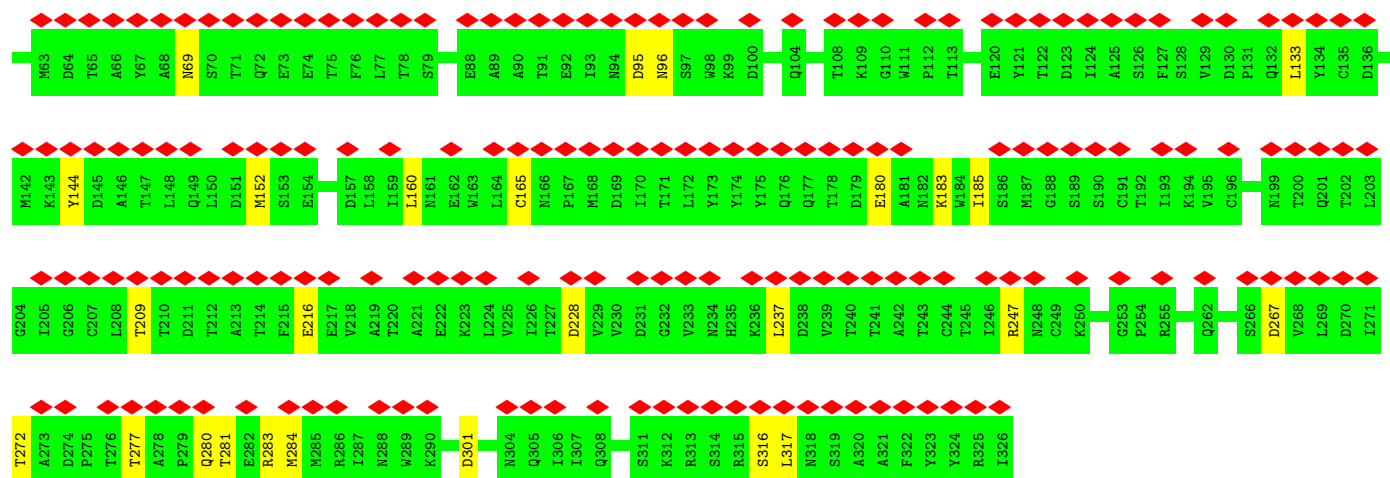


- Molecule 3: Outer capsid glycoprotein VP7

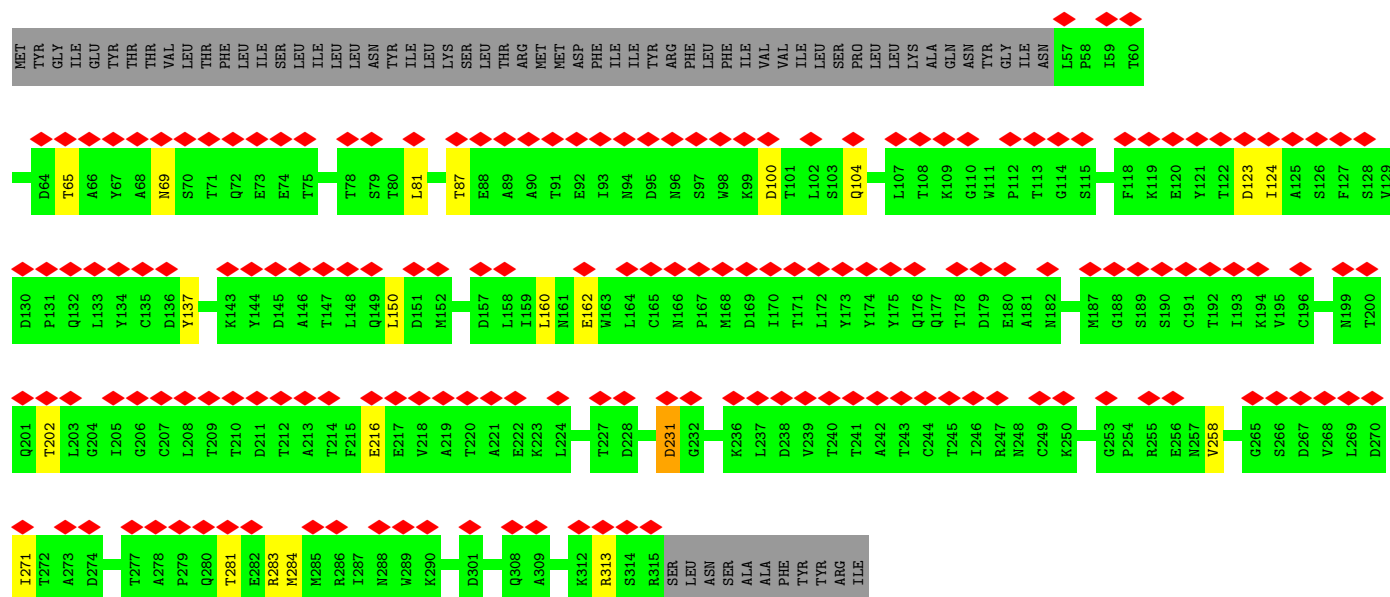
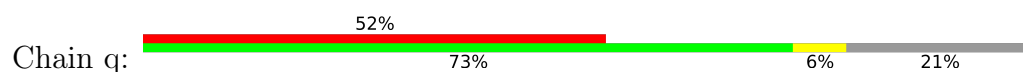


- Molecule 3: Outer capsid glycoprotein VP7

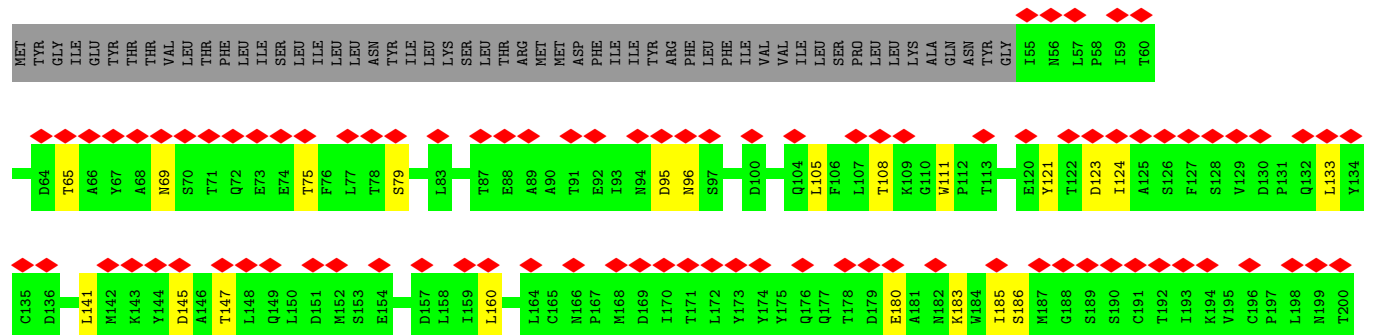
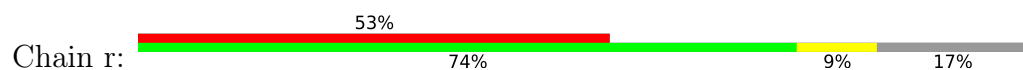


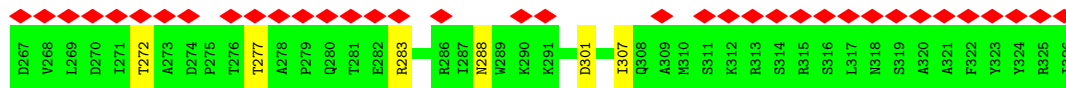


• Molecule 3: Outer capsid glycoprotein VP7



• Molecule 3: Outer capsid glycoprotein VP7





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	252548	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI POLARA 300	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	33	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	3000	Depositor
Magnification	40605	Depositor
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.157	Depositor
Minimum map value	-0.100	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.010	Depositor
Recommended contour level	0.05	Depositor
Map size (Å)	393.91998, 393.91998, 393.91998	wwPDB
Map dimensions	320, 320, 320	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.231, 1.231, 1.231	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: CA, NAG

