Full wwPDB EM Validation Report ⓘ

Feb 25, 2024 – 12:55 PM EST

PDB ID	:	5IPI
EMDB ID	:	EMD-8099
Title	:	Structure of Adeno-associated virus type 2 VLP
Authors	:	Drouin, L.M.; Lins, B.; Janssen, M.E.; Bennet, A.; Chipman, P.; McKenna, R.; Chen, W.; Muzyczka, N.; Cardone, G.; Baker, T.S.; Agbandje-McKenna, M.
Deposited on	:	2016-03-09
Resolution	:	3.80 Å(reported)

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

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/EMValidationReportHelp
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

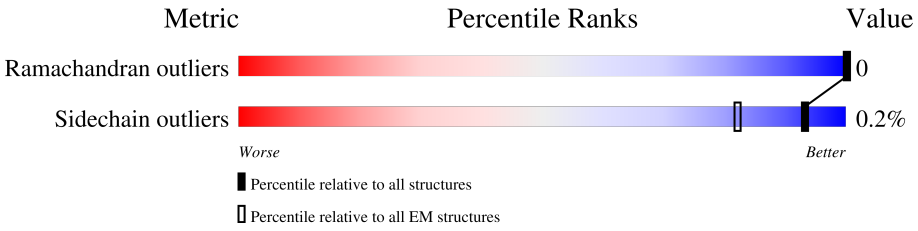
EMDB validation analysis	:	0.0.1.dev70
MolProbity	:	4.02b-467
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ	:	1.9.13
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 3.80 Å.

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



Metric	Whole archive (#Entries)	EM structures (#Entries)
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826

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	1	735	17% 69% 30%
1	2	735	17% 69% 30%
1	3	735	16% 69% 30%
1	4	735	17% 69% 30%
1	5	735	17% 69% 30%
1	6	735	17% 69% 30%
1	7	735	16% 69% 30%
1	8	735	16% 69% 30%
1	A	735	17% 69% 30%

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Mol	Chain	Length	Quality of chain
1	B	735	17% 69% 30%
1	C	735	16% 69% 30%
1	D	735	17% 69% 30%
1	E	735	16% 69% 30%
1	F	735	17% 69% 30%
1	G	735	17% 69% 30%
1	H	735	17% 69% 30%
1	I	735	17% 69% 30%
1	J	735	17% 69% 30%
1	K	735	17% 69% 30%
1	L	735	17% 69% 30%
1	M	735	16% 69% 30%
1	N	735	17% 69% 30%
1	O	735	17% 69% 30%
1	P	735	17% 69% 30%
1	Q	735	16% 69% 30%
1	R	735	16% 69% 30%
1	S	735	17% 69% 30%
1	T	735	17% 69% 30%
1	U	735	17% 69% 30%
1	V	735	17% 69% 30%
1	W	735	17% 69% 30%
1	X	735	17% 69% 30%
1	Y	735	16% 69% 30%
1	Z	735	17% 69% 30%

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Mol	Chain	Length	Quality of chain
1	a	735	16% 69% 30%
1	b	735	17% 69% 30%
1	c	735	17% 69% 30%
1	d	735	17% 69% 30%
1	e	735	17% 69% 30%
1	f	735	16% 69% 30%
1	g	735	17% 69% 30%
1	h	735	16% 69% 30%
1	i	735	17% 69% 30%
1	j	735	16% 69% 30%
1	k	735	17% 69% 30%
1	l	735	17% 69% 30%
1	m	735	17% 69% 30%
1	n	735	17% 69% 30%
1	o	735	17% 69% 30%
1	p	735	17% 69% 30%
1	q	735	17% 69% 30%
1	r	735	17% 69% 30%
1	s	735	17% 69% 30%
1	t	735	17% 69% 30%
1	u	735	17% 69% 30%
1	v	735	16% 69% 30%
1	w	735	17% 69% 30%
1	x	735	17% 69% 30%
1	y	735	17% 69% 30%

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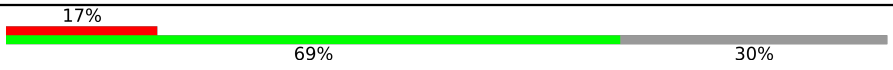
Mol	Chain	Length	Quality of chain
1	z	735	

2 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 246360 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 Capsid protein VP1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	511	Total 4106	C 2588	N 716	O 789	S 13	0	0
1	B	511	Total 4106	C 2588	N 716	O 789	S 13	0	0
1	C	511	Total 4106	C 2588	N 716	O 789	S 13	0	0
1	D	511	Total 4106	C 2588	N 716	O 789	S 13	0	0
1	E	511	Total 4106	C 2588	N 716	O 789	S 13	0	0
1	F	511	Total 4106	C 2588	N 716	O 789	S 13	0	0
1	G	511	Total 4106	C 2588	N 716	O 789	S 13	0	0
1	H	511	Total 4106	C 2588	N 716	O 789	S 13	0	0
1	I	511	Total 4106	C 2588	N 716	O 789	S 13	0	0
1	J	511	Total 4106	C 2588	N 716	O 789	S 13	0	0
1	K	511	Total 4106	C 2588	N 716	O 789	S 13	0	0
1	L	511	Total 4106	C 2588	N 716	O 789	S 13	0	0
1	M	511	Total 4106	C 2588	N 716	O 789	S 13	0	0
1	N	511	Total 4106	C 2588	N 716	O 789	S 13	0	0
1	O	511	Total 4106	C 2588	N 716	O 789	S 13	0	0
1	P	511	Total 4106	C 2588	N 716	O 789	S 13	0	0
1	Q	511	Total 4106	C 2588	N 716	O 789	S 13	0	0

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Mol	Chain	Residues	Atoms					AltConf	Trace
1	R	511	Total	C	N	O	S	0	0
			4106	2588	716	789	13		
1	S	511	Total	C	N	O	S	0	0
			4106	2588	716	789	13		
1	T	511	Total	C	N	O	S	0	0
			4106	2588	716	789	13		
1	U	511	Total	C	N	O	S	0	0
			4106	2588	716	789	13		
1	V	511	Total	C	N	O	S	0	0
			4106	2588	716	789	13		
1	W	511	Total	C	N	O	S	0	0
			4106	2588	716	789	13		
1	X	511	Total	C	N	O	S	0	0
			4106	2588	716	789	13		
1	Y	511	Total	C	N	O	S	0	0
			4106	2588	716	789	13		
1	Z	511	Total	C	N	O	S	0	0
			4106	2588	716	789	13		
1	a	511	Total	C	N	O	S	0	0
			4106	2588	716	789	13		
1	b	511	Total	C	N	O	S	0	0
			4106	2588	716	789	13		
1	c	511	Total	C	N	O	S	0	0
			4106	2588	716	789	13		
1	d	511	Total	C	N	O	S	0	0
			4106	2588	716	789	13		
1	e	511	Total	C	N	O	S	0	0
			4106	2588	716	789	13		
1	f	511	Total	C	N	O	S	0	0
			4106	2588	716	789	13		
1	g	511	Total	C	N	O	S	0	0
			4106	2588	716	789	13		
1	h	511	Total	C	N	O	S	0	0
			4106	2588	716	789	13		
1	i	511	Total	C	N	O	S	0	0
			4106	2588	716	789	13		
1	j	511	Total	C	N	O	S	0	0
			4106	2588	716	789	13		
1	k	511	Total	C	N	O	S	0	0
			4106	2588	716	789	13		
1	l	511	Total	C	N	O	S	0	0
			4106	2588	716	789	13		

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Mol	Chain	Residues	Atoms					AltConf	Trace
1	m	511	Total	C	N	O	S	0	0
			4106	2588	716	789	13		
1	n	511	Total	C	N	O	S	0	0
			4106	2588	716	789	13		
1	o	511	Total	C	N	O	S	0	0
			4106	2588	716	789	13		
1	p	511	Total	C	N	O	S	0	0
			4106	2588	716	789	13		
1	q	511	Total	C	N	O	S	0	0
			4106	2588	716	789	13		
1	r	511	Total	C	N	O	S	0	0
			4106	2588	716	789	13		
1	s	511	Total	C	N	O	S	0	0
			4106	2588	716	789	13		
1	t	511	Total	C	N	O	S	0	0
			4106	2588	716	789	13		
1	u	511	Total	C	N	O	S	0	0
			4106	2588	716	789	13		
1	v	511	Total	C	N	O	S	0	0
			4106	2588	716	789	13		
1	w	511	Total	C	N	O	S	0	0
			4106	2588	716	789	13		
1	x	511	Total	C	N	O	S	0	0
			4106	2588	716	789	13		
1	y	511	Total	C	N	O	S	0	0
			4106	2588	716	789	13		
1	z	511	Total	C	N	O	S	0	0
			4106	2588	716	789	13		
1	1	511	Total	C	N	O	S	0	0
			4106	2588	716	789	13		
1	2	511	Total	C	N	O	S	0	0
			4106	2588	716	789	13		
1	3	511	Total	C	N	O	S	0	0
			4106	2588	716	789	13		
1	4	511	Total	C	N	O	S	0	0
			4106	2588	716	789	13		
1	5	511	Total	C	N	O	S	0	0
			4106	2588	716	789	13		
1	6	511	Total	C	N	O	S	0	0
			4106	2588	716	789	13		
1	7	511	Total	C	N	O	S	0	0
			4106	2588	716	789	13		

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Mol	Chain	Residues	Atoms					AltConf	Trace
1	8	511	Total	C	N	O	S	0	0
			4106	2588	716	789	13		

There are 60 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	452	THR	SER	conflict	UNP P03135
B	452	THR	SER	conflict	UNP P03135
C	452	THR	SER	conflict	UNP P03135
D	452	THR	SER	conflict	UNP P03135
E	452	THR	SER	conflict	UNP P03135
F	452	THR	SER	conflict	UNP P03135
G	452	THR	SER	conflict	UNP P03135
H	452	THR	SER	conflict	UNP P03135
I	452	THR	SER	conflict	UNP P03135
J	452	THR	SER	conflict	UNP P03135
K	452	THR	SER	conflict	UNP P03135
L	452	THR	SER	conflict	UNP P03135
M	452	THR	SER	conflict	UNP P03135
N	452	THR	SER	conflict	UNP P03135
O	452	THR	SER	conflict	UNP P03135
P	452	THR	SER	conflict	UNP P03135
Q	452	THR	SER	conflict	UNP P03135
R	452	THR	SER	conflict	UNP P03135
S	452	THR	SER	conflict	UNP P03135
T	452	THR	SER	conflict	UNP P03135
U	452	THR	SER	conflict	UNP P03135
V	452	THR	SER	conflict	UNP P03135
W	452	THR	SER	conflict	UNP P03135
X	452	THR	SER	conflict	UNP P03135
Y	452	THR	SER	conflict	UNP P03135
Z	452	THR	SER	conflict	UNP P03135
a	452	THR	SER	conflict	UNP P03135
b	452	THR	SER	conflict	UNP P03135
c	452	THR	SER	conflict	UNP P03135
d	452	THR	SER	conflict	UNP P03135
e	452	THR	SER	conflict	UNP P03135
f	452	THR	SER	conflict	UNP P03135
g	452	THR	SER	conflict	UNP P03135
h	452	THR	SER	conflict	UNP P03135
i	452	THR	SER	conflict	UNP P03135
j	452	THR	SER	conflict	UNP P03135

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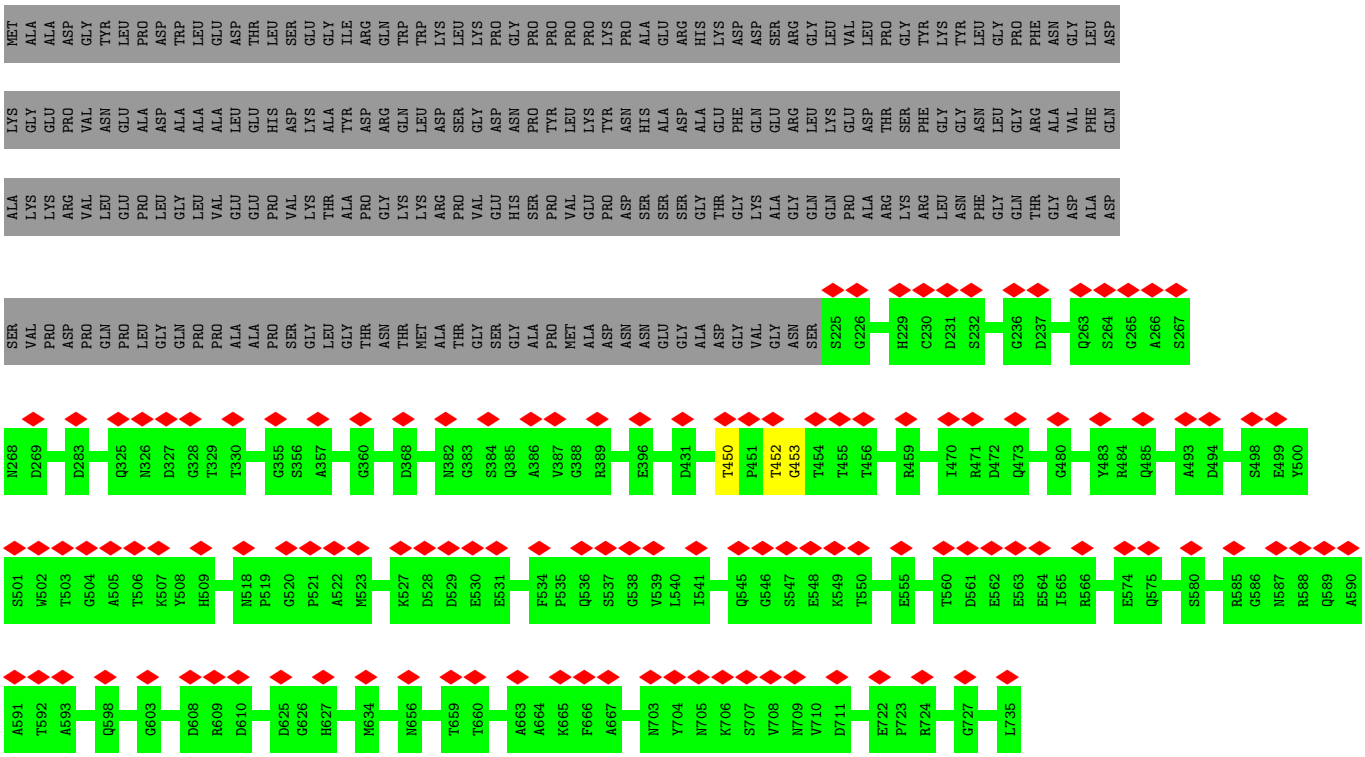
Chain	Residue	Modelled	Actual	Comment	Reference
k	452	THR	SER	conflict	UNP P03135
l	452	THR	SER	conflict	UNP P03135
m	452	THR	SER	conflict	UNP P03135
n	452	THR	SER	conflict	UNP P03135
o	452	THR	SER	conflict	UNP P03135
p	452	THR	SER	conflict	UNP P03135
q	452	THR	SER	conflict	UNP P03135
r	452	THR	SER	conflict	UNP P03135
s	452	THR	SER	conflict	UNP P03135
t	452	THR	SER	conflict	UNP P03135
u	452	THR	SER	conflict	UNP P03135
v	452	THR	SER	conflict	UNP P03135
w	452	THR	SER	conflict	UNP P03135
x	452	THR	SER	conflict	UNP P03135
y	452	THR	SER	conflict	UNP P03135
z	452	THR	SER	conflict	UNP P03135
1	452	THR	SER	conflict	UNP P03135
2	452	THR	SER	conflict	UNP P03135
3	452	THR	SER	conflict	UNP P03135
4	452	THR	SER	conflict	UNP P03135
5	452	THR	SER	conflict	UNP P03135
6	452	THR	SER	conflict	UNP P03135
7	452	THR	SER	conflict	UNP P03135
8	452	THR	SER	conflict	UNP P03135

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.