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	1	0.19	0/2232	0.40	0/3032
1	2	0.19	0/2232	0.41	0/3032
1	3	0.21	0/2232	0.42	0/3032
2	A	0.21	0/3233	0.48	0/4397
2	B	0.21	0/3233	0.47	0/4397
2	C	0.20	0/3233	0.46	0/4397
2	D	0.21	0/3233	0.47	0/4397
2	E	0.21	0/3233	0.47	0/4397
2	F	0.22	0/3233	0.48	0/4397
2	G	0.21	0/3233	0.48	0/4397
2	H	0.22	0/3233	0.46	0/4397
2	I	0.21	0/3233	0.47	0/4397
2	J	0.21	0/3233	0.47	0/4397
2	K	0.21	0/3233	0.46	0/4397
2	L	0.21	0/3233	0.46	0/4397
2	M	0.21	0/3233	0.47	0/4397
2	N	0.20	0/3233	0.45	0/4397
2	O	0.21	0/3233	0.47	0/4397
2	P	0.21	0/3233	0.46	0/4397
2	Q	0.21	0/3233	0.48	0/4397
2	R	0.21	0/3233	0.47	0/4397
3	a	0.19	0/2089	0.43	0/2854
3	b	0.18	0/2200	0.44	0/3005
3	c	0.19	0/2105	0.44	0/2876
3	d	0.20	0/2139	0.44	0/2922
3	e	0.20	0/2094	0.45	0/2862
3	f	0.18	0/2139	0.43	0/2922
3	g	0.19	0/2200	0.45	0/3005
3	h	0.19	0/2089	0.42	0/2854
3	i	0.20	0/2200	0.44	0/3005
3	j	0.19	0/2200	0.45	0/3005
3	k	0.19	0/2139	0.45	0/2922

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
3	l	0.19	0/2139	0.44	0/2922
3	m	0.19	0/2200	0.45	0/3005
3	n	0.18	0/2139	0.43	0/2922
3	o	0.19	0/2139	0.44	0/2922
3	p	0.19	0/2200	0.44	0/3005
3	q	0.18	0/2089	0.43	0/2854
3	r	0.18	0/2200	0.44	0/3005
All	All	0.20	0/103590	0.45	0/141109

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
3	e	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	e	311	SER	Peptide

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	1	2184	2107	2107	21	0
1	2	2184	2107	2107	14	0
1	3	2184	2107	2107	20	0
2	A	3163	3112	3112	32	0
2	B	3163	3112	3112	25	0
2	C	3163	3112	3112	25	0
2	D	3163	3112	3112	29	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	E	3163	3113	3112	33	0
2	F	3163	3113	3112	31	0
2	G	3163	3113	3112	30	0
2	H	3163	3113	3112	24	0
2	I	3163	3112	3112	27	0
2	J	3163	3112	3112	29	0
2	K	3163	3112	3112	28	0
2	L	3163	3113	3112	29	0
2	M	3163	3113	3112	27	0
2	N	3163	3112	3112	22	0
2	O	3163	3113	3112	25	0
2	P	3163	3112	3112	34	0
2	Q	3163	3113	3112	25	0
2	R	3163	3112	3112	25	0
3	a	2047	1994	1993	21	0
3	b	2155	2099	2099	27	0
3	c	2063	2011	2010	19	0
3	d	2096	2037	2036	15	0
3	e	2052	1998	1997	21	0
3	f	2096	2037	2036	21	0
3	g	2155	2099	2098	17	0
3	h	2047	1994	1993	17	0
3	i	2155	2099	2098	16	0
3	j	2155	2099	2098	18	0
3	k	2096	2037	2036	17	0
3	l	2096	2037	2036	22	0
3	m	2155	2099	2098	21	0
3	n	2096	2037	2036	17	0
3	o	2096	2037	2036	15	0
3	p	2155	2099	2098	14	0
3	q	2047	1994	1993	13	0
3	r	2155	2099	2098	18	0
4	a	14	14	13	1	0
4	b	14	14	13	0	0
4	c	14	14	13	1	0
4	d	14	14	13	0	0
4	e	14	14	13	0	0
4	f	14	14	13	0	0
4	g	14	14	13	0	0
4	h	14	14	13	1	0
4	i	14	14	13	1	0
4	j	14	14	13	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	k	14	14	13	1	0
4	l	14	14	13	0	0
4	m	14	14	13	1	0
4	n	14	14	13	0	0
4	o	14	14	13	0	0
4	p	14	14	13	1	0
4	q	14	14	13	0	0
4	r	14	14	13	1	0
5	a	5	0	0	0	0
5	b	3	0	0	0	0
5	c	1	0	0	0	0
5	d	5	0	0	0	0
5	e	3	0	0	0	0
5	f	1	0	0	0	0
5	g	5	0	0	0	0
5	h	3	0	0	0	0
5	i	1	0	0	0	0
5	j	5	0	0	0	0
5	k	3	0	0	0	0
5	l	1	0	0	0	0
5	m	5	0	0	0	0
5	n	3	0	0	0	0
5	o	1	0	0	0	0
5	p	5	0	0	0	0
5	q	3	0	0	0	0
5	r	1	0	0	0	0
All	All	101709	99503	99460	769	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 4.

The worst 5 of 769 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)
1:1:485:THR:HG1	1:2:412:SER:HG	1.26	0.80
2:G:255:ARG:NH1	3:h:65:THR:O	2.18	0.77
2:E:255:ARG:NH1	3:f:65:THR:O	2.19	0.76
2:Q:255:ARG:NH1	3:r:65:THR:O	2.19	0.75
2:H:255:ARG:NH1	3:i:65:THR:O	2.19	0.74

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	1	273/776 (35%)	273 (100%)	0	0	100	100
1	2	273/776 (35%)	273 (100%)	0	0	100	100
1	3	273/776 (35%)	273 (100%)	0	0	100	100
2	A	395/397 (100%)	392 (99%)	3 (1%)	0	100	100
2	B	395/397 (100%)	392 (99%)	3 (1%)	0	100	100
2	C	395/397 (100%)	393 (100%)	2 (0%)	0	100	100
2	D	395/397 (100%)	392 (99%)	3 (1%)	0	100	100
2	E	395/397 (100%)	392 (99%)	3 (1%)	0	100	100
2	F	395/397 (100%)	391 (99%)	4 (1%)	0	100	100
2	G	395/397 (100%)	393 (100%)	2 (0%)	0	100	100
2	H	395/397 (100%)	392 (99%)	3 (1%)	0	100	100
2	I	395/397 (100%)	392 (99%)	3 (1%)	0	100	100
2	J	395/397 (100%)	392 (99%)	3 (1%)	0	100	100
2	K	395/397 (100%)	391 (99%)	4 (1%)	0	100	100
2	L	395/397 (100%)	392 (99%)	3 (1%)	0	100	100
2	M	395/397 (100%)	392 (99%)	3 (1%)	0	100	100
2	N	395/397 (100%)	392 (99%)	3 (1%)	0	100	100
2	O	395/397 (100%)	393 (100%)	2 (0%)	0	100	100
2	P	395/397 (100%)	392 (99%)	3 (1%)	0	100	100
2	Q	395/397 (100%)	393 (100%)	2 (0%)	0	100	100
2	R	395/397 (100%)	393 (100%)	2 (0%)	0	100	100
3	a	257/326 (79%)	257 (100%)	0	0	100	100
3	b	270/326 (83%)	269 (100%)	1 (0%)	0	100	100
3	c	259/326 (79%)	258 (100%)	1 (0%)	0	100	100
3	d	263/326 (81%)	263 (100%)	0	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	e	258/326 (79%)	257 (100%)	1 (0%)	0	100	100
3	f	263/326 (81%)	263 (100%)	0	0	100	100
3	g	270/326 (83%)	269 (100%)	1 (0%)	0	100	100
3	h	257/326 (79%)	257 (100%)	0	0	100	100
3	i	270/326 (83%)	270 (100%)	0	0	100	100
3	j	270/326 (83%)	268 (99%)	2 (1%)	0	100	100
3	k	263/326 (81%)	263 (100%)	0	0	100	100
3	l	263/326 (81%)	263 (100%)	0	0	100	100
3	m	270/326 (83%)	269 (100%)	1 (0%)	0	100	100
3	n	263/326 (81%)	263 (100%)	0	0	100	100
3	o	263/326 (81%)	263 (100%)	0	0	100	100
3	p	270/326 (83%)	270 (100%)	0	0	100	100
3	q	257/326 (79%)	257 (100%)	0	0	100	100
3	r	270/326 (83%)	269 (100%)	1 (0%)	0	100	100
All	All	12685/15342 (83%)	12626 (100%)	59 (0%)	0	100	100

There are no Ramachandran outliers to report.

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	1	241/688 (35%)	234 (97%)	7 (3%)	37	62
1	2	241/688 (35%)	235 (98%)	6 (2%)	42	64
1	3	241/688 (35%)	232 (96%)	9 (4%)	30	58
2	A	350/350 (100%)	345 (99%)	5 (1%)	59	73
2	B	350/350 (100%)	344 (98%)	6 (2%)	53	71
2	C	350/350 (100%)	344 (98%)	6 (2%)	53	71
2	D	350/350 (100%)	344 (98%)	6 (2%)	53	71

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	E	350/350 (100%)	346 (99%)	4 (1%)	65	76
2	F	350/350 (100%)	341 (97%)	9 (3%)	40	64
2	G	350/350 (100%)	346 (99%)	4 (1%)	65	76
2	H	350/350 (100%)	343 (98%)	7 (2%)	48	68
2	I	350/350 (100%)	346 (99%)	4 (1%)	65	76
2	J	350/350 (100%)	346 (99%)	4 (1%)	65	76
2	K	350/350 (100%)	346 (99%)	4 (1%)	65	76
2	L	350/350 (100%)	345 (99%)	5 (1%)	59	73
2	M	350/350 (100%)	347 (99%)	3 (1%)	70	78
2	N	350/350 (100%)	346 (99%)	4 (1%)	65	76
2	O	350/350 (100%)	344 (98%)	6 (2%)	53	71
2	P	350/350 (100%)	345 (99%)	5 (1%)	59	73
2	Q	350/350 (100%)	347 (99%)	3 (1%)	70	78
2	R	350/350 (100%)	347 (99%)	3 (1%)	70	78
3	a	233/295 (79%)	231 (99%)	2 (1%)	70	78
3	b	244/295 (83%)	238 (98%)	6 (2%)	42	64
3	c	235/295 (80%)	232 (99%)	3 (1%)	61	74
3	d	238/295 (81%)	233 (98%)	5 (2%)	47	67
3	e	234/295 (79%)	229 (98%)	5 (2%)	47	67
3	f	238/295 (81%)	234 (98%)	4 (2%)	53	71
3	g	244/295 (83%)	239 (98%)	5 (2%)	48	68
3	h	233/295 (79%)	230 (99%)	3 (1%)	61	74
3	i	244/295 (83%)	239 (98%)	5 (2%)	48	68
3	j	244/295 (83%)	238 (98%)	6 (2%)	42	64
3	k	238/295 (81%)	233 (98%)	5 (2%)	47	67
3	l	238/295 (81%)	230 (97%)	8 (3%)	32	59
3	m	244/295 (83%)	241 (99%)	3 (1%)	63	75
3	n	238/295 (81%)	235 (99%)	3 (1%)	61	74
3	o	238/295 (81%)	231 (97%)	7 (3%)	37	62
3	p	244/295 (83%)	238 (98%)	6 (2%)	42	64
3	q	233/295 (79%)	229 (98%)	4 (2%)	53	71

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
3	r	244/295 (83%)	239 (98%)	5 (2%)	48 68
All	All	11327/13674 (83%)	11132 (98%)	195 (2%)	52 71

5 of 195 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	b	208	LEU
3	i	87	THR
3	c	281	THR
3	e	281	THR
3	j	225	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 113 such sidechains are listed below:

Mol	Chain	Res	Type
2	M	257	ASN
3	p	235	HIS
2	Q	383	GLN
3	p	166	ASN
3	k	262	GLN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 72 ligands modelled in this entry, 54 are monoatomic - leaving 18 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and

the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	NAG	m	401	3	14,14,15	0.62	1 (7%)	17,19,21	0.73	1 (5%)
4	NAG	e	401	3	14,14,15	0.21	0	17,19,21	0.57	0
4	NAG	a	401	3	14,14,15	0.67	1 (7%)	17,19,21	0.78	1 (5%)
4	NAG	c	401	3	14,14,15	0.62	1 (7%)	17,19,21	0.72	1 (5%)
4	NAG	b	401	-	14,14,15	0.24	0	17,19,21	0.50	0
4	NAG	g	401	3	14,14,15	0.45	0	17,19,21	0.46	0
4	NAG	j	401	3	14,14,15	0.37	0	17,19,21	0.37	0
4	NAG	n	401	3	14,14,15	0.32	0	17,19,21	0.36	0
4	NAG	q	401	3	14,14,15	0.47	0	17,19,21	0.72	1 (5%)
4	NAG	d	401	3	14,14,15	0.39	0	17,19,21	0.40	0
4	NAG	o	401	3	14,14,15	0.50	0	17,19,21	0.47	0
4	NAG	f	401	3	14,14,15	0.38	0	17,19,21	0.45	0
4	NAG	h	401	3	14,14,15	0.64	1 (7%)	17,19,21	0.74	1 (5%)
4	NAG	l	401	3	14,14,15	0.41	0	17,19,21	0.47	0
4	NAG	p	401	3	14,14,15	0.55	0	17,19,21	0.76	1 (5%)
4	NAG	r	401	3	14,14,15	0.65	1 (7%)	17,19,21	0.78	1 (5%)
4	NAG	k	401	3	14,14,15	0.64	1 (7%)	17,19,21	0.81	1 (5%)
4	NAG	i	401	3	14,14,15	0.64	1 (7%)	17,19,21	0.75	1 (5%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	m	401	3	-	0/6/23/26	0/1/1/1
4	NAG	e	401	3	-	0/6/23/26	0/1/1/1
4	NAG	a	401	3	-	0/6/23/26	0/1/1/1
4	NAG	c	401	3	-	2/6/23/26	0/1/1/1
4	NAG	b	401	-	-	0/6/23/26	0/1/1/1
4	NAG	g	401	3	-	1/6/23/26	0/1/1/1
4	NAG	j	401	3	-	0/6/23/26	0/1/1/1
4	NAG	n	401	3	-	0/6/23/26	0/1/1/1
4	NAG	q	401	3	-	0/6/23/26	0/1/1/1
4	NAG	d	401	3	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	o	401	3	-	0/6/23/26	0/1/1/1
4	NAG	f	401	3	-	0/6/23/26	0/1/1/1
4	NAG	h	401	3	-	0/6/23/26	0/1/1/1
4	NAG	l	401	3	-	0/6/23/26	0/1/1/1
4	NAG	p	401	3	-	2/6/23/26	0/1/1/1
4	NAG	r	401	3	-	0/6/23/26	0/1/1/1
4	NAG	k	401	3	-	0/6/23/26	0/1/1/1
4	NAG	i	401	3	-	0/6/23/26	0/1/1/1

The worst 5 of 7 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	a	401	NAG	C1-C2	2.30	1.55	1.52
4	r	401	NAG	C1-C2	2.22	1.55	1.52
4	h	401	NAG	C1-C2	2.21	1.55	1.52
4	k	401	NAG	C1-C2	2.21	1.55	1.52
4	i	401	NAG	C1-C2	2.19	1.55	1.52

The worst 5 of 9 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	k	401	NAG	C1-O5-C5	2.88	116.04	112.19
4	r	401	NAG	C1-O5-C5	2.75	115.88	112.19
4	a	401	NAG	C1-O5-C5	2.72	115.83	112.19
4	p	401	NAG	C1-O5-C5	2.61	115.68	112.19
4	i	401	NAG	C1-O5-C5	2.59	115.65	112.19

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	c	401	NAG	C4-C5-C6-O6
4	c	401	NAG	O5-C5-C6-O6
4	p	401	NAG	O5-C5-C6-O6
4	p	401	NAG	C4-C5-C6-O6
4	g	401	NAG	C1-C2-N2-C7

There are no ring outliers.

8 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	m	401	NAG	1	0
4	a	401	NAG	1	0
4	c	401	NAG	1	0
4	h	401	NAG	1	0
4	p	401	NAG	1	0
4	r	401	NAG	1	0
4	k	401	NAG	1	0
4	i	401	NAG	1	0

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.

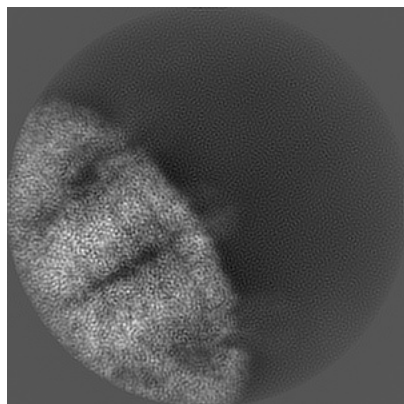
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-21957. These allow visual inspection of the internal detail of the map and identification of artifacts.