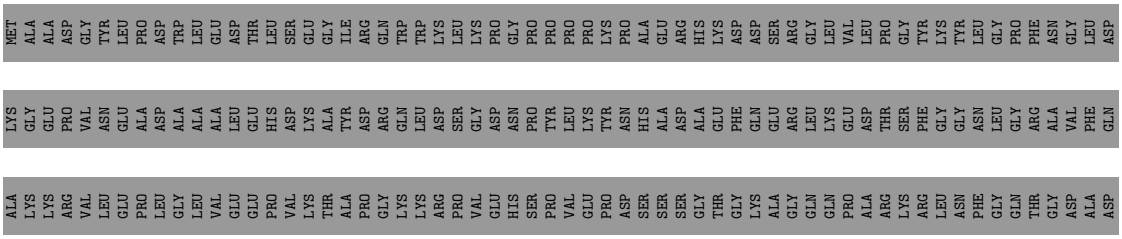
• Molecule 1: Capsid protein VP1

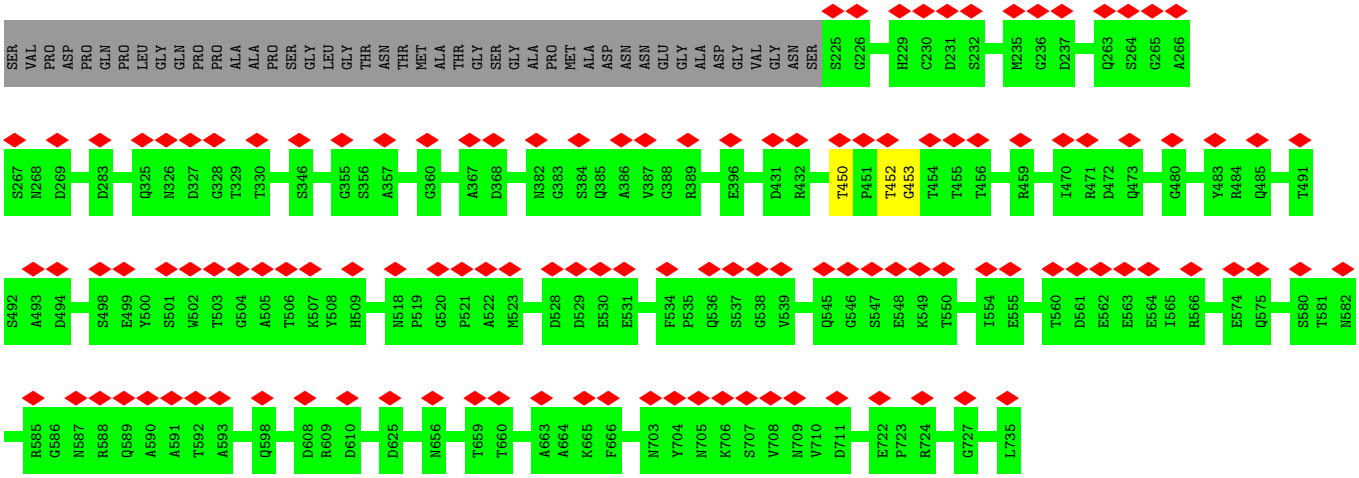
Chain A:



• Molecule 1: Capsid protein VP1

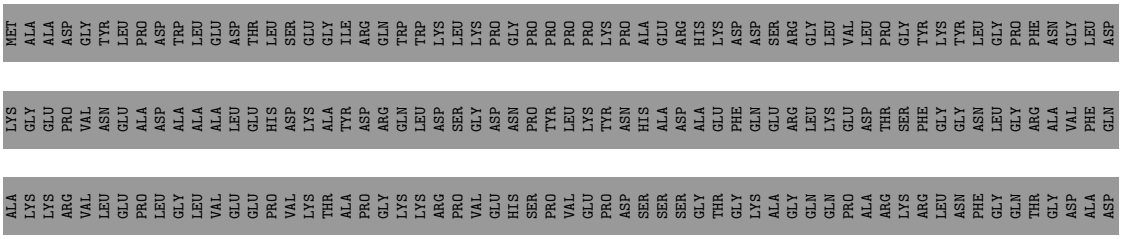
Chain B:

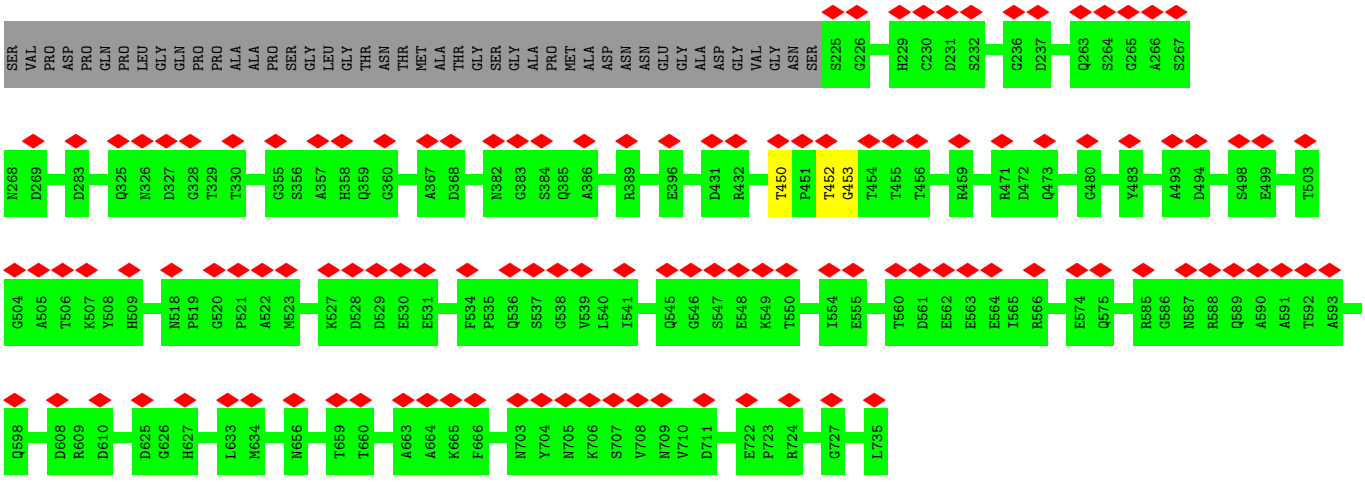




• Molecule 1: Capsid protein VP1

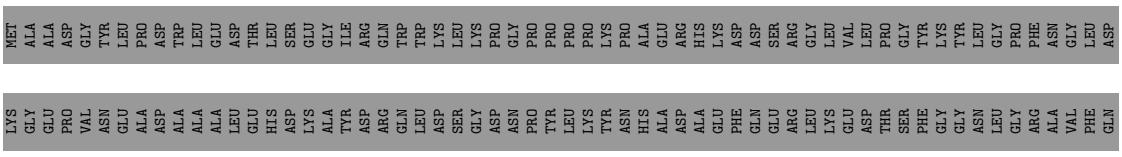
Chain C:

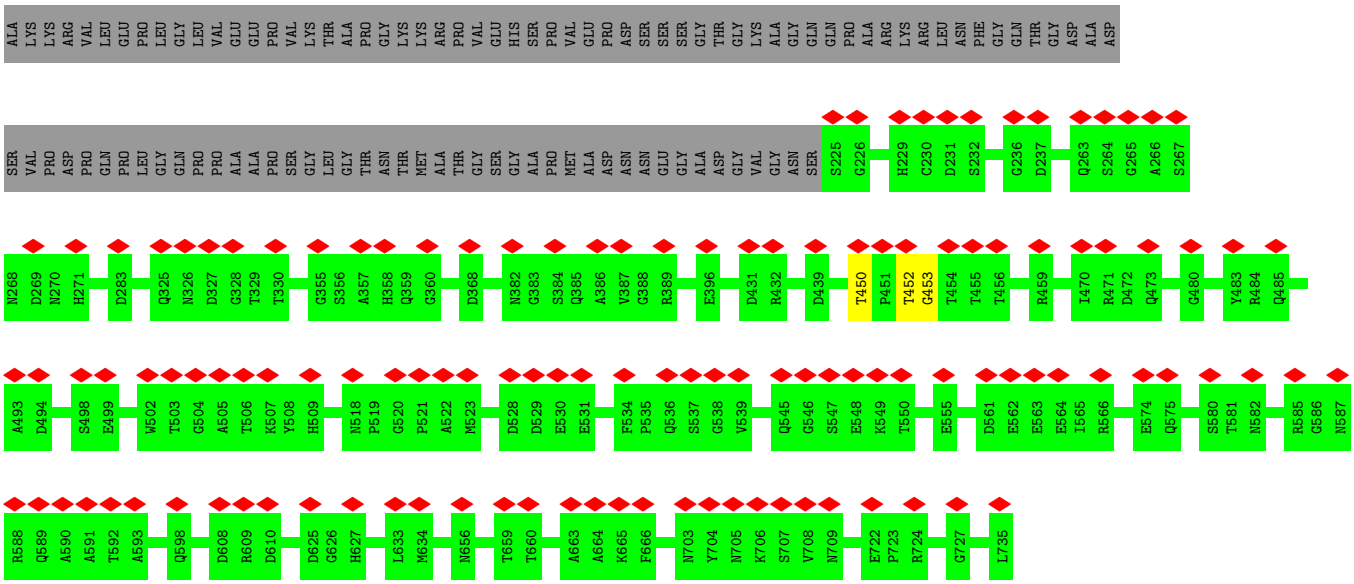




• Molecule 1: Capsid protein VP1

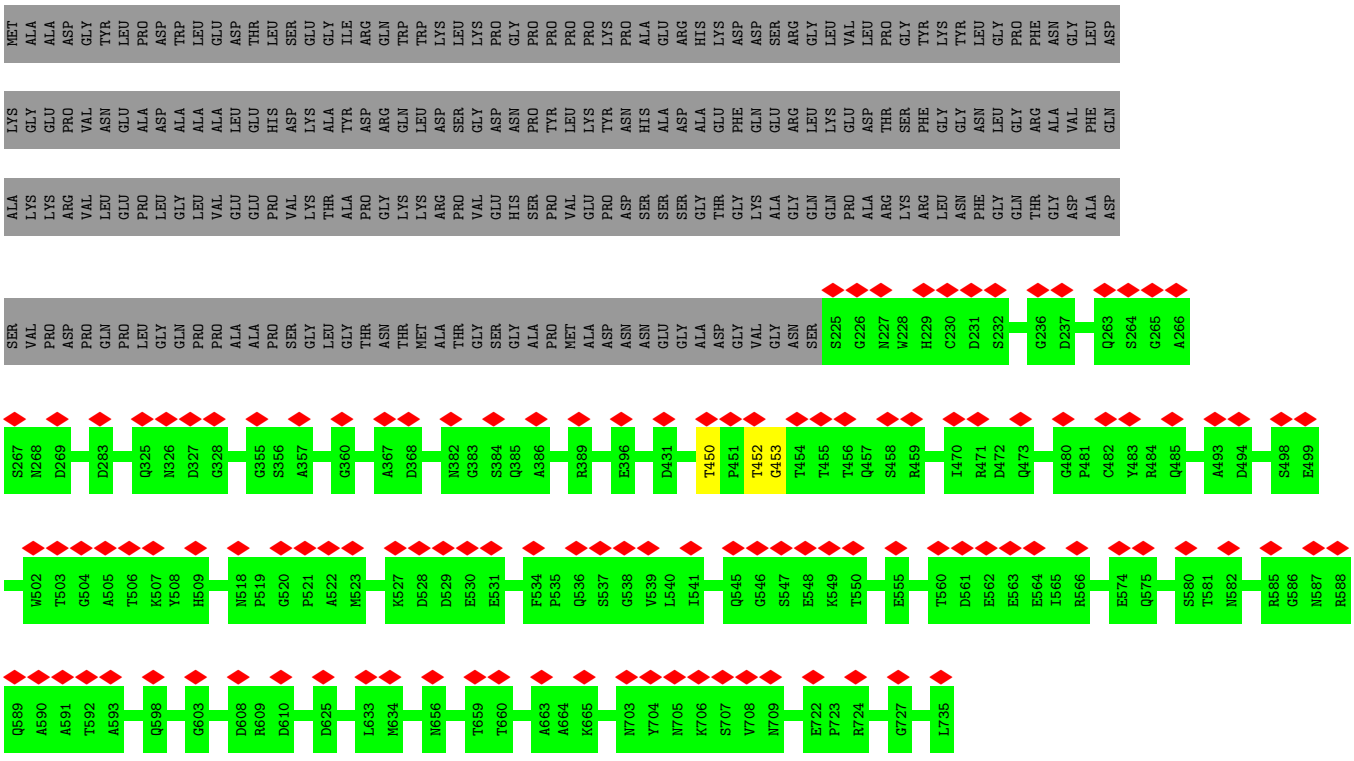
Chain D:





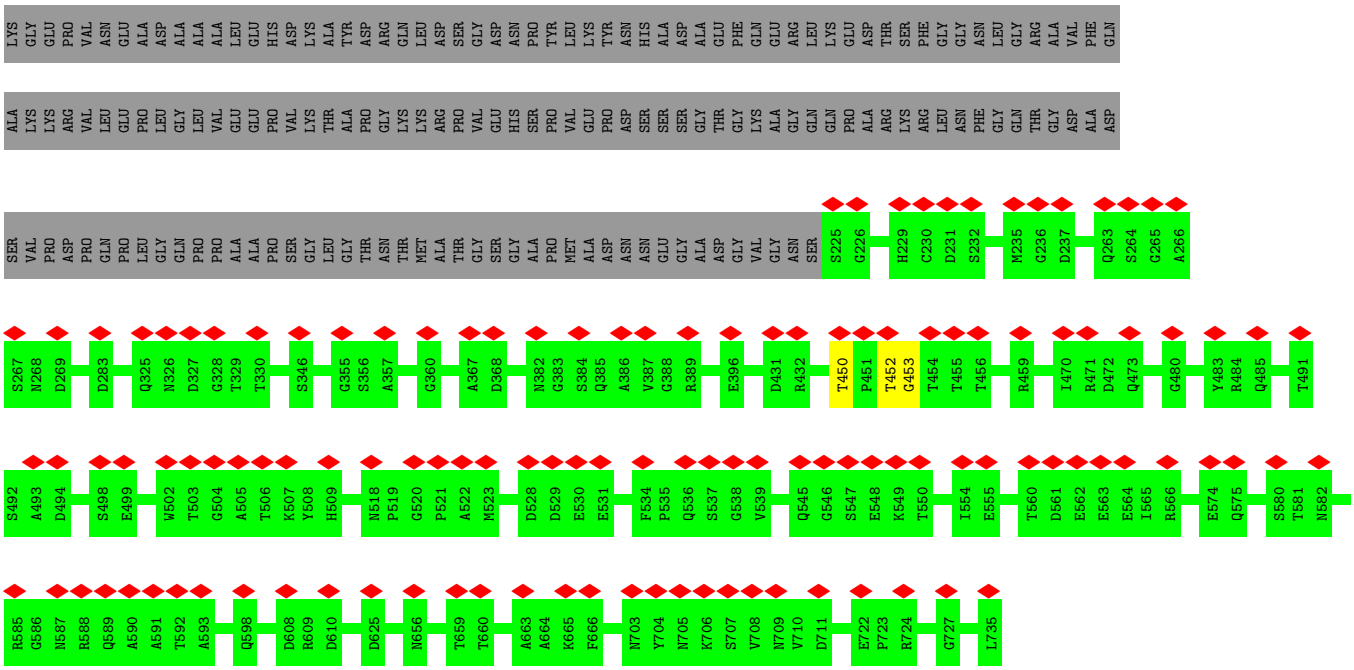
• Molecule 1: Capsid protein VP1

Chain E:



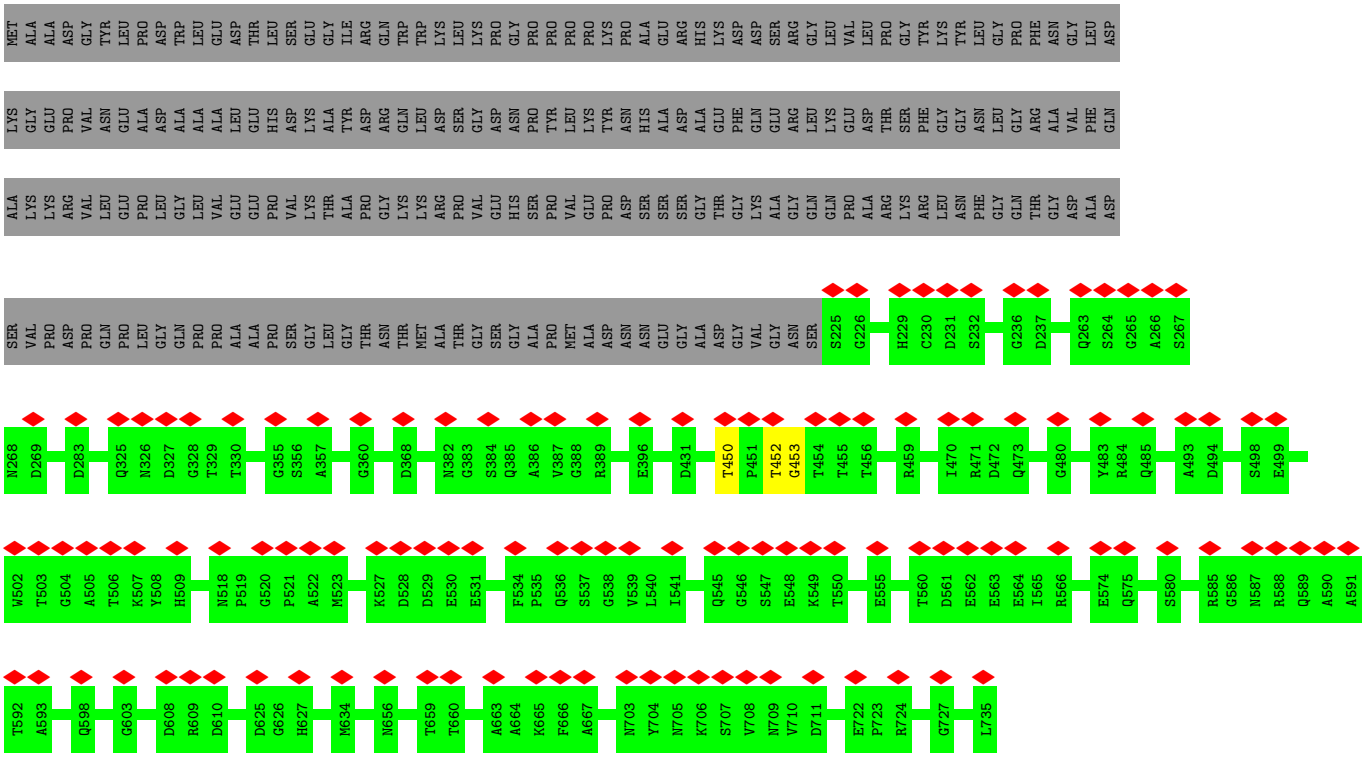
• Molecule 1: Capsid protein VP1

Chain F:



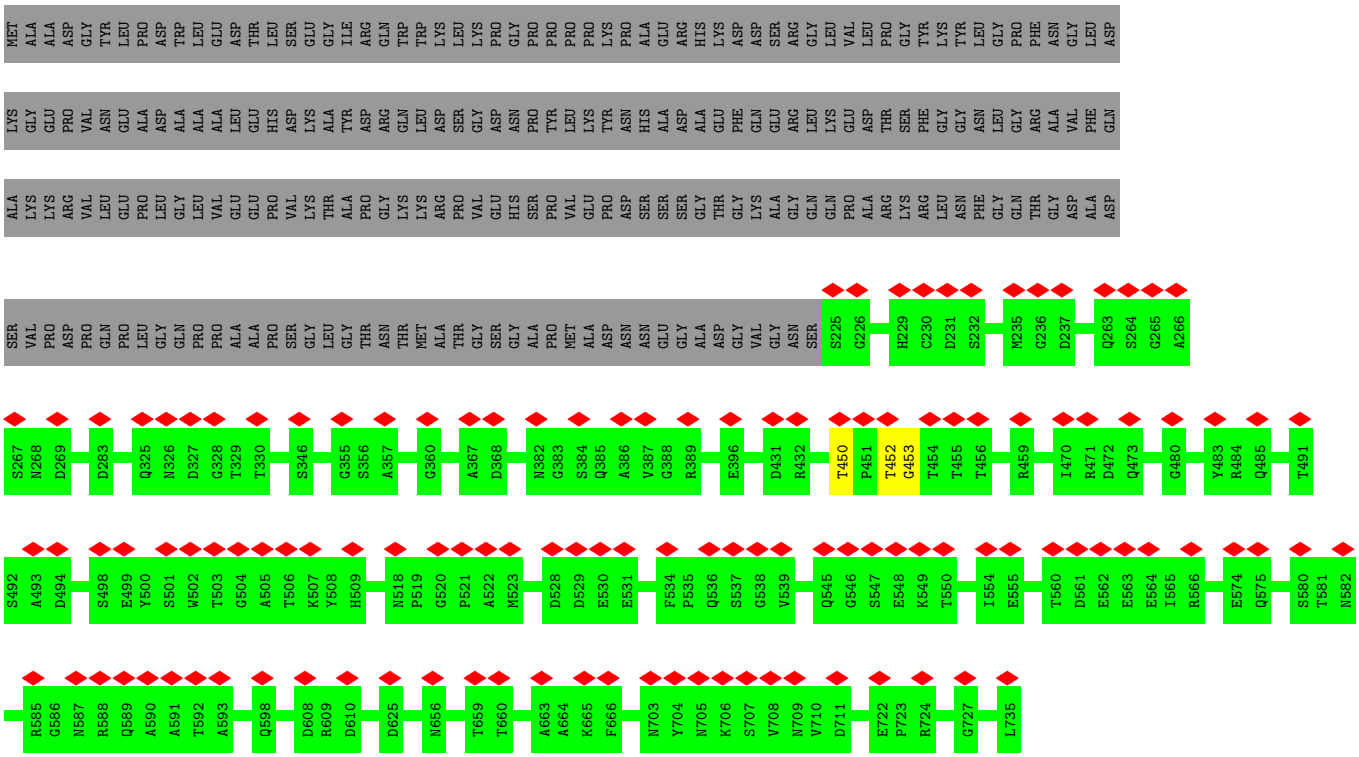
• Molecule 1: Capsid protein VP1

Chain G:



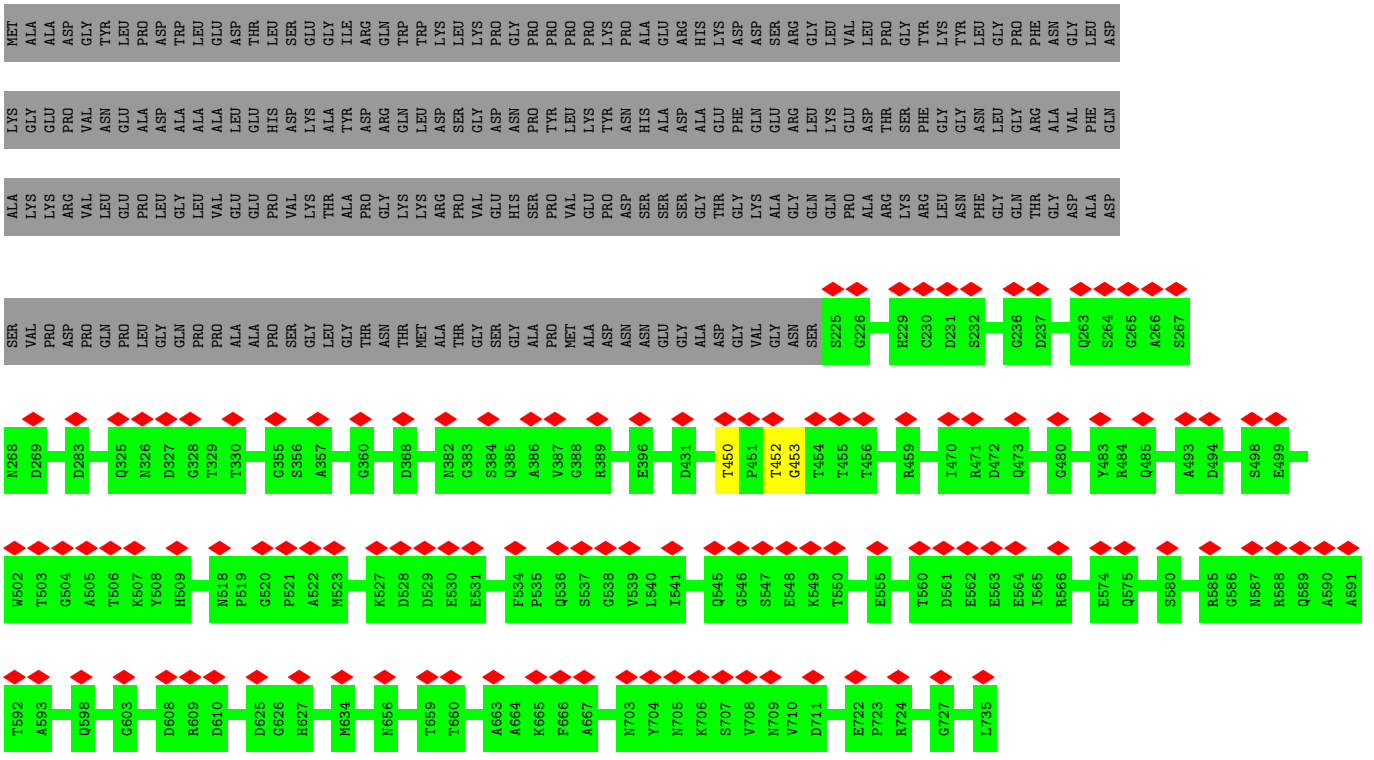
• Molecule 1: Capsid protein VP1

Chain H:



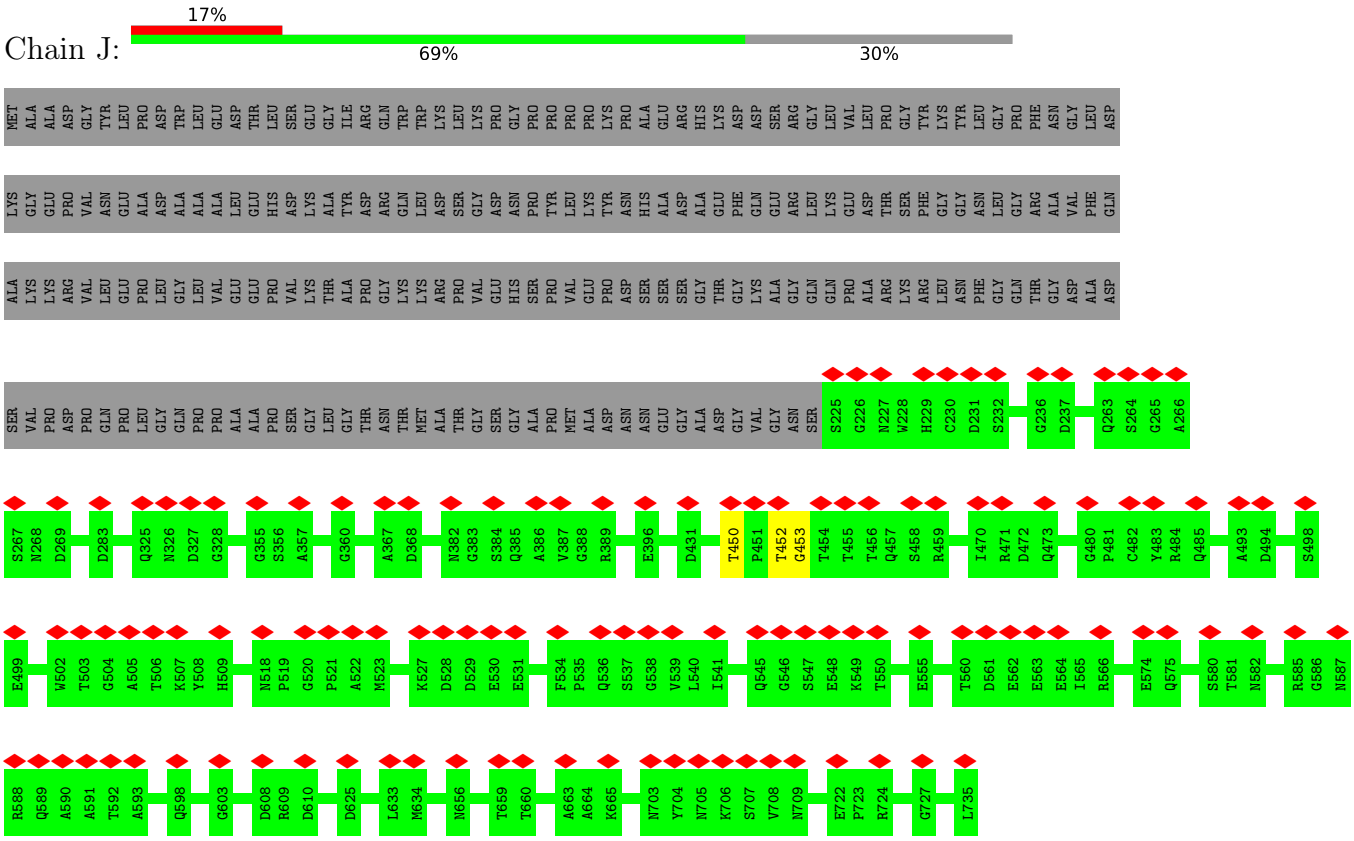
• Molecule 1: Capsid protein VP1

Chain I:



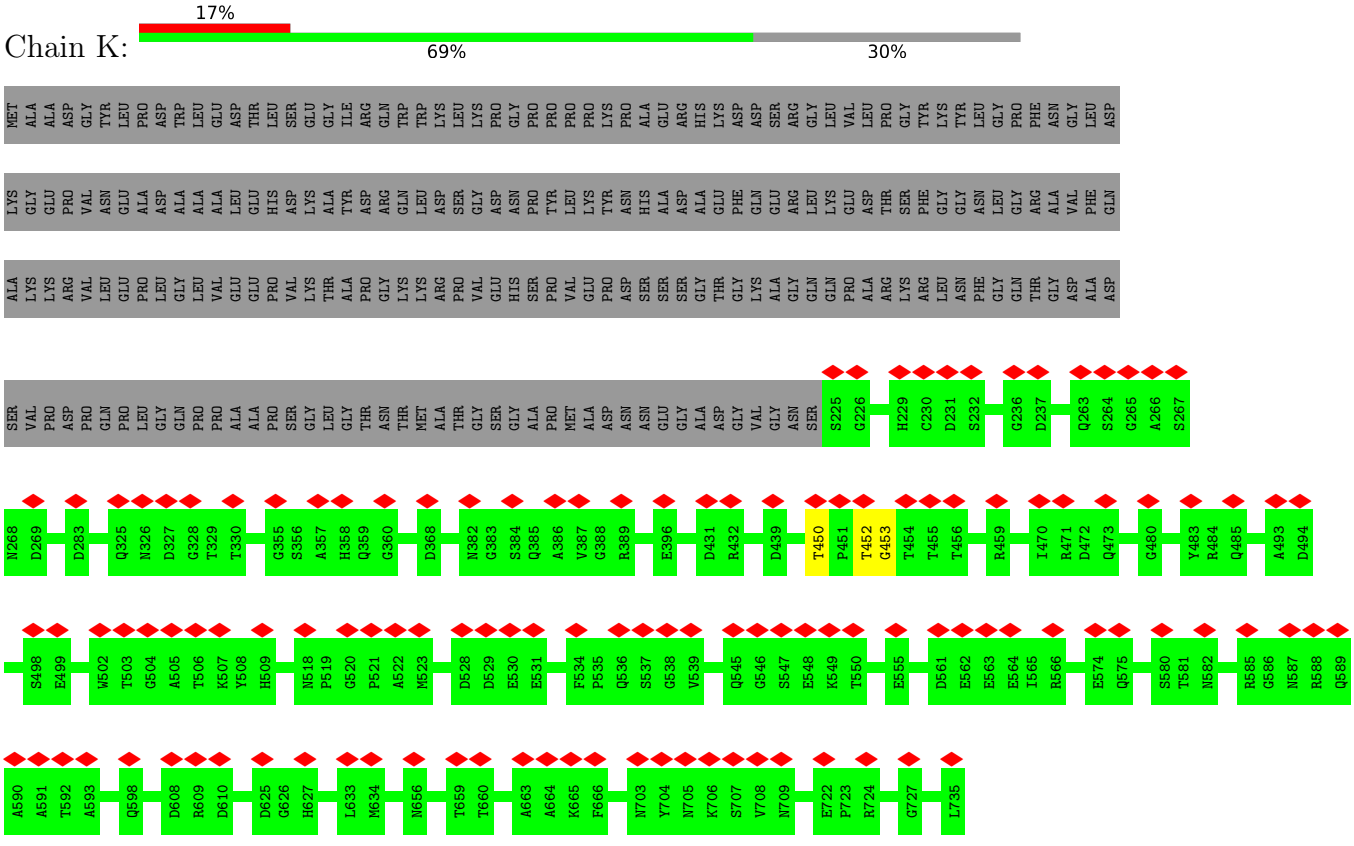
• Molecule 1: Capsid protein VP1

Chain J:



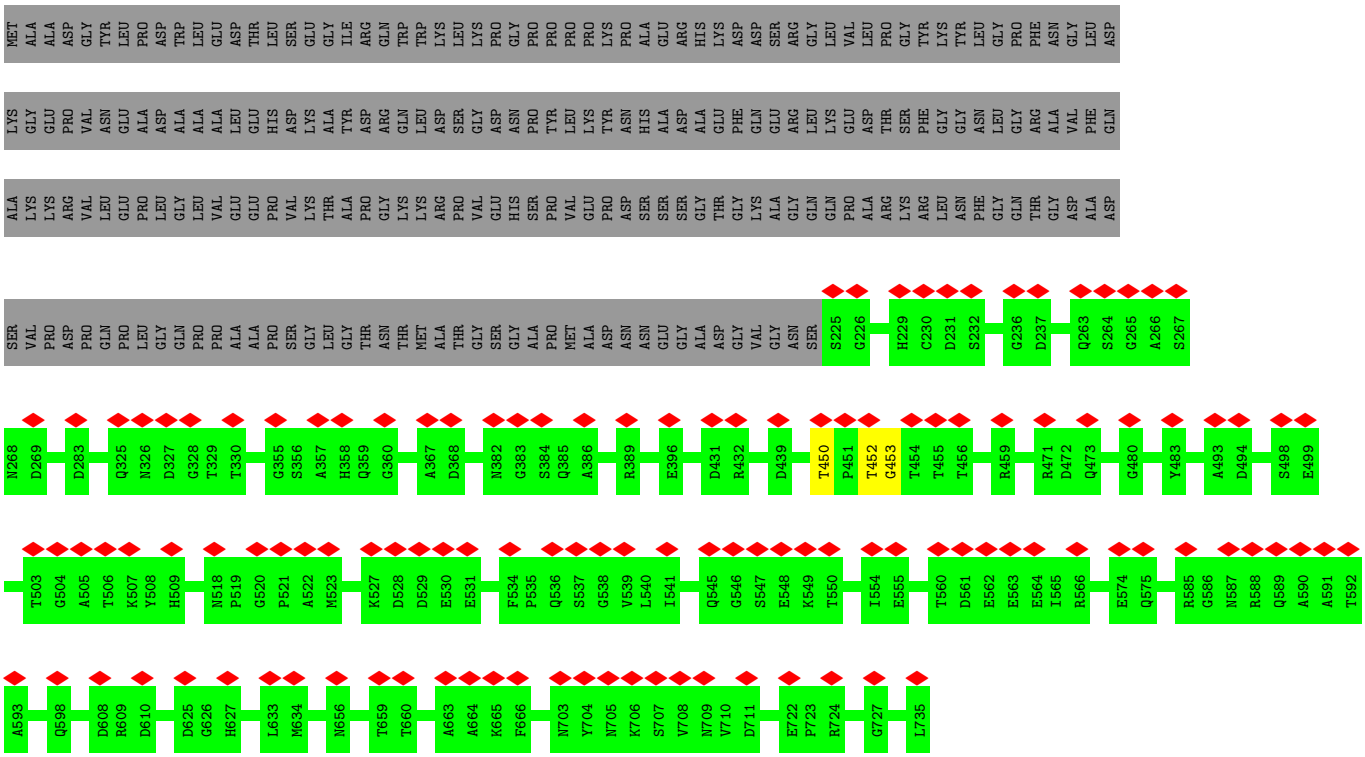
• Molecule 1: Capsid protein VP1

Chain K:



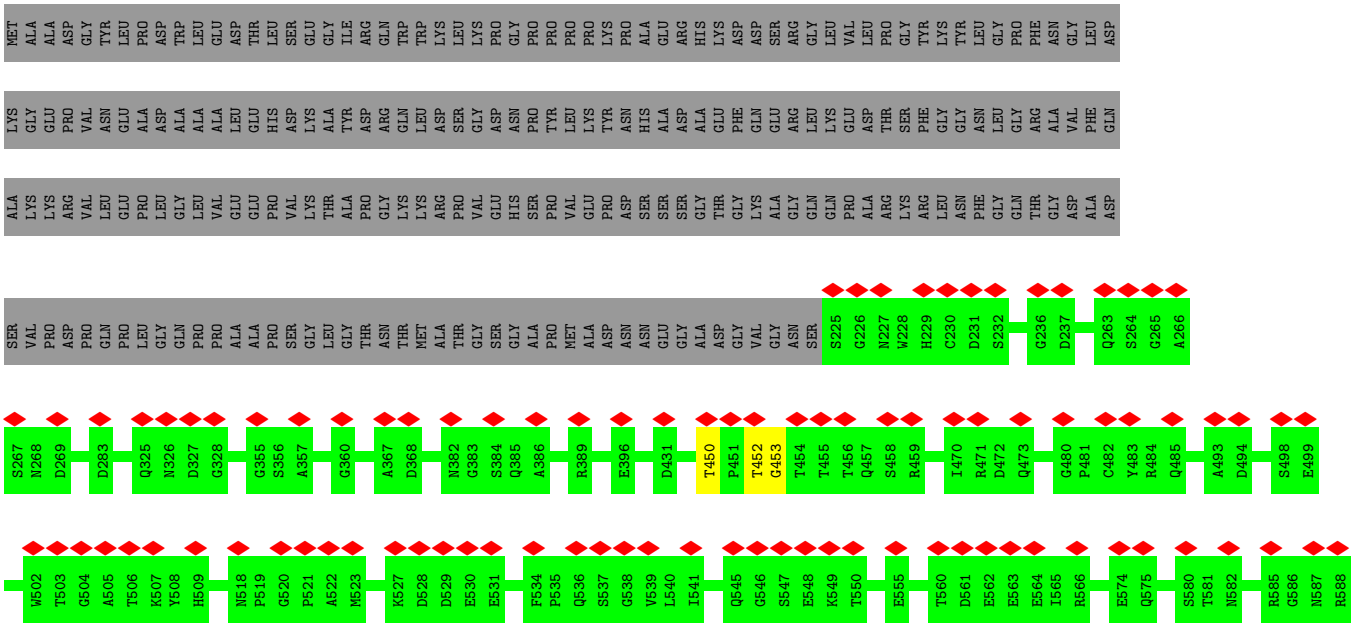
• Molecule 1: Capsid protein VP1

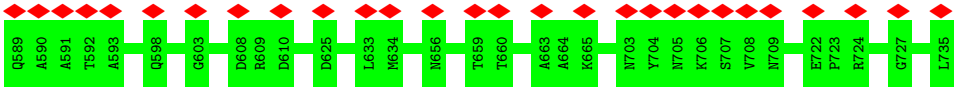
Chain L:



• Molecule 1: Capsid protein VP1

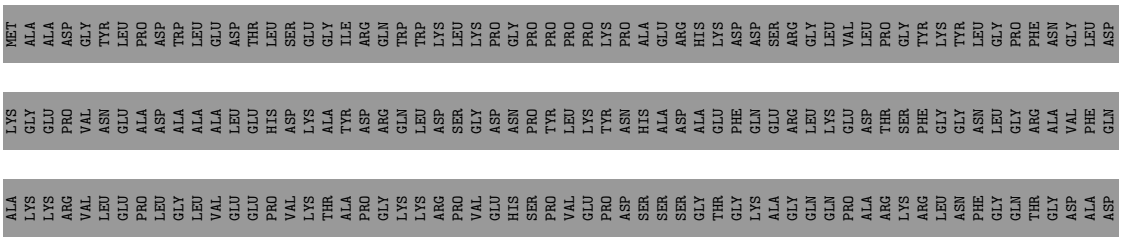
Chain M:





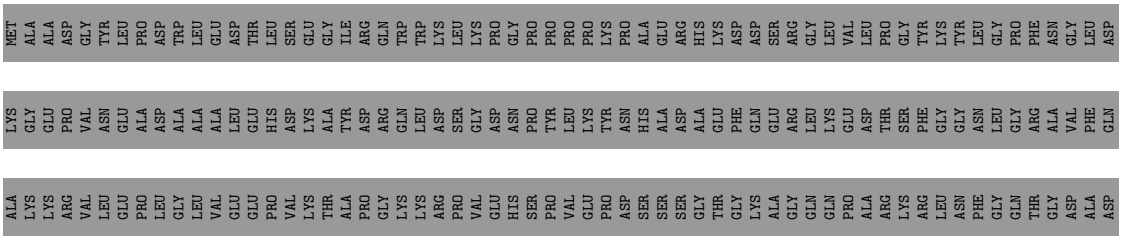
• Molecule 1: Capsid protein VP1

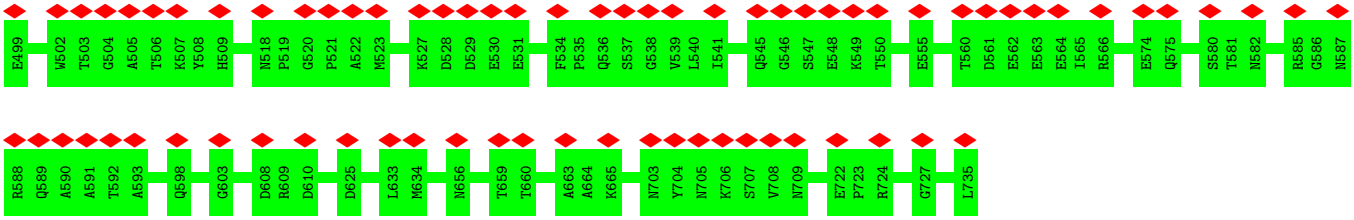
Chain N:



• Molecule 1: Capsid protein VP1

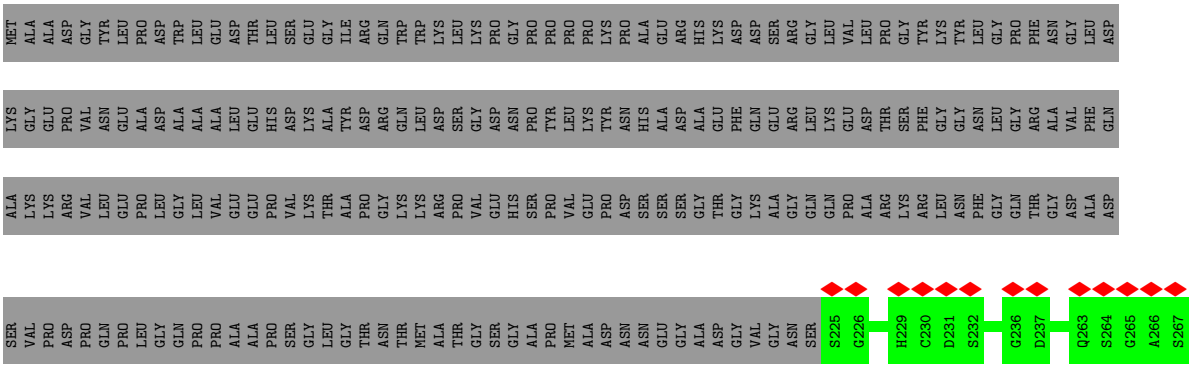
Chain O:

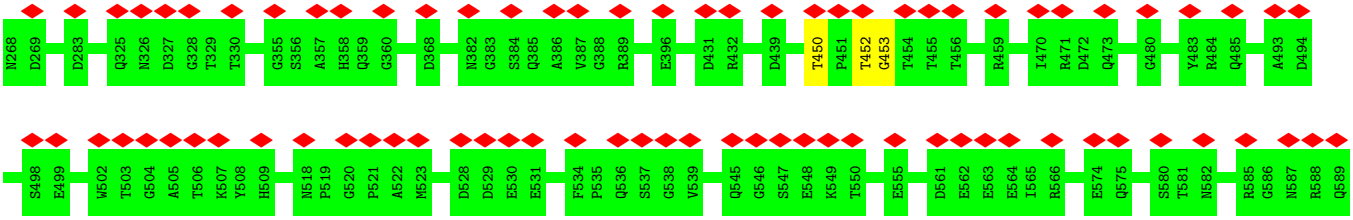




• Molecule 1: Capsid protein VP1

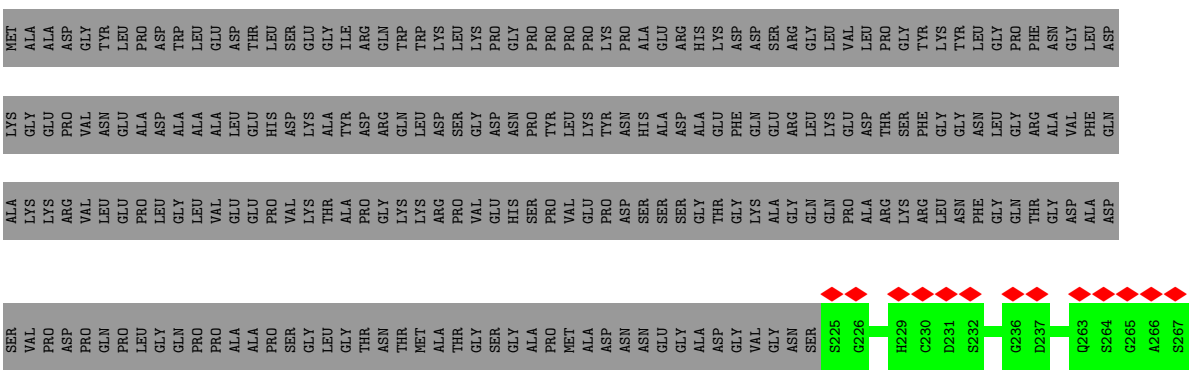
Chain P:

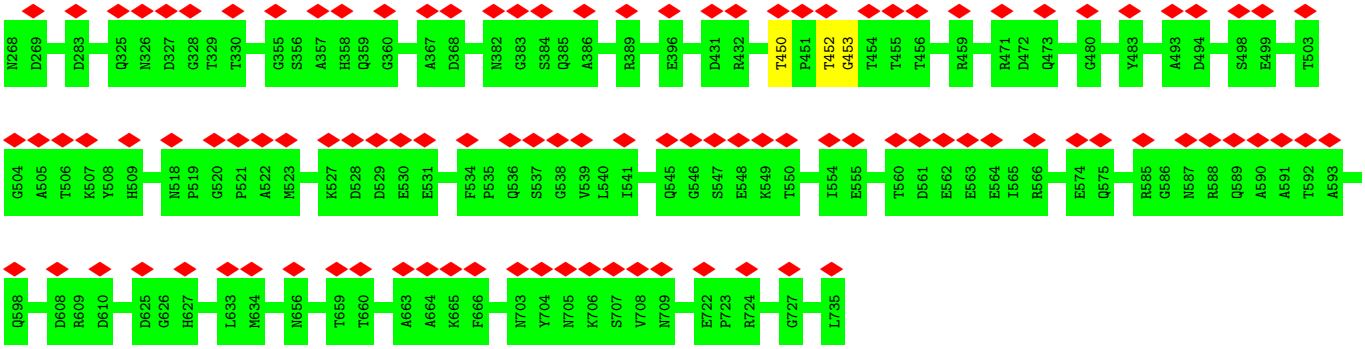




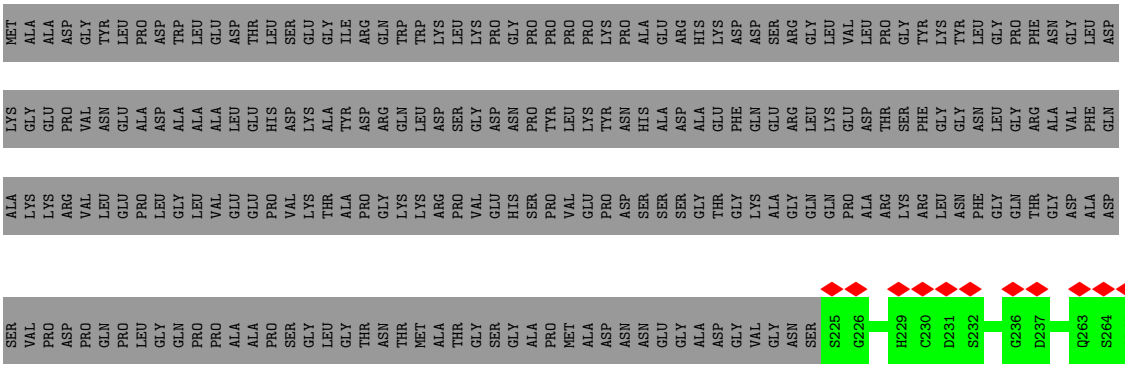
• Molecule 1: Capsid protein VP1

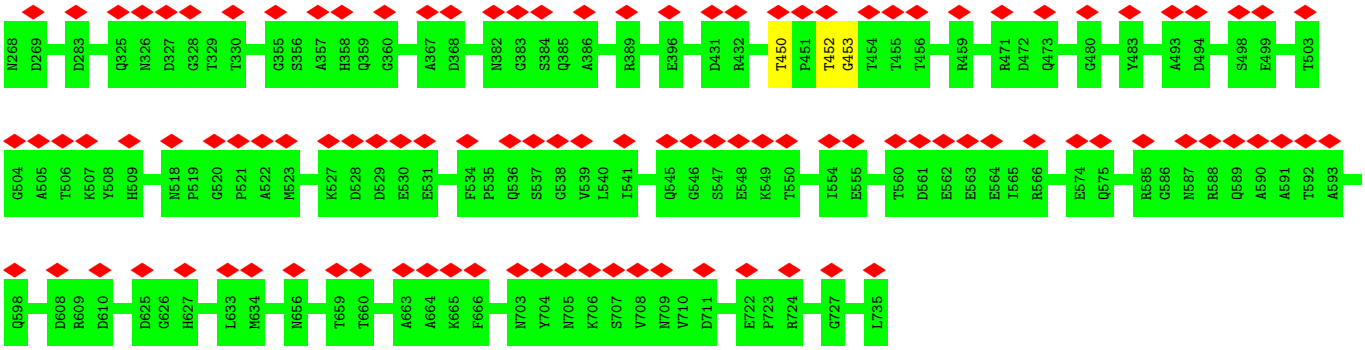
Chain Q:



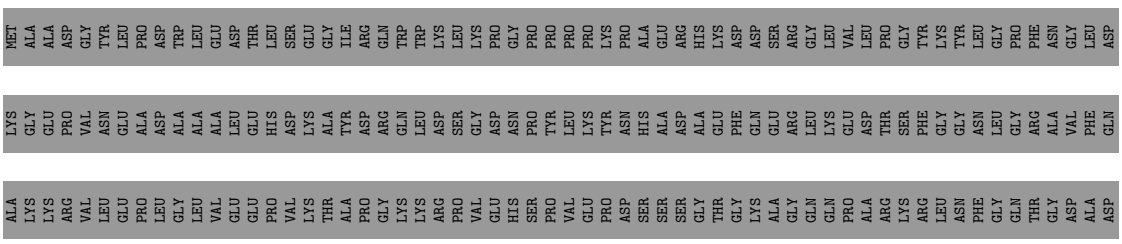


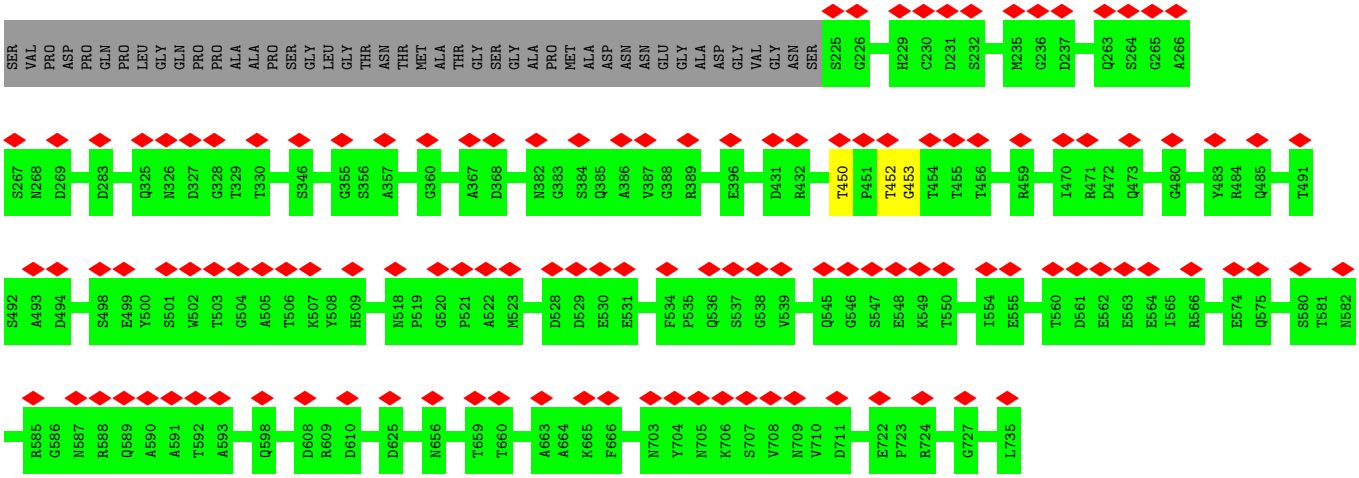
• Molecule 1: Capsid protein VP1





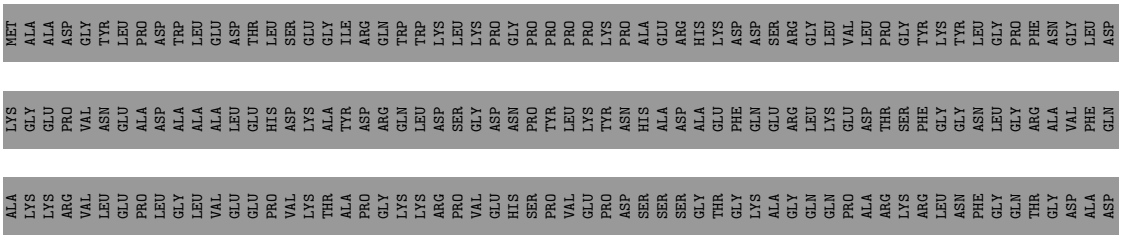
• Molecule 1: Capsid protein VP1

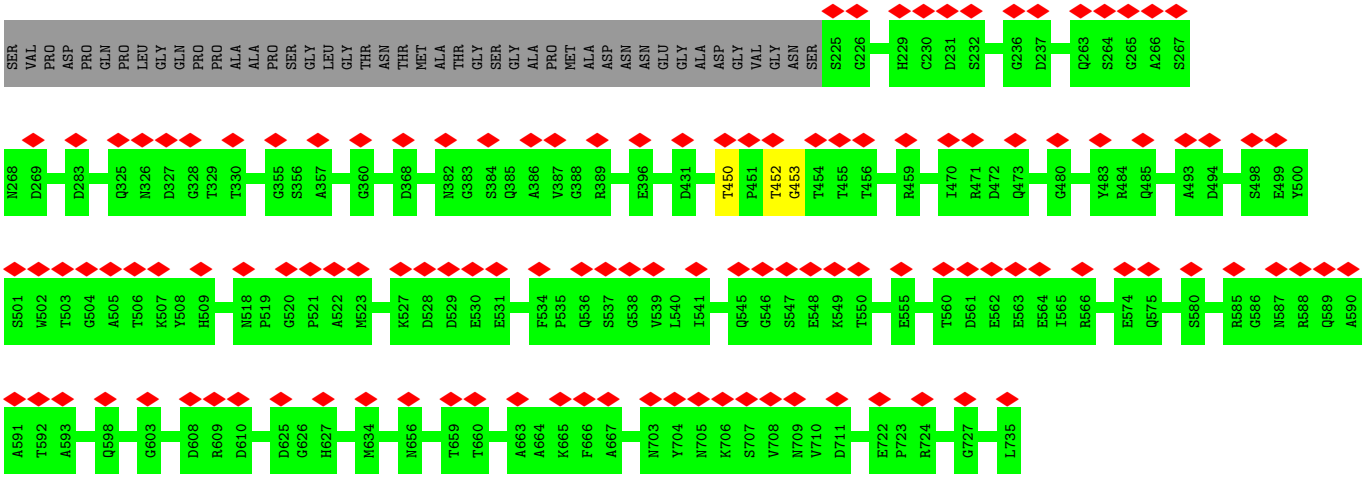


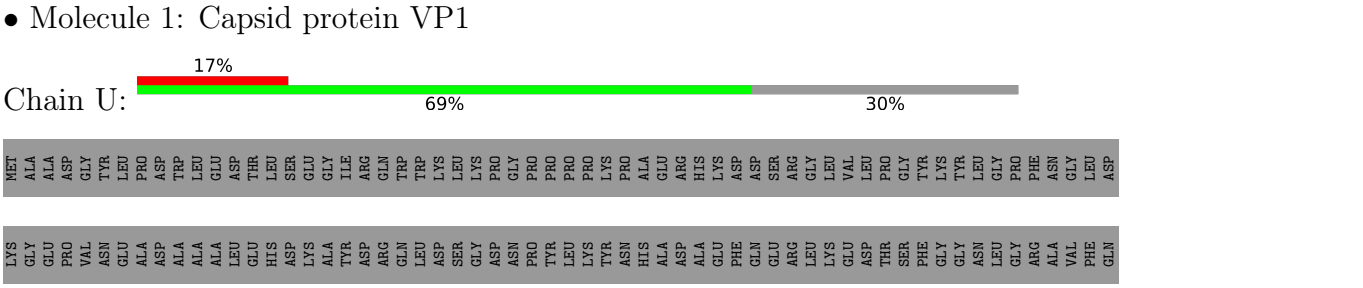


• Molecule 1: Capsid protein VP1

Chain T:

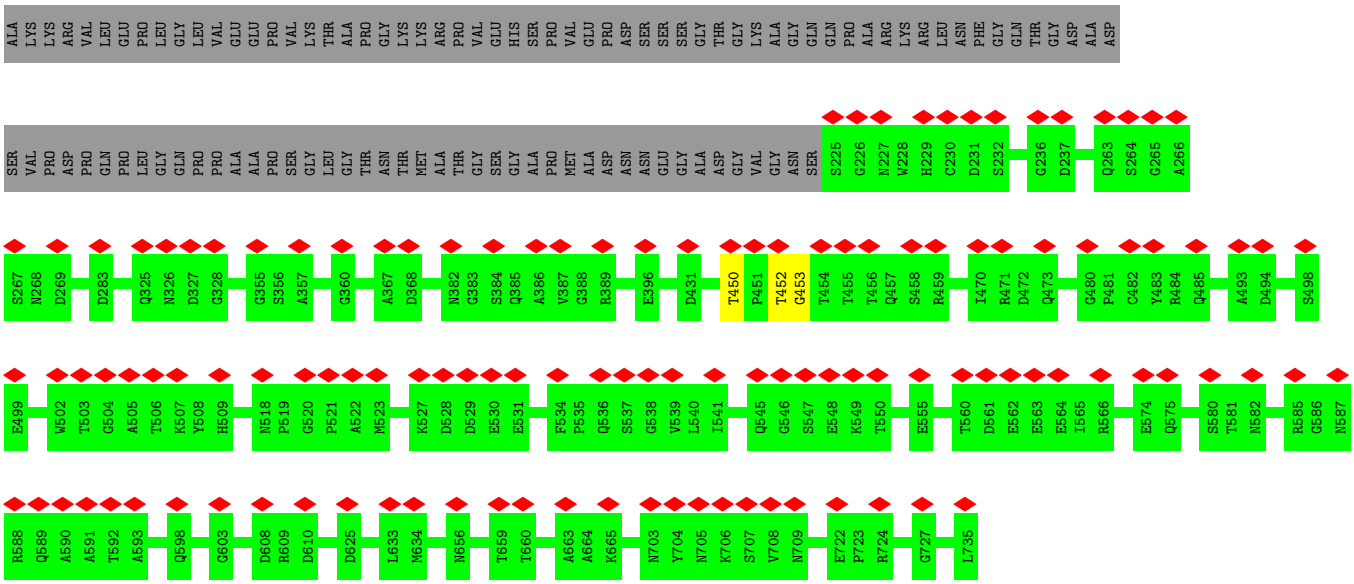






• Molecule 1: Capsid protein VP1

Chain U:



• Molecule 1: Capsid protein VP1

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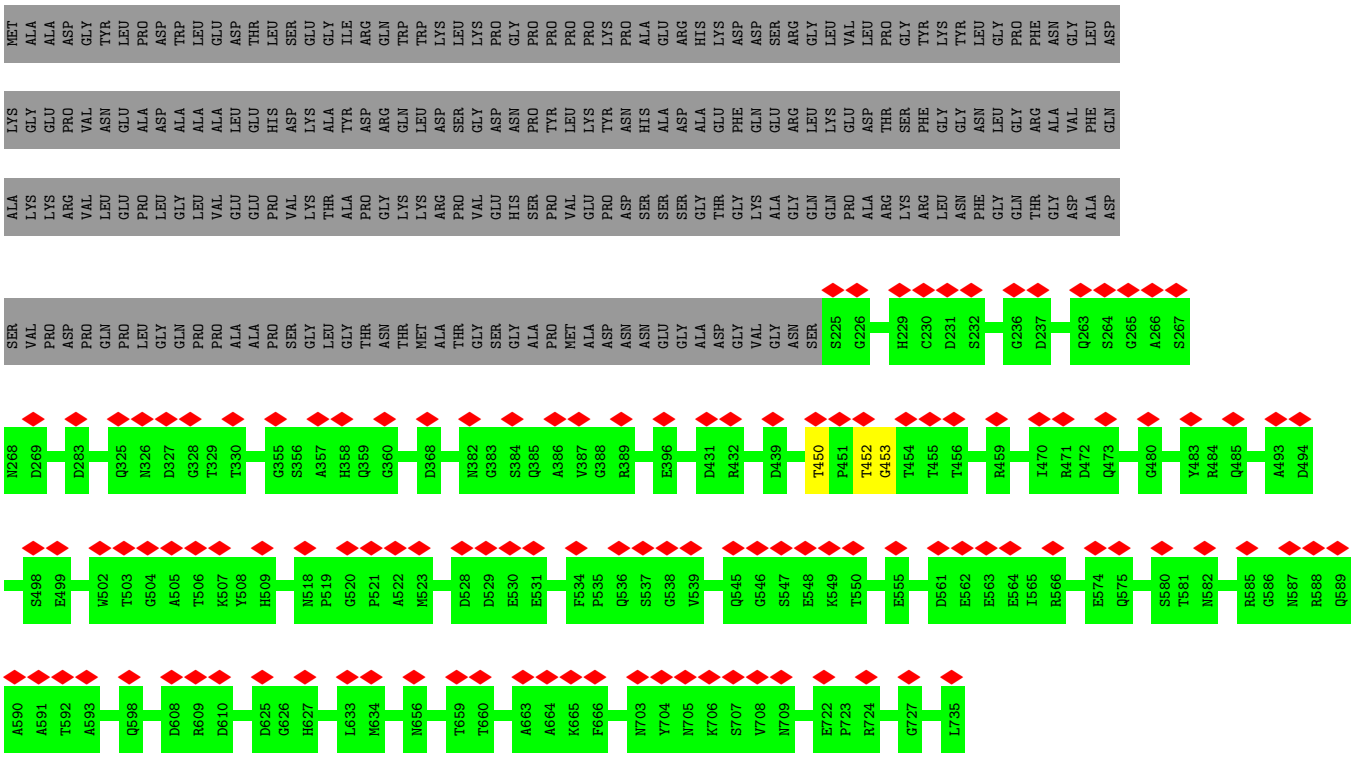
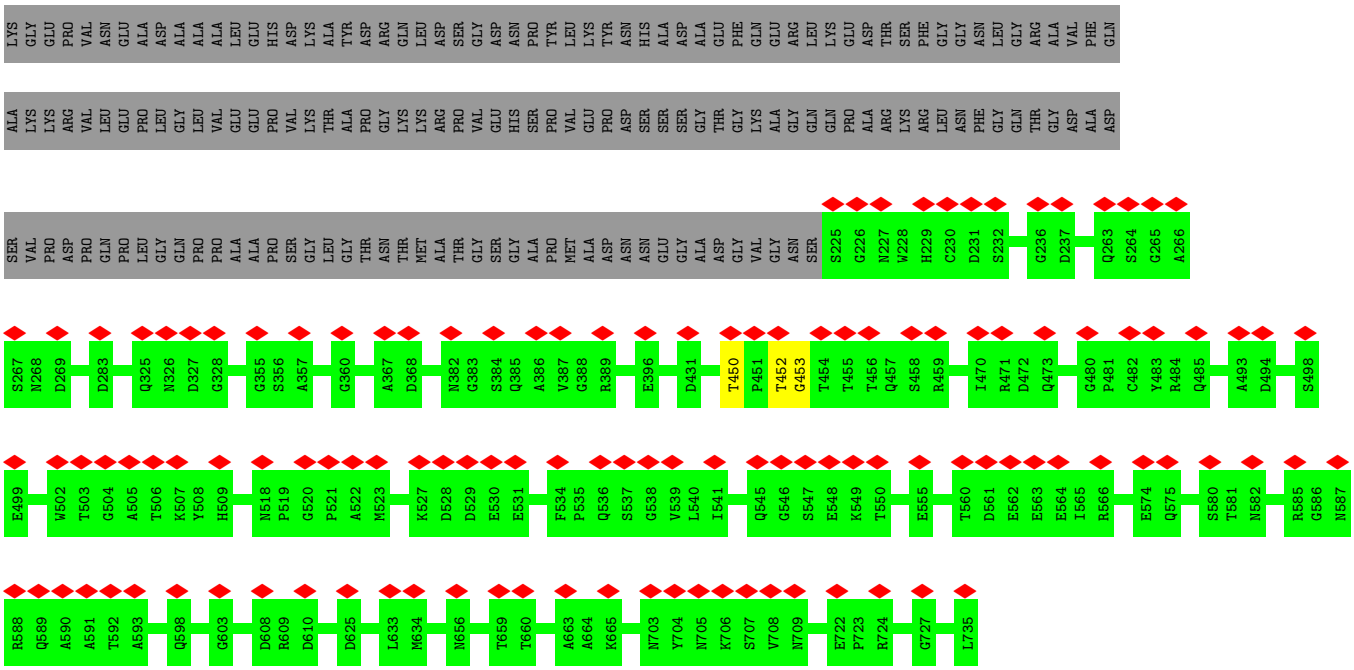
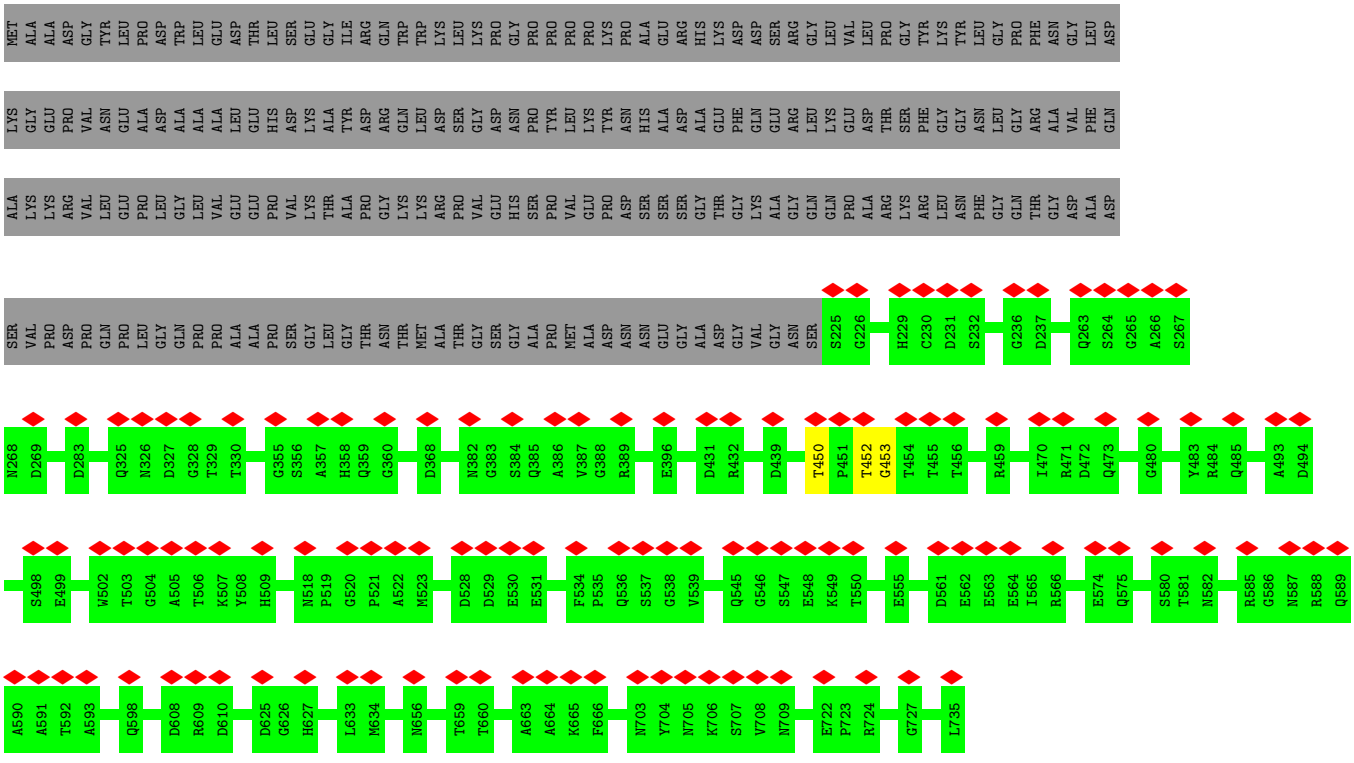