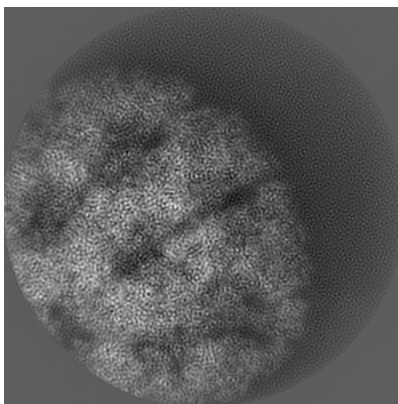
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

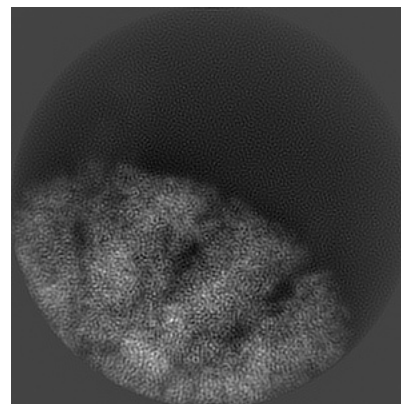
6.1.1 Primary map



X

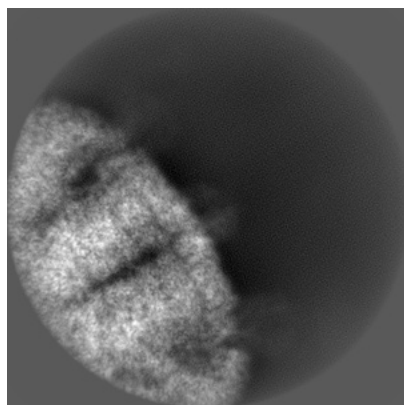


Y

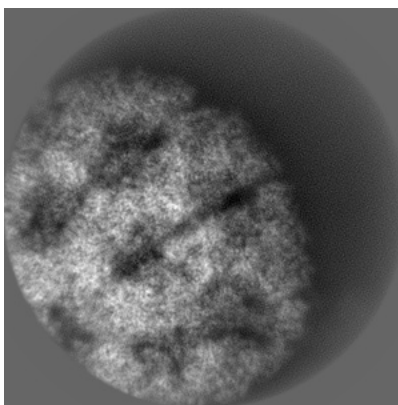


Z

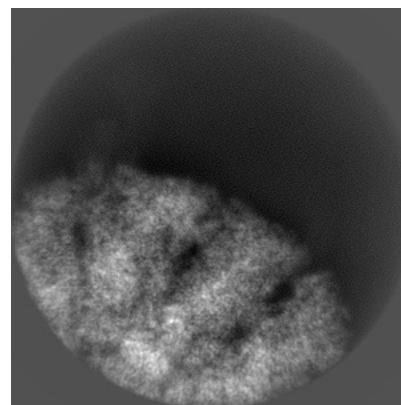
6.1.2 Raw map



X



Y

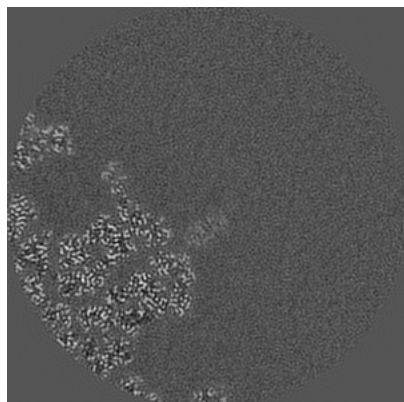


Z

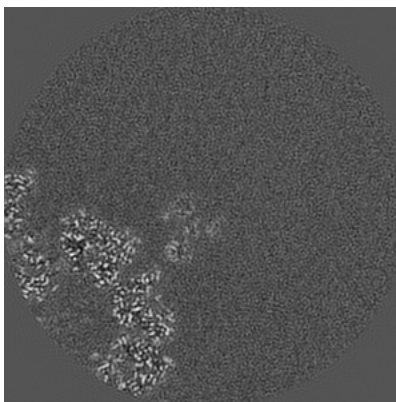
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

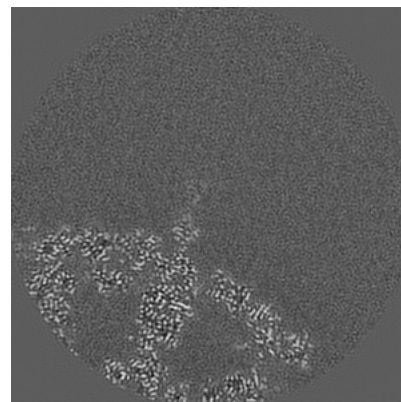
6.2.1 Primary map



X Index: 160

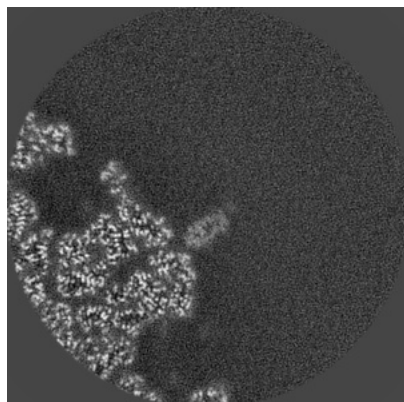


Y Index: 160

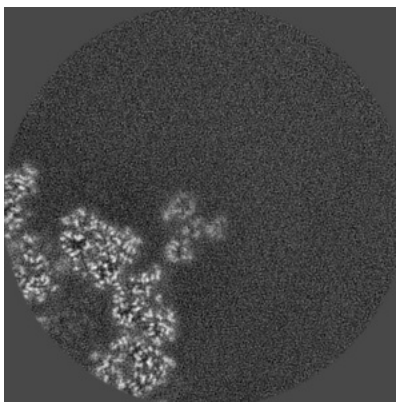


Z Index: 160

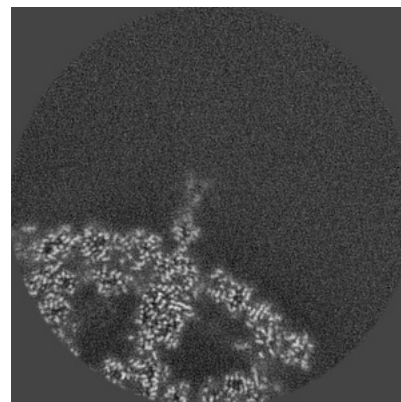
6.2.2 Raw map



X Index: 160



Y Index: 160

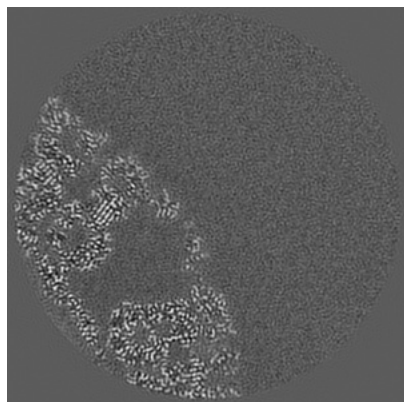


Z Index: 160

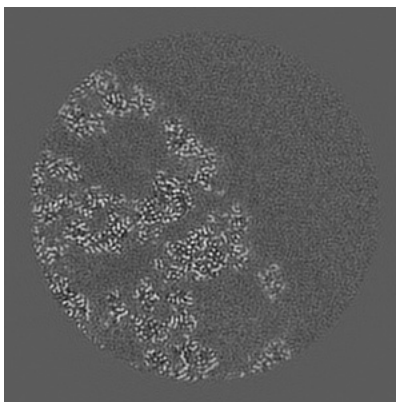
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

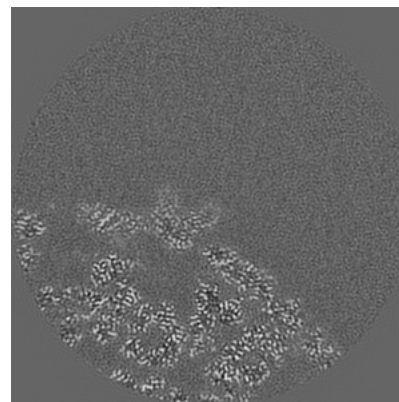
6.3.1 Primary map



X Index: 112

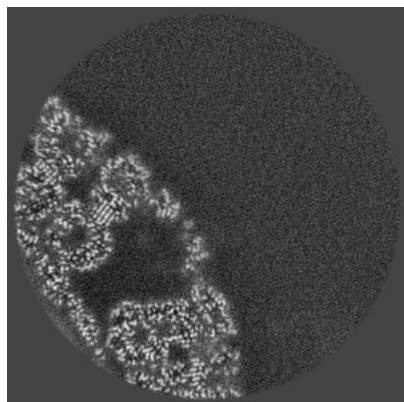


Y Index: 79

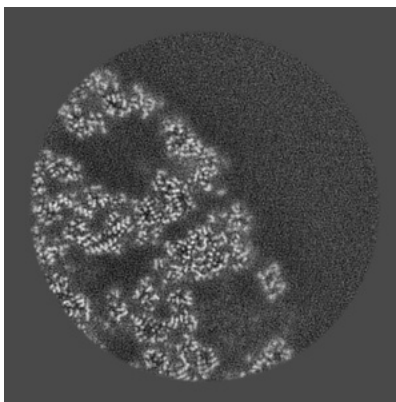


Z Index: 135

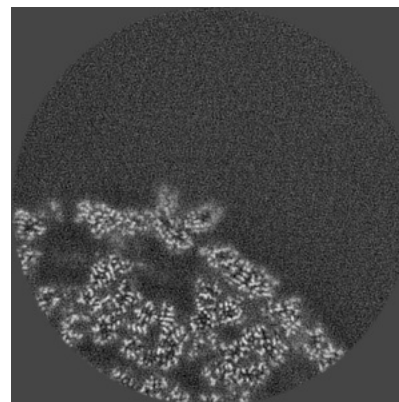
6.3.2 Raw map



X Index: 112



Y Index: 79

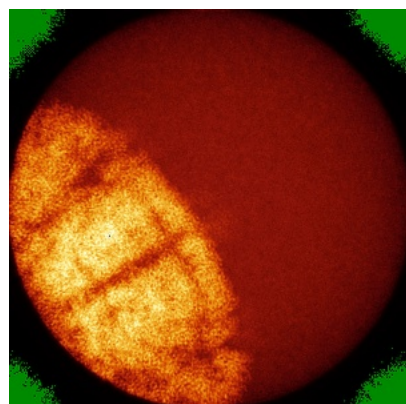


Z Index: 134

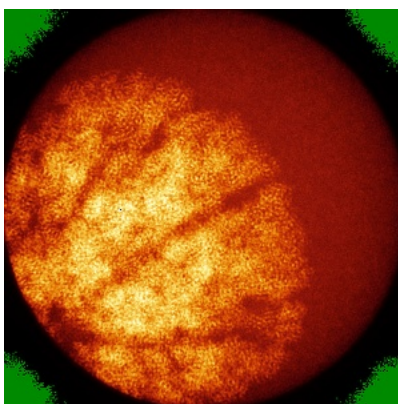
The images above show the largest variance slices of the map in three orthogonal directions.

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

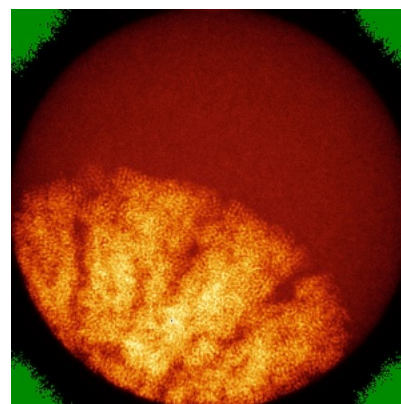
6.4.1 Primary map



X

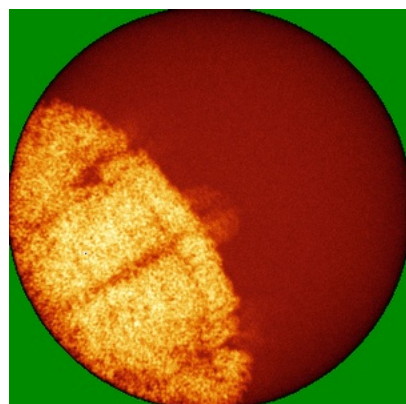


Y

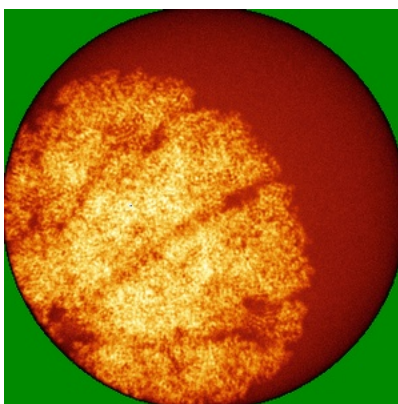


Z

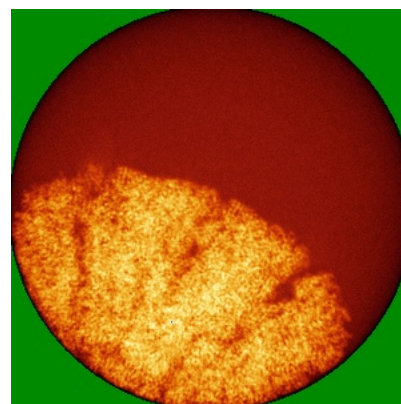
6.4.2 Raw map



X



Y



Z

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

6.5 Orthogonal surface views [i](#)

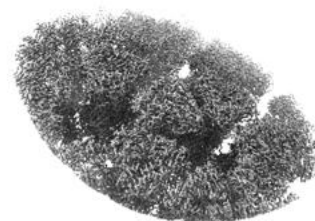
6.5.1 Primary map



X



Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.05. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



X



Y



Z

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

6.6 Mask visualisation [i](#)

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

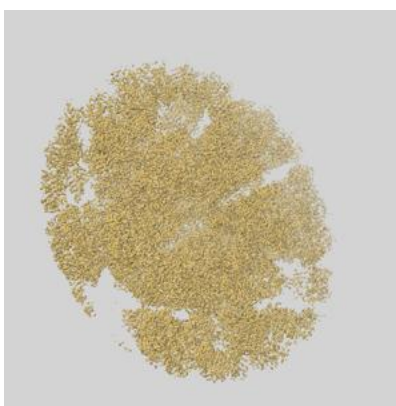
A mask typically either:

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

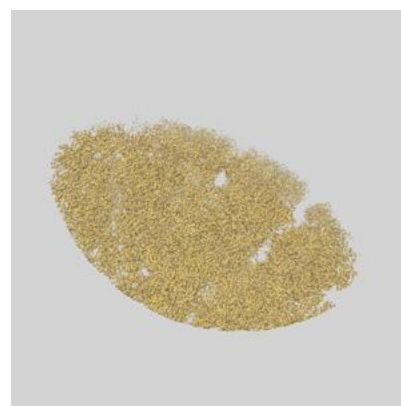
6.6.1 emd_21957_msk_1.map [i](#)