• Molecule 1: Capsid protein VP1

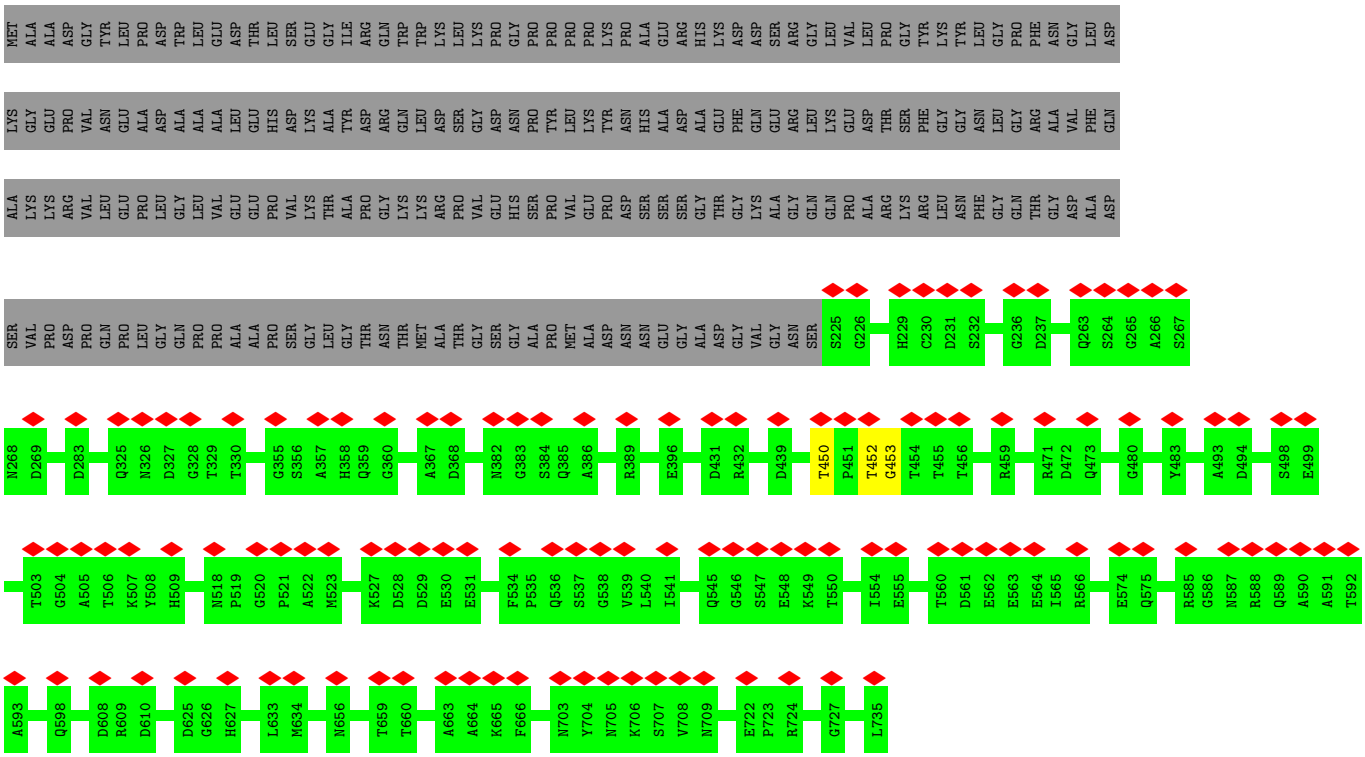
Chain W:



• Molecule 1: Capsid protein VP1

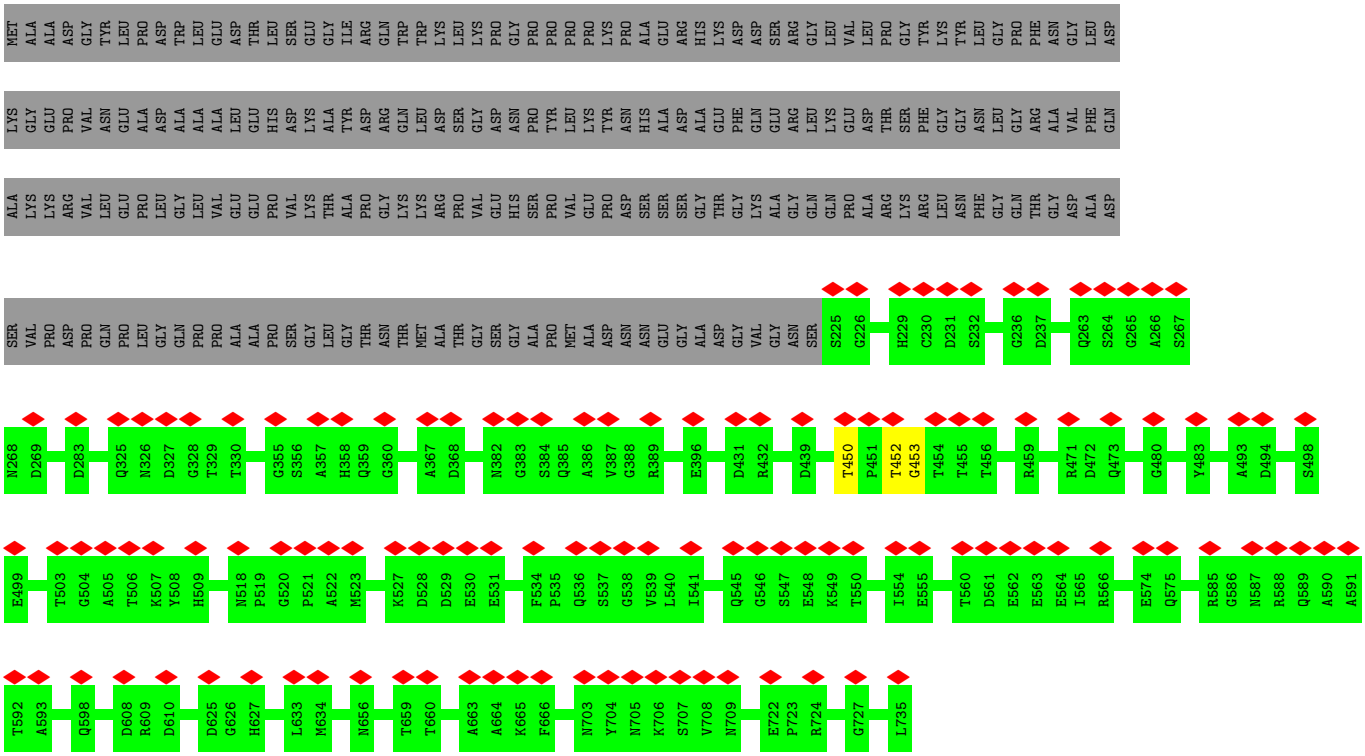


• Molecule 1: Capsid protein VP1



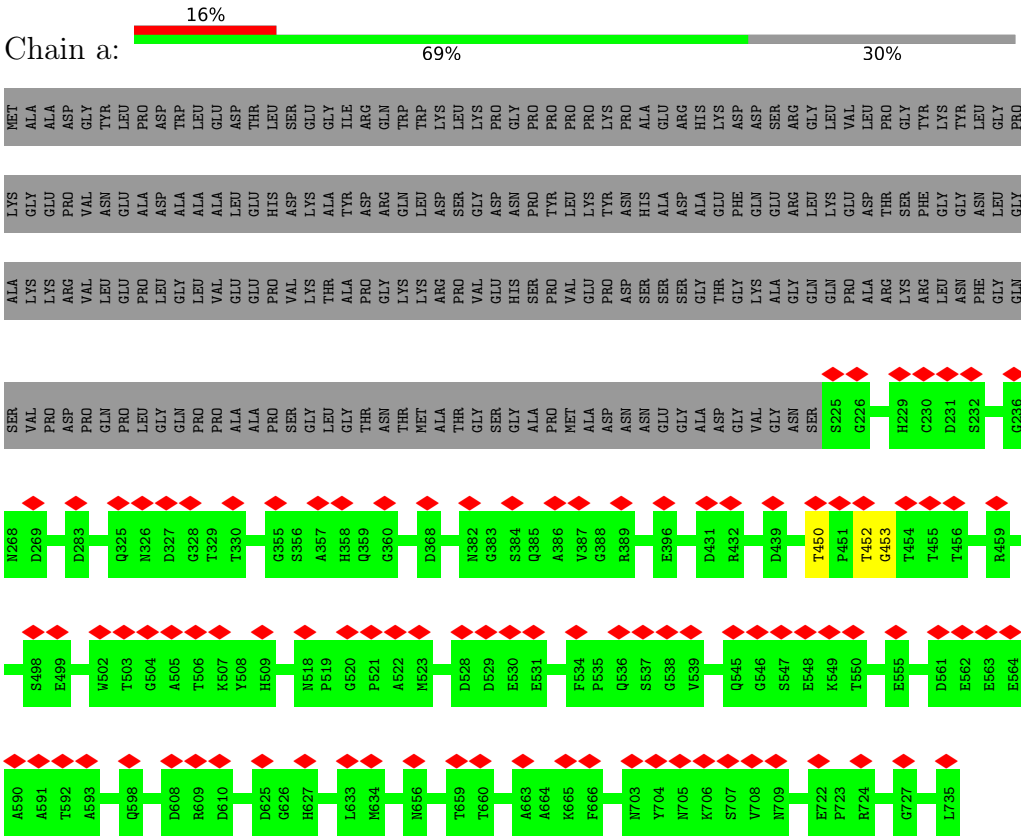
• Molecule 1: Capsid protein VP1

Chain Z:



• Molecule 1: Capsid protein VP1

Chain a:



• Molecule 1: Capsid protein VP1

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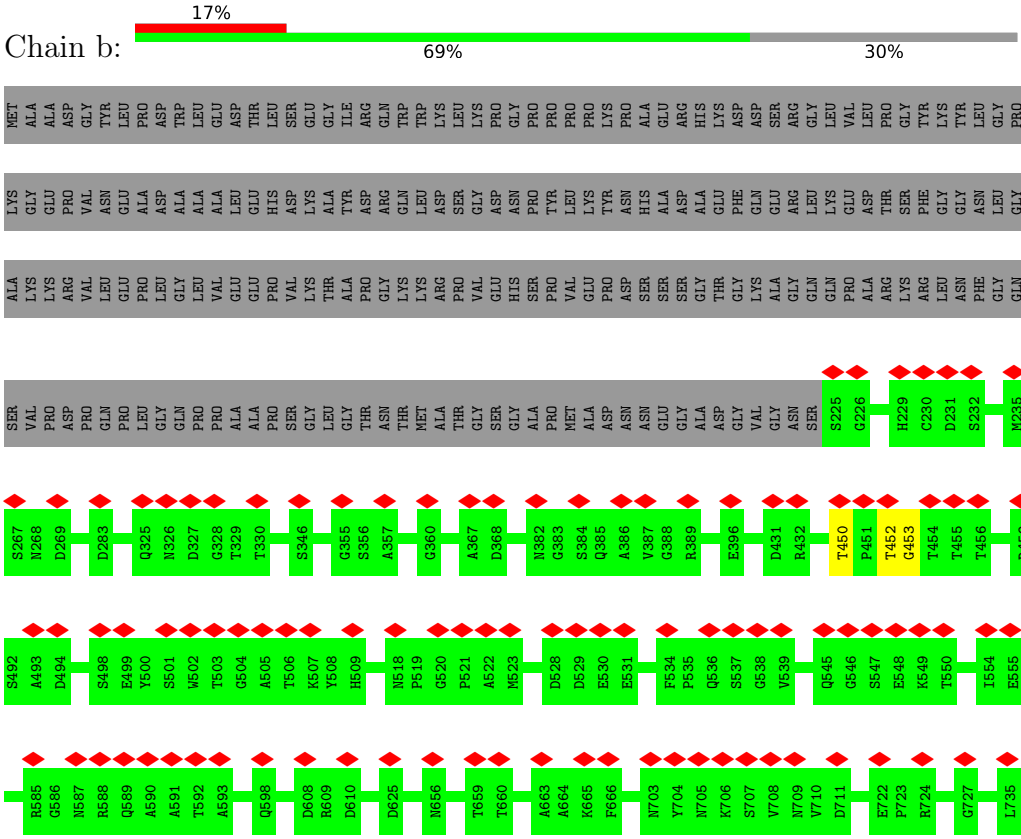
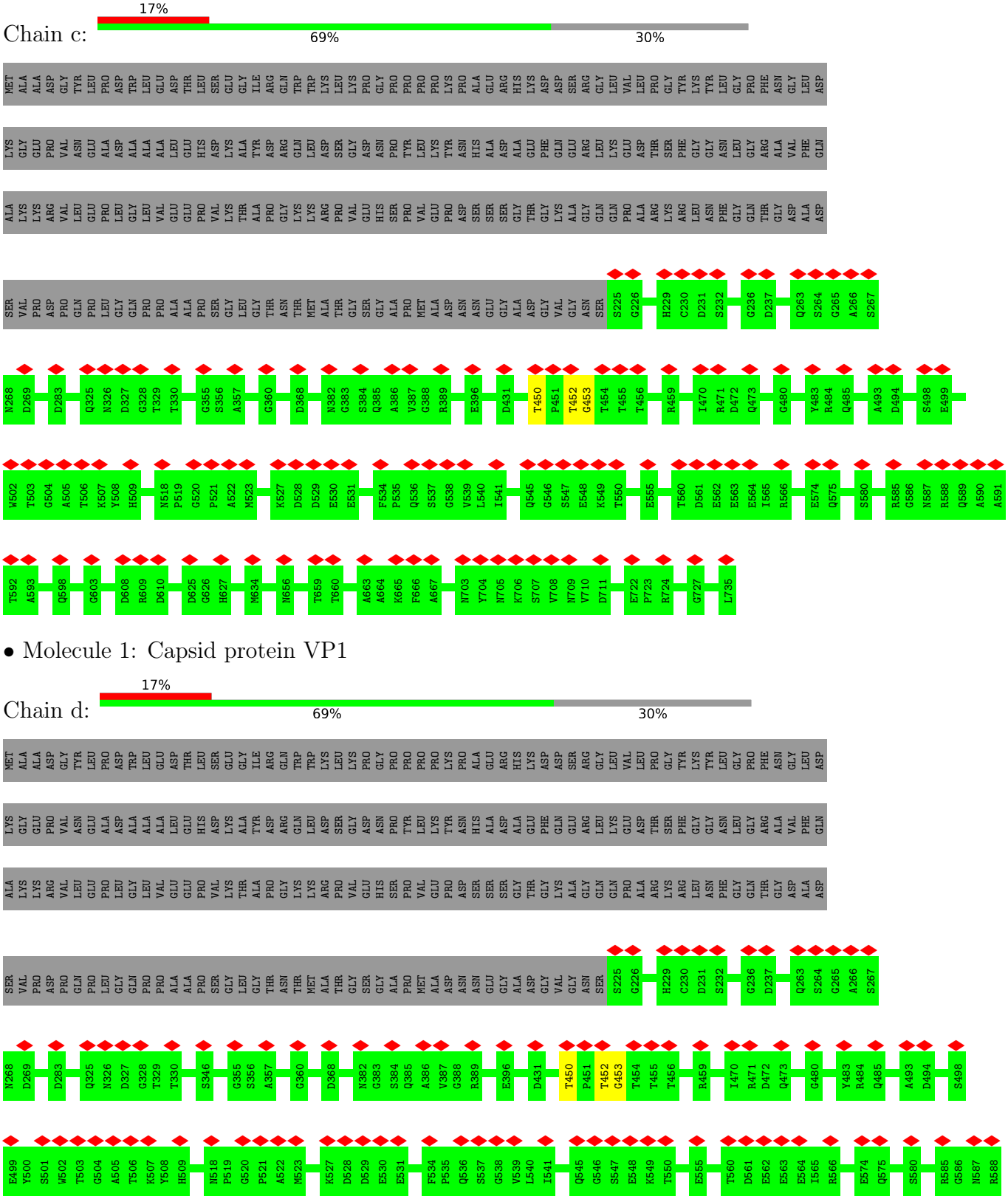