X



Y

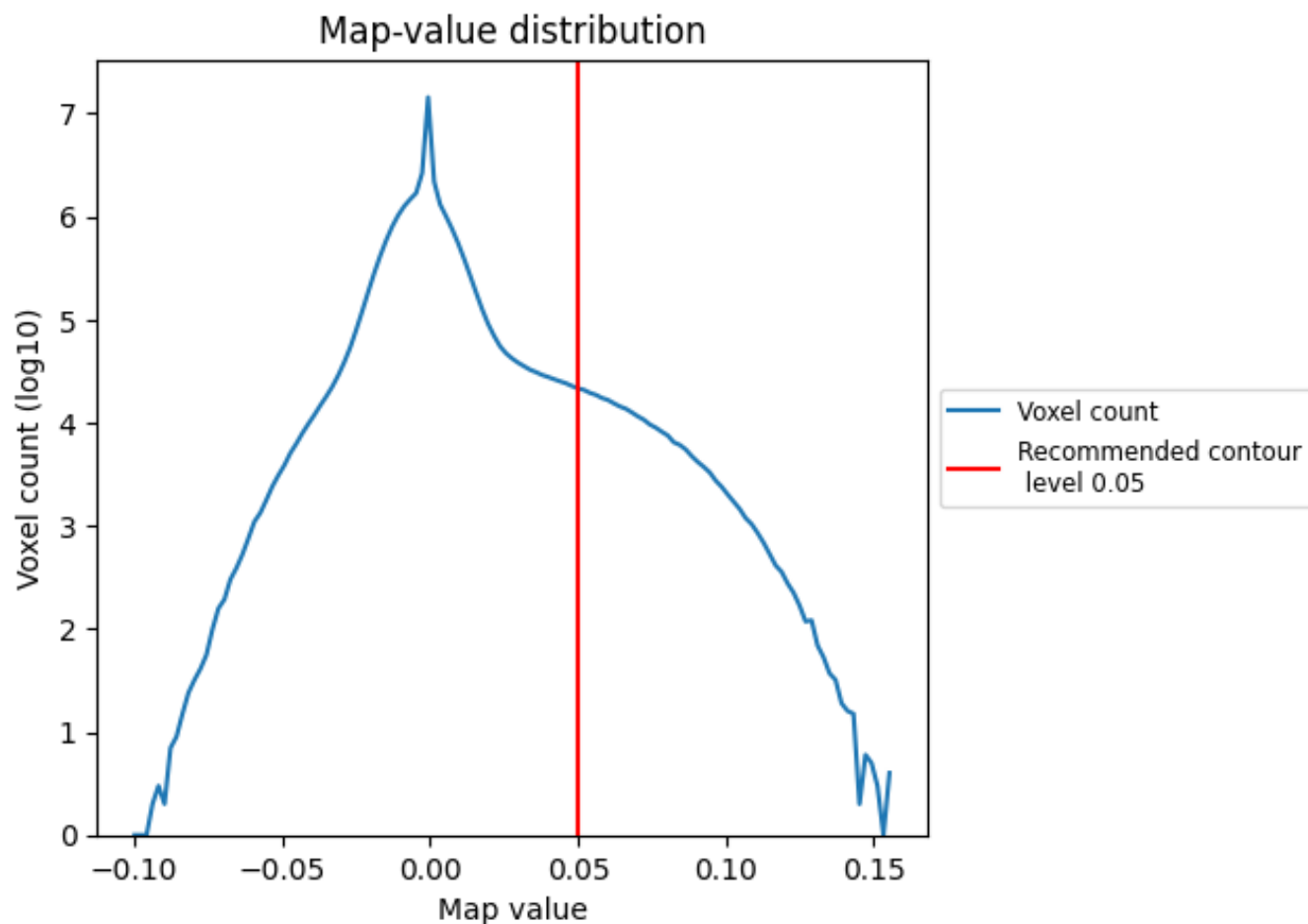


Z

7 Map analysis [i](#)

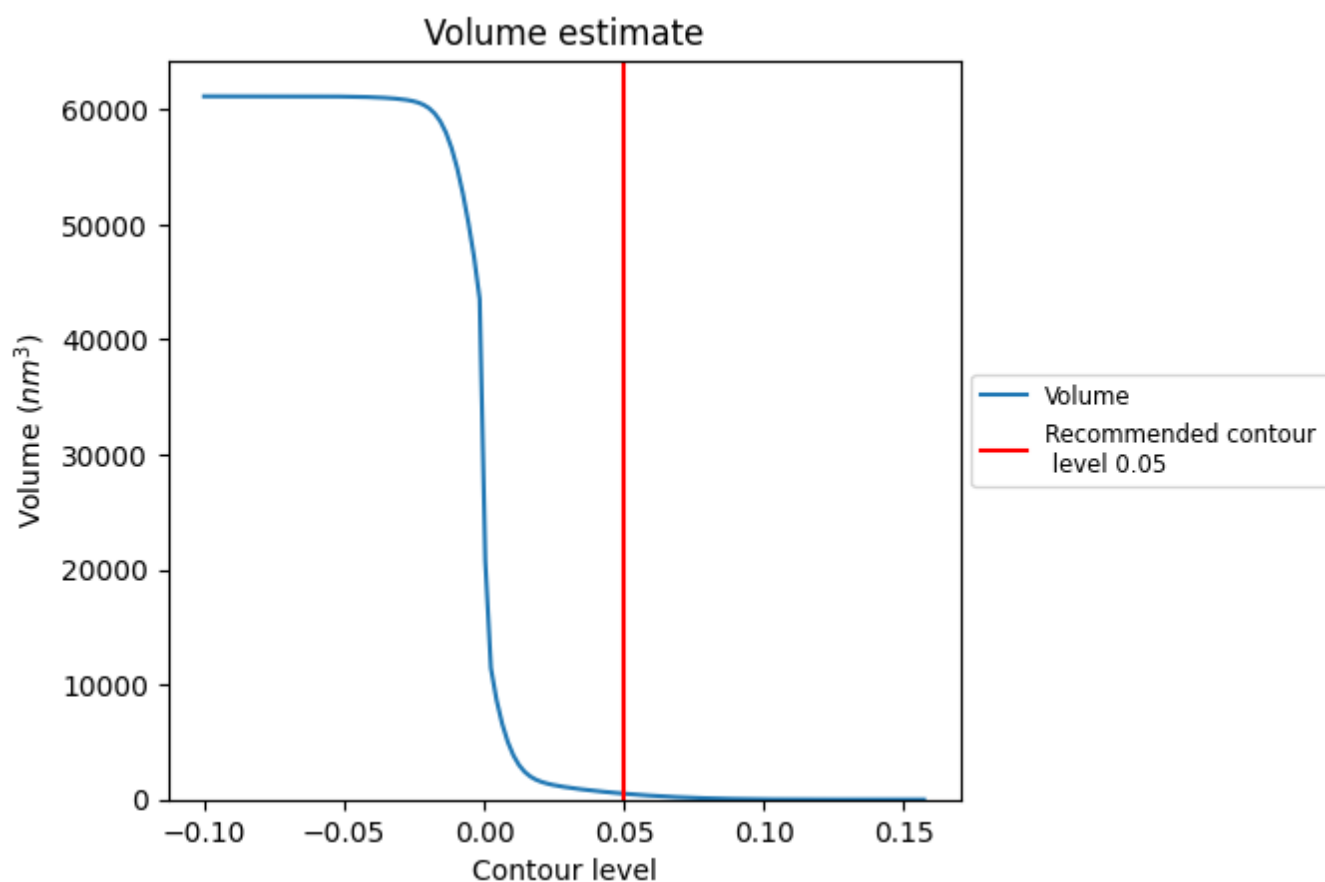
This section contains the results of statistical analysis of the map.

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.

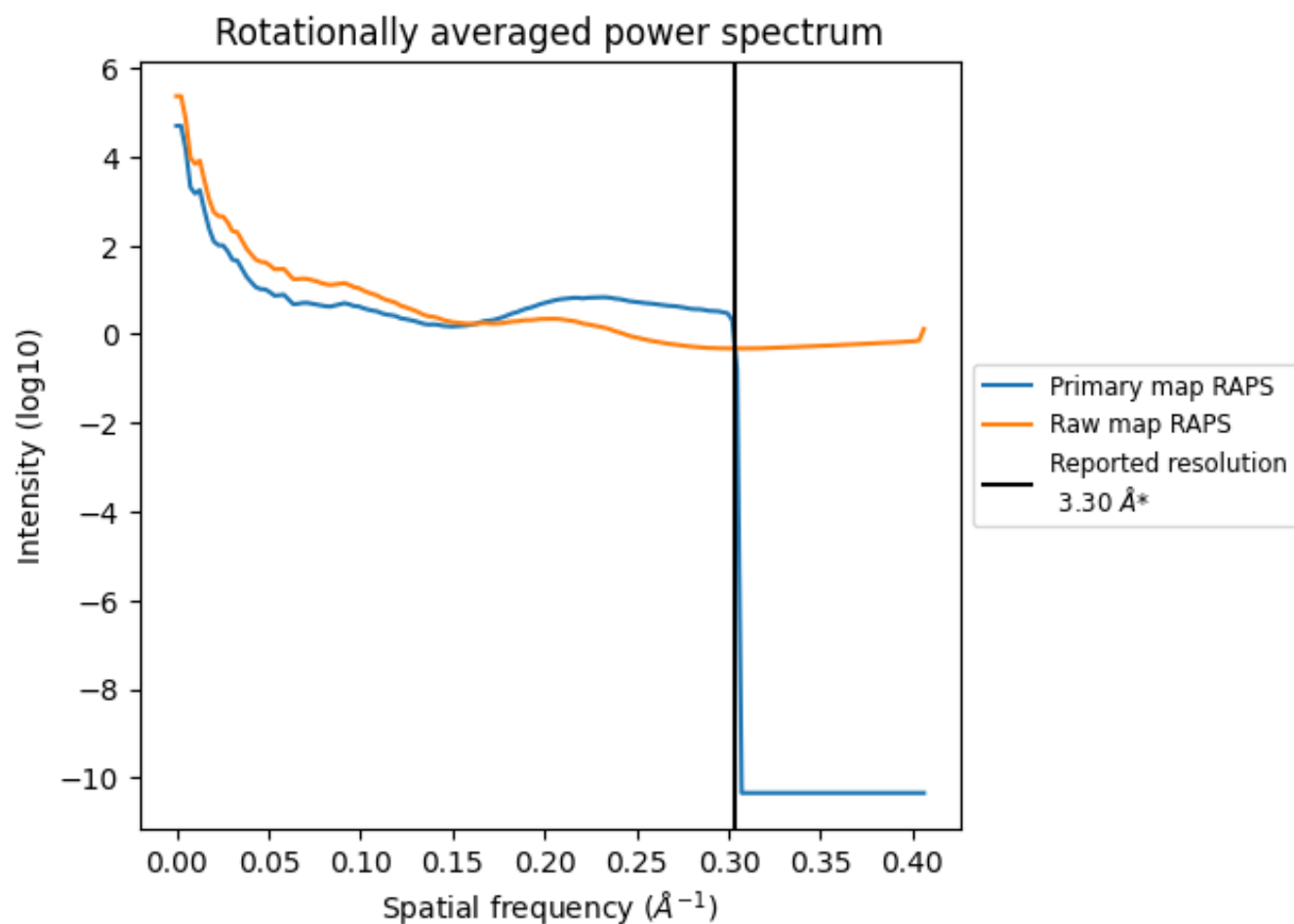
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 516 nm³; this corresponds to an approximate mass of 466 kDa.

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

7.3 Rotationally averaged power spectrum ⓘ

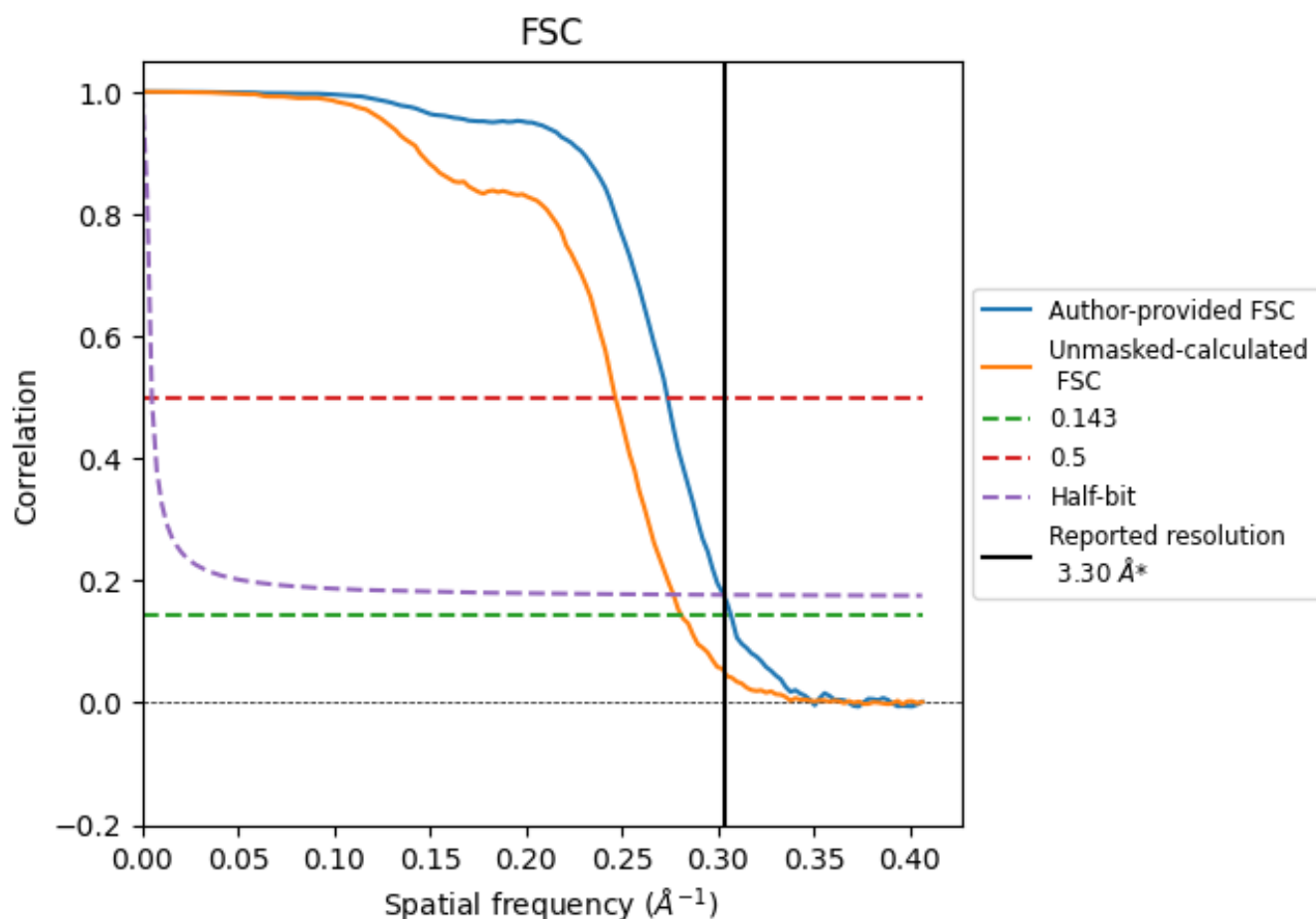


*Reported resolution corresponds to spatial frequency of 0.303 Å⁻¹

8 Fourier-Shell correlation [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 $0.303 \text{ \AA}^{-1}$

8.2 Resolution estimates [i](#)

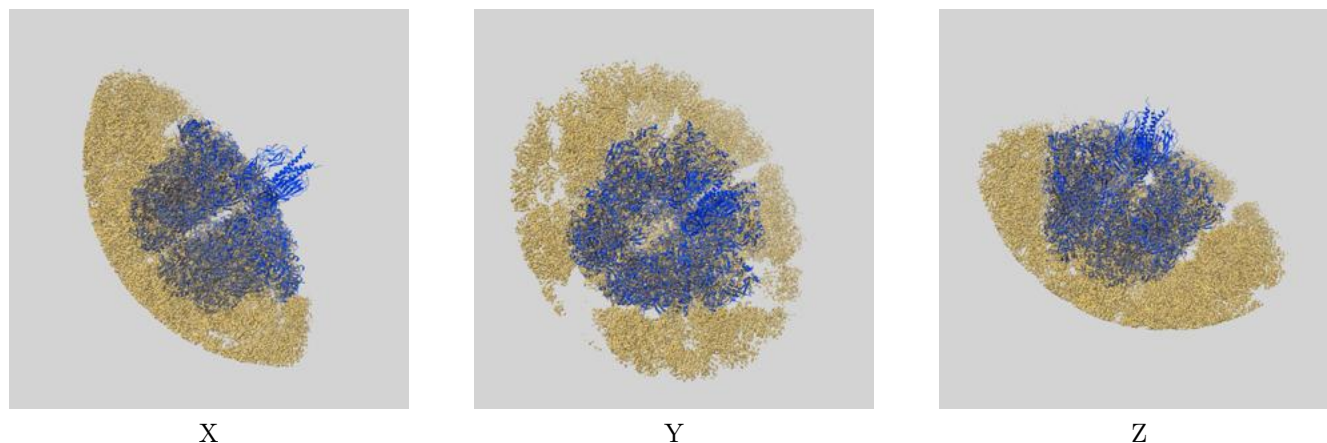
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.30	-	-
Author-provided FSC curve	3.26	3.66	3.30
Unmasked-calculated*	3.56	4.06	3.61

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

9 Map-model fit [i](#)

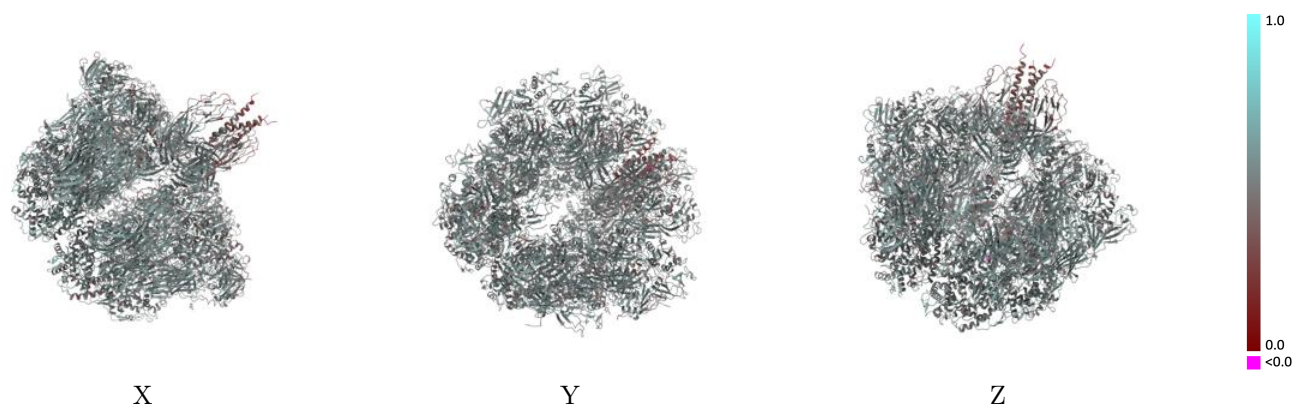
This section contains information regarding the fit between EMDB map EMD-21957 and PDB model 6WXG. Per-residue inclusion information can be found in [section 3](#) on [page 10](#).

9.1 Map-model overlay [i](#)



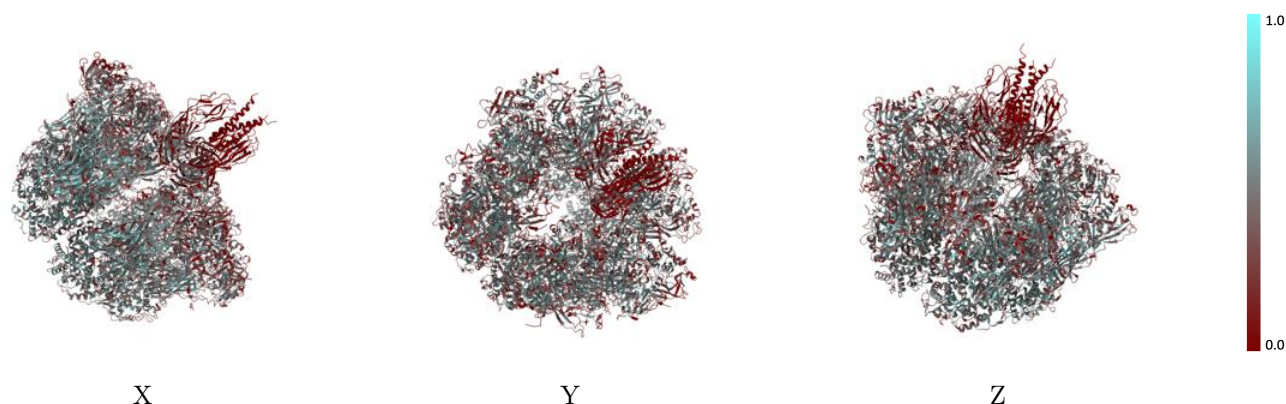
The images above show the 3D surface view of the map at the recommended contour level 0.05 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



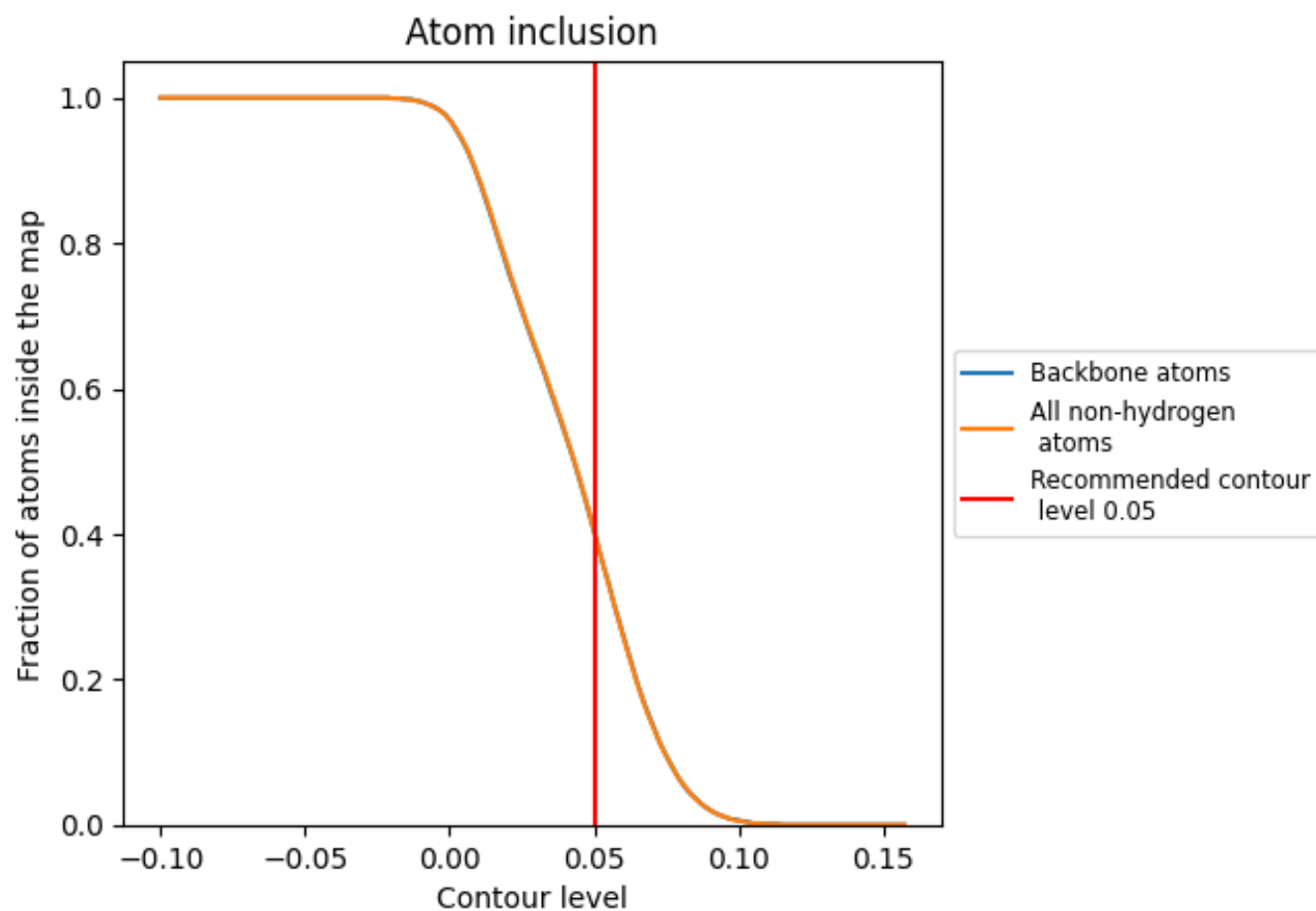
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.05).

9.4 Atom inclusion [i](#)



At the recommended contour level, 40% of all backbone atoms, 40% 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 (0.05) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.4040	 0.5090
1	 0.0880	 0.4590
2	 0.1000	 0.4570
3	 0.1000	 0.4490
A	 0.5210	 0.5290
B	 0.4960	 0.5150
C	 0.5130	 0.5250
D	 0.5190	 0.5280
E	 0.5100	 0.5220
F	 0.5070	 0.5260
G	 0.5100	 0.5220
H	 0.4900	 0.5220
I	 0.5020	 0.5280
J	 0.4770	 0.5180
K	 0.4730	 0.5200
L	 0.4750	 0.5190
M	 0.4800	 0.5200
N	 0.4770	 0.5230
O	 0.4580	 0.5190
P	 0.4820	 0.5230
Q	 0.4910	 0.5220
R	 0.4660	 0.5170
a	 0.3450	 0.5060
b	 0.3500	 0.4970
c	 0.3600	 0.5020
d	 0.3950	 0.5090
e	 0.4000	 0.5100
f	 0.3760	 0.5070
g	 0.3240	 0.4960
h	 0.3180	 0.5000
i	 0.3640	 0.4970
j	 0.3280	 0.4920
k	 0.3170	 0.4960
l	 0.3630	 0.5110
m	 0.3470	 0.4960



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Chain	Atom inclusion	Q-score
n	 0.3270	 0.5000
o	 0.3480	 0.5000
p	 0.2940	 0.4930
q	 0.3140	 0.4980
r	 0.3220	 0.4930