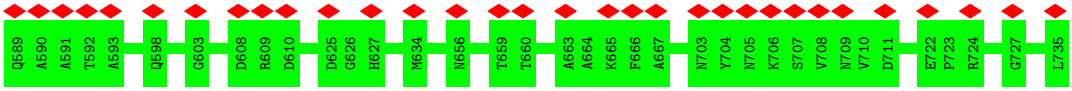
• Molecule 1: Capsid protein VP1

Chain c:



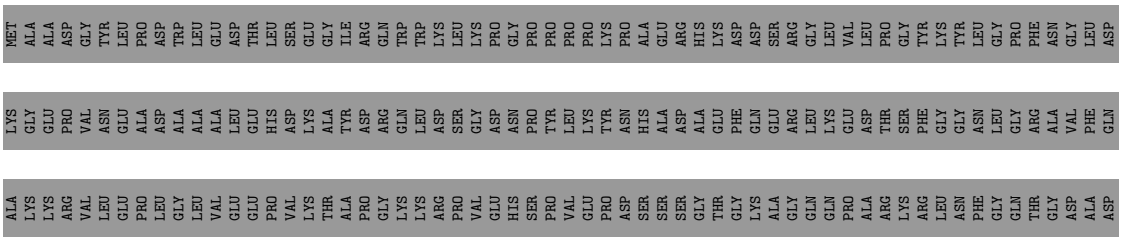
• Molecule 1: Capsid protein VP1

Chain d:



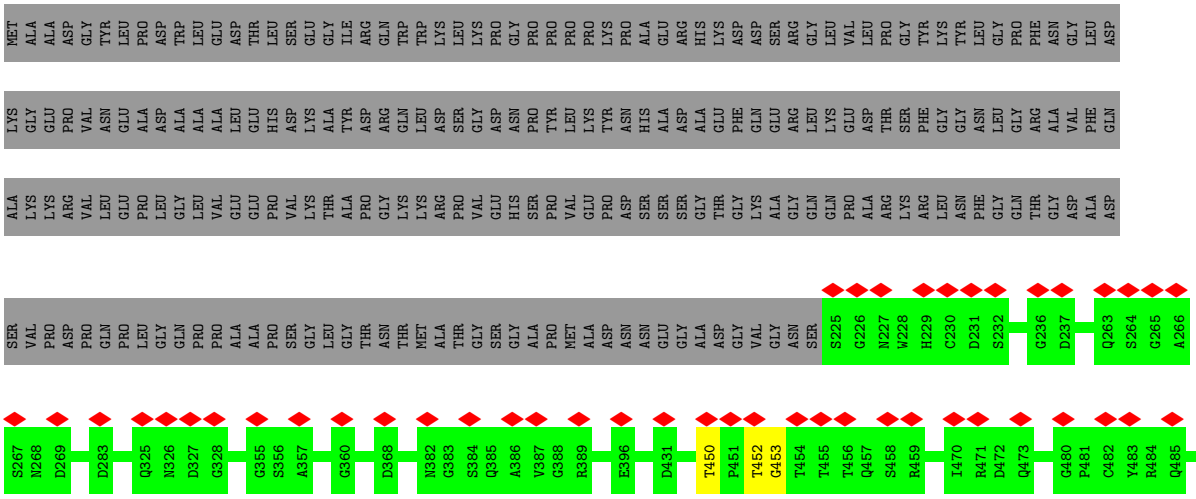
• Molecule 1: Capsid protein VP1

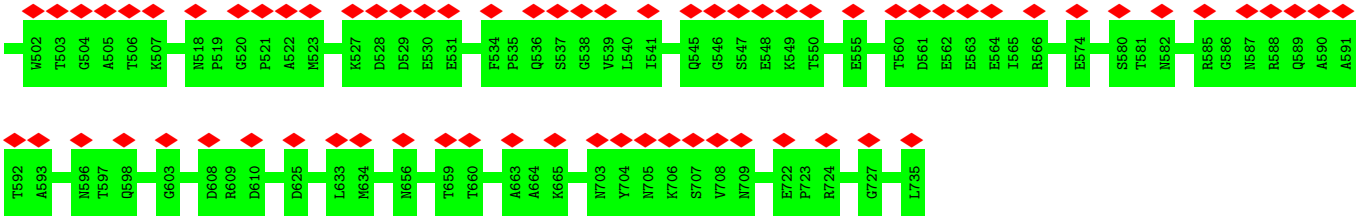
Chain e:



• Molecule 1: Capsid protein VP1

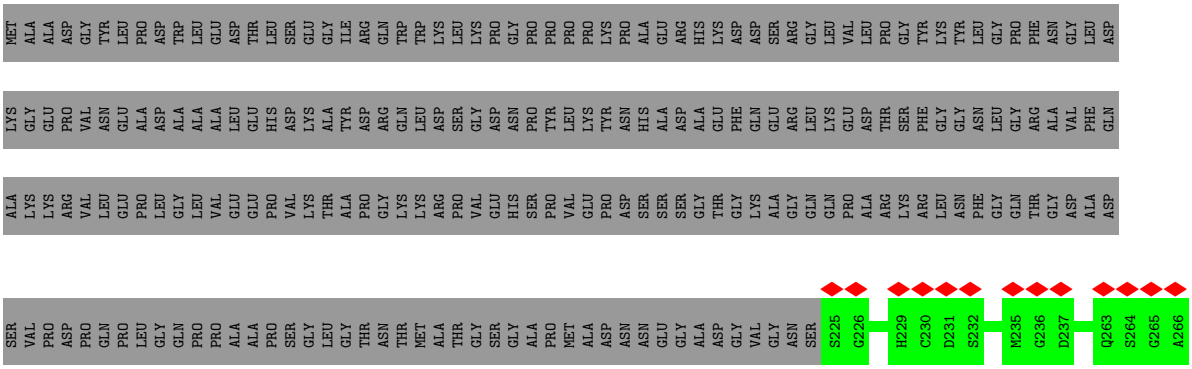
Chain f:





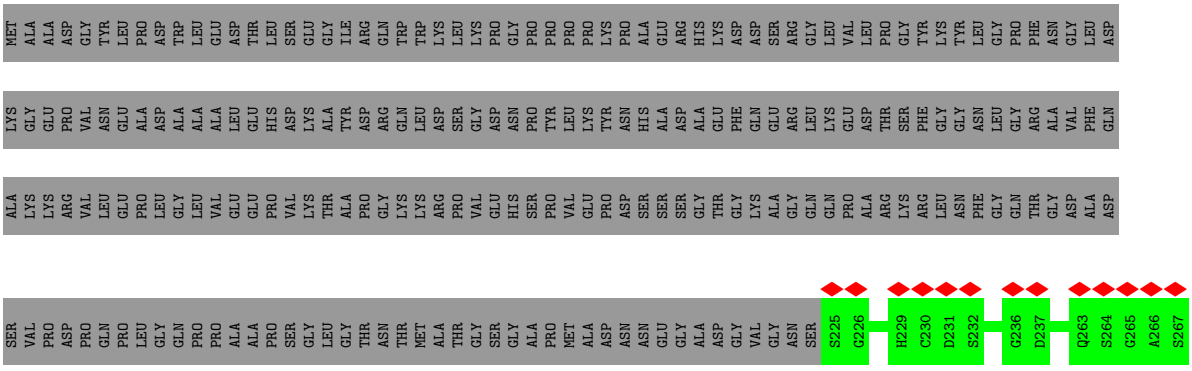
• Molecule 1: Capsid protein VP1

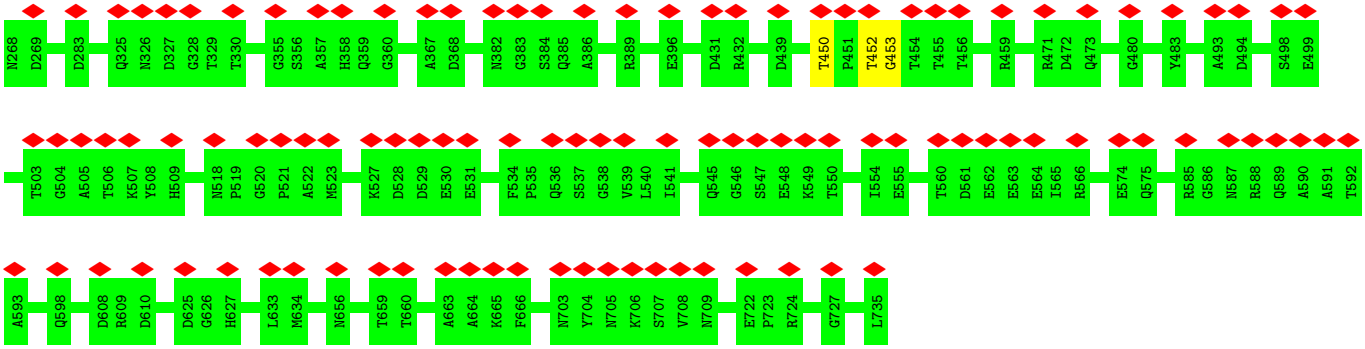
Chain g:



• Molecule 1: Capsid protein VP1

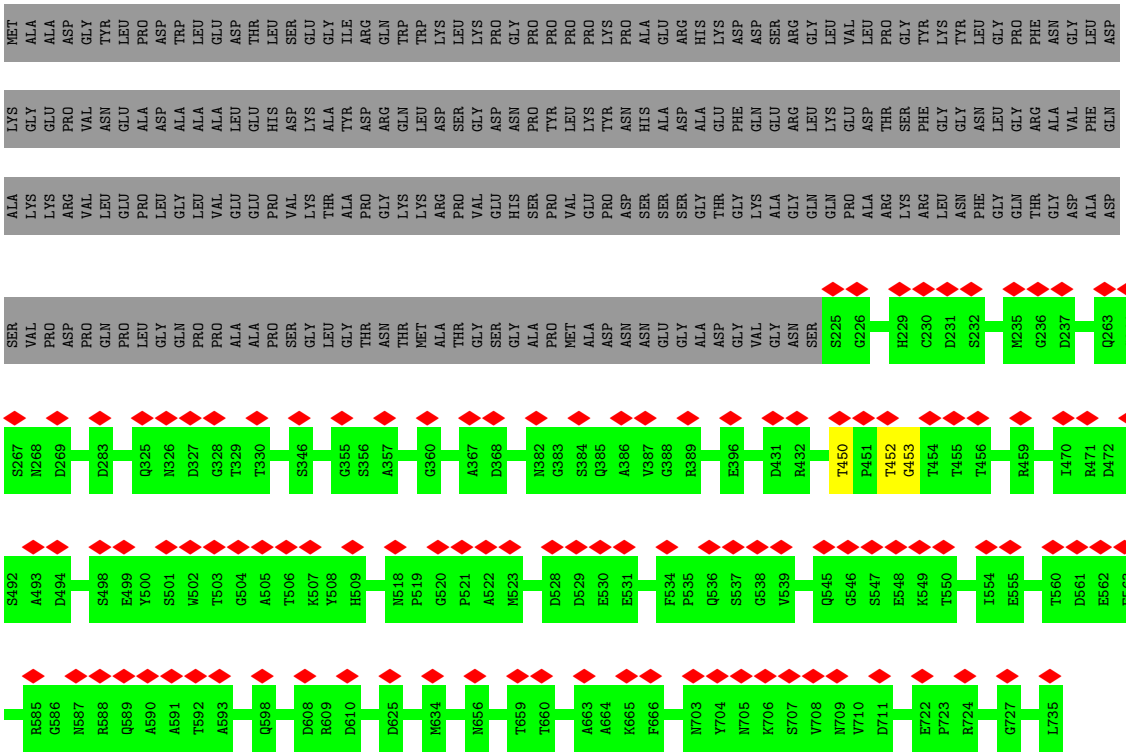
Chain h:





• Molecule 1: Capsid protein VP1

Chain i:



• Molecule 1: Capsid protein VP1

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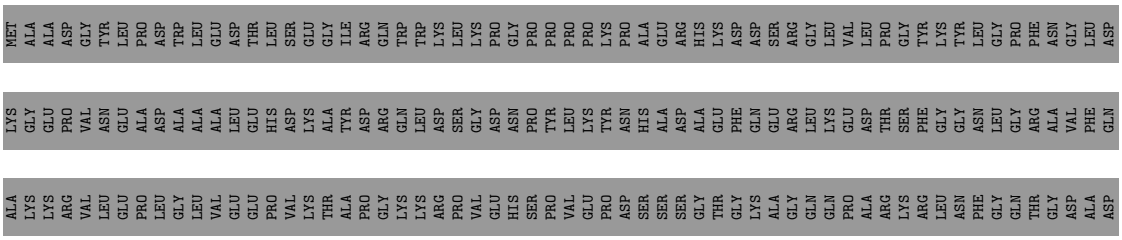
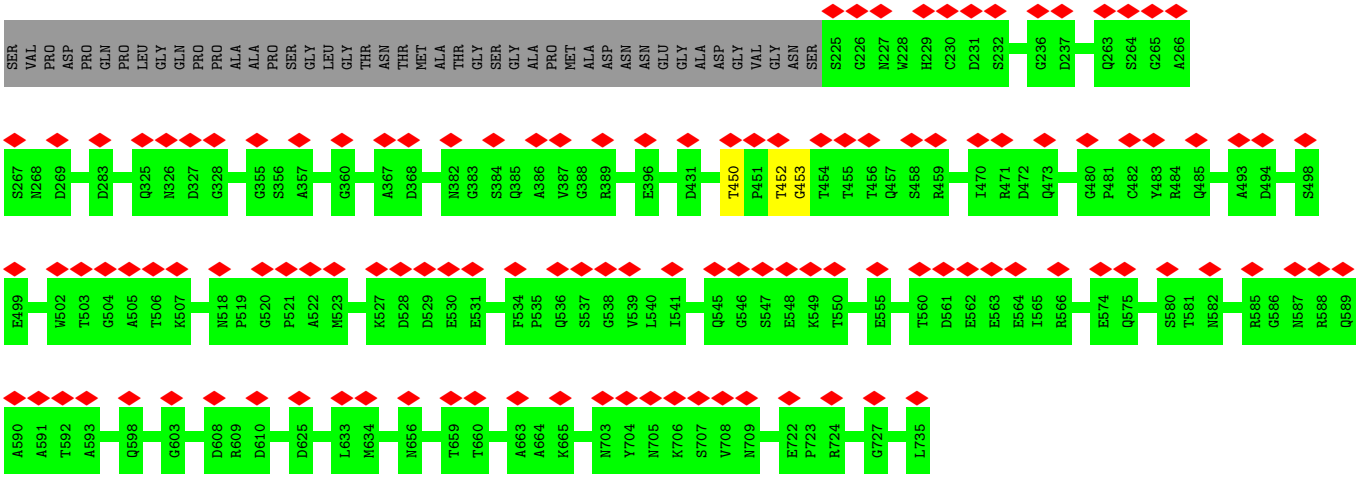
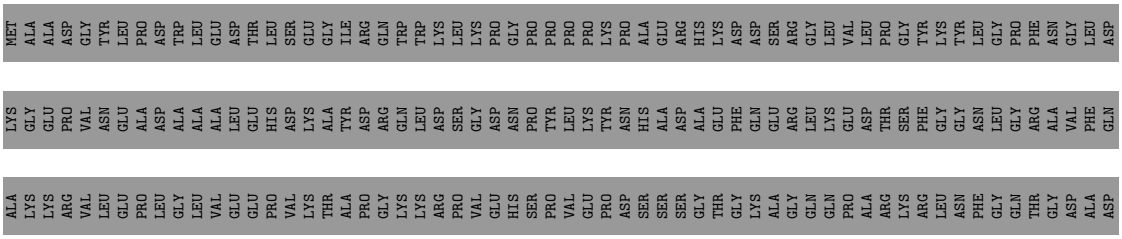
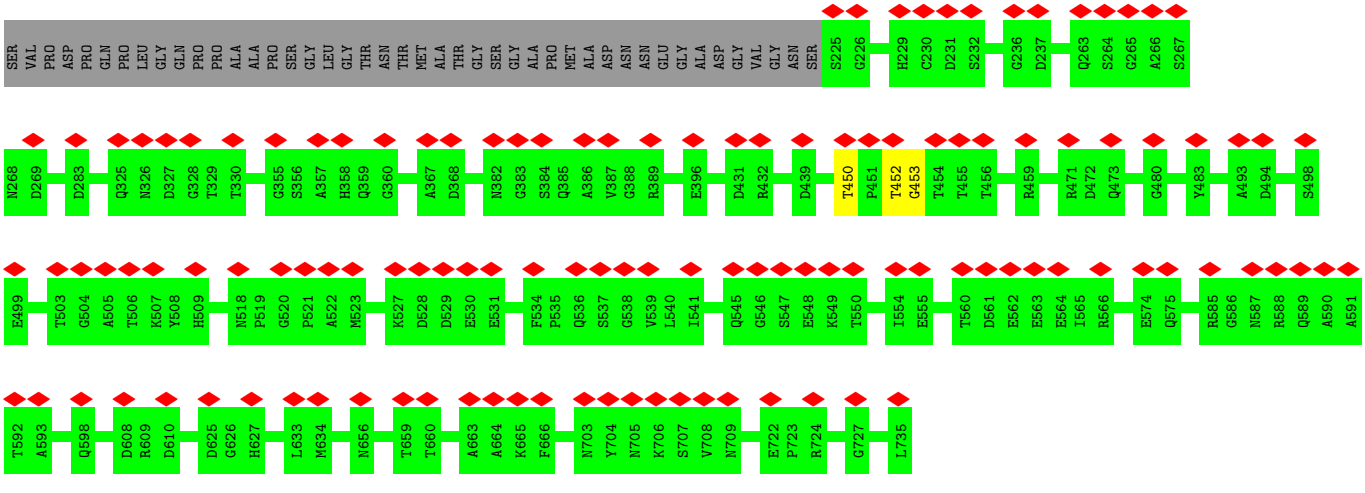




• Molecule 1: Capsid protein VP1

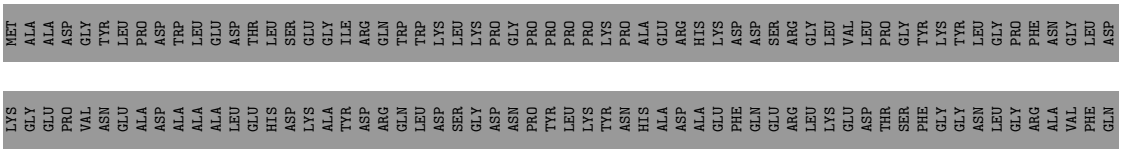
Chain k:

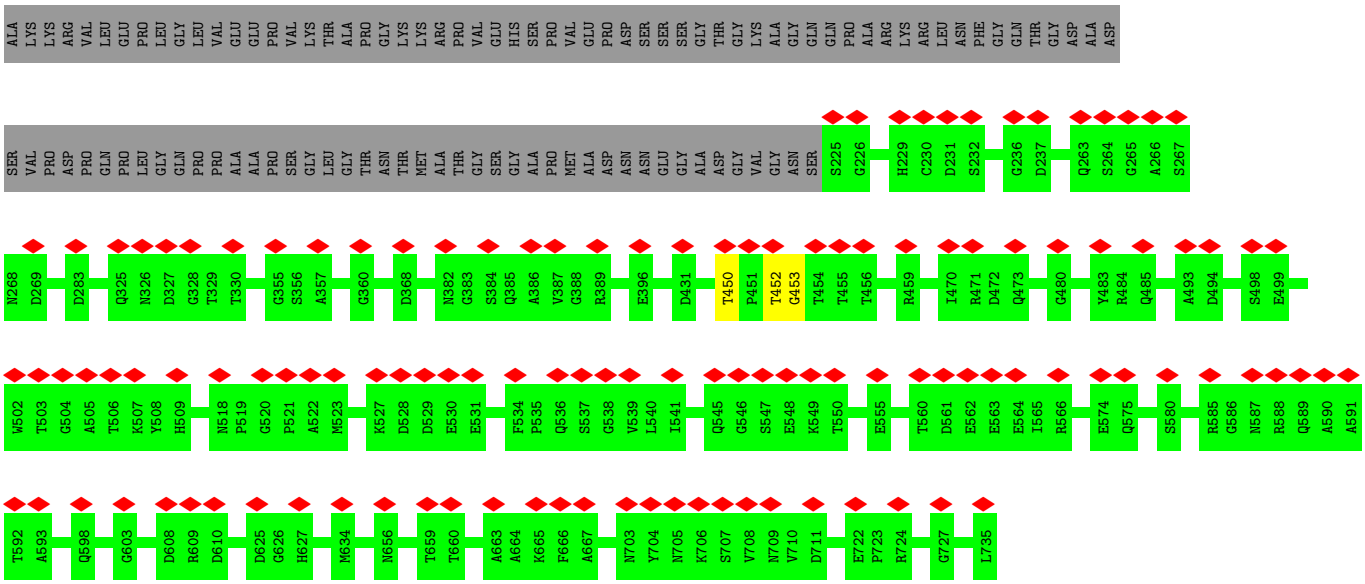




• Molecule 1: Capsid protein VP1

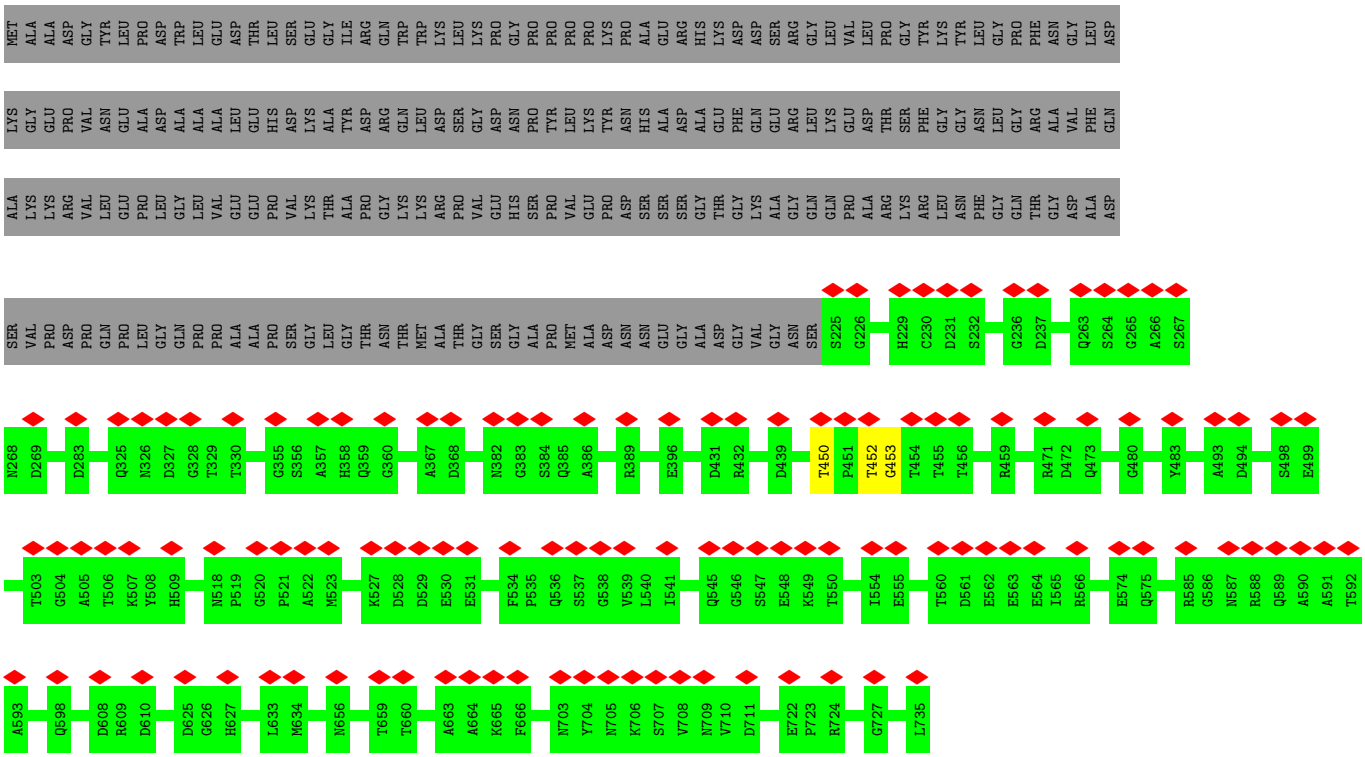
Chain l:





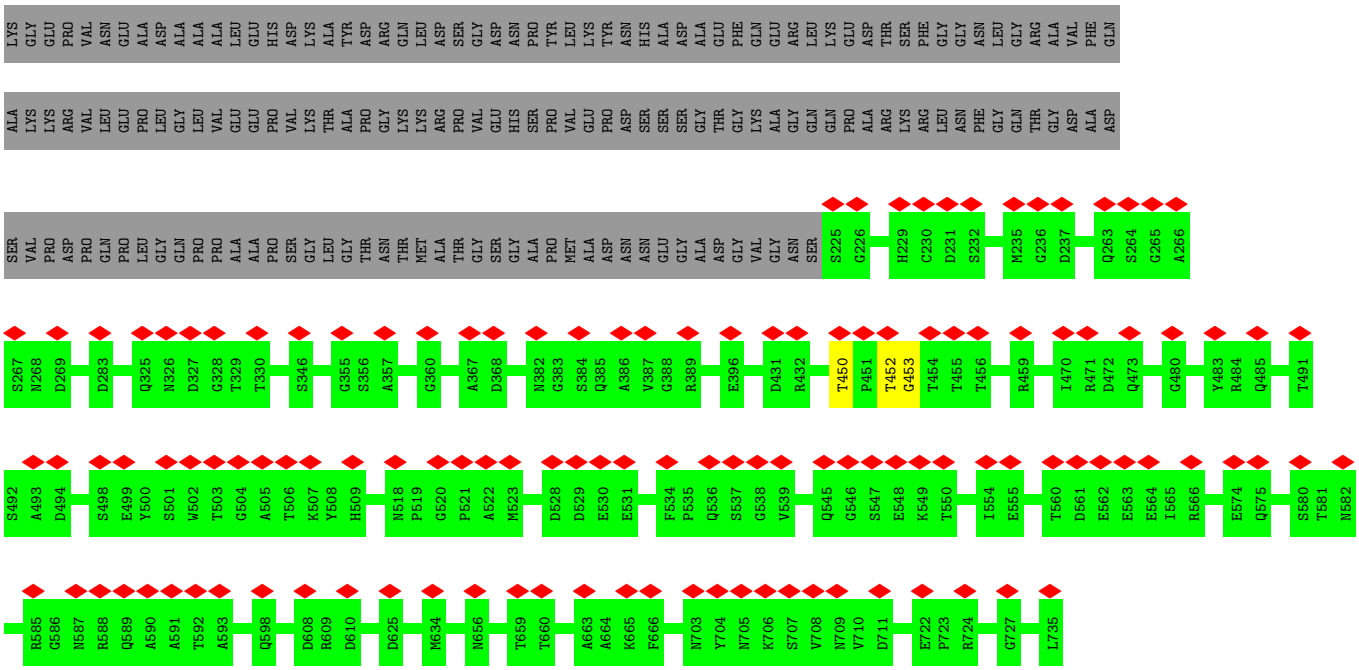
• Molecule 1: Capsid protein VP1

Chain m:

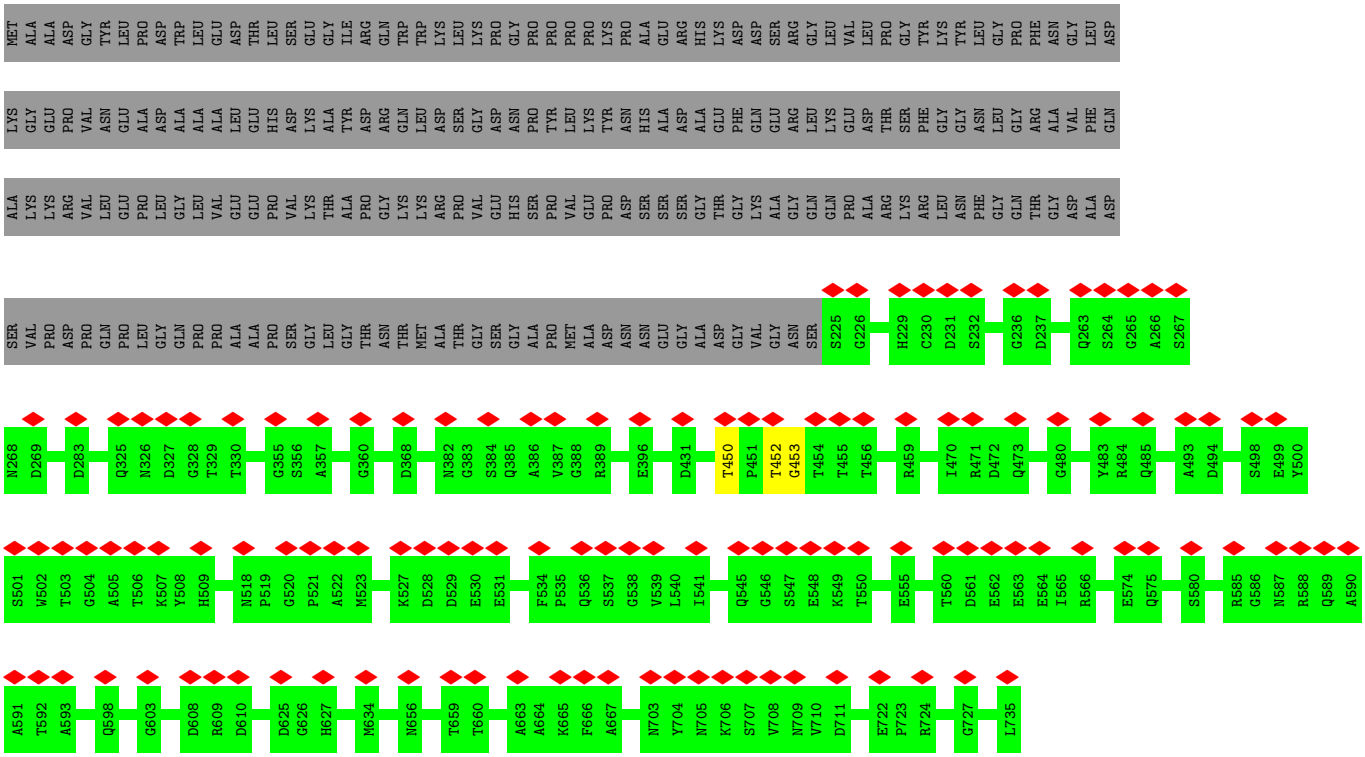


• Molecule 1: Capsid protein VP1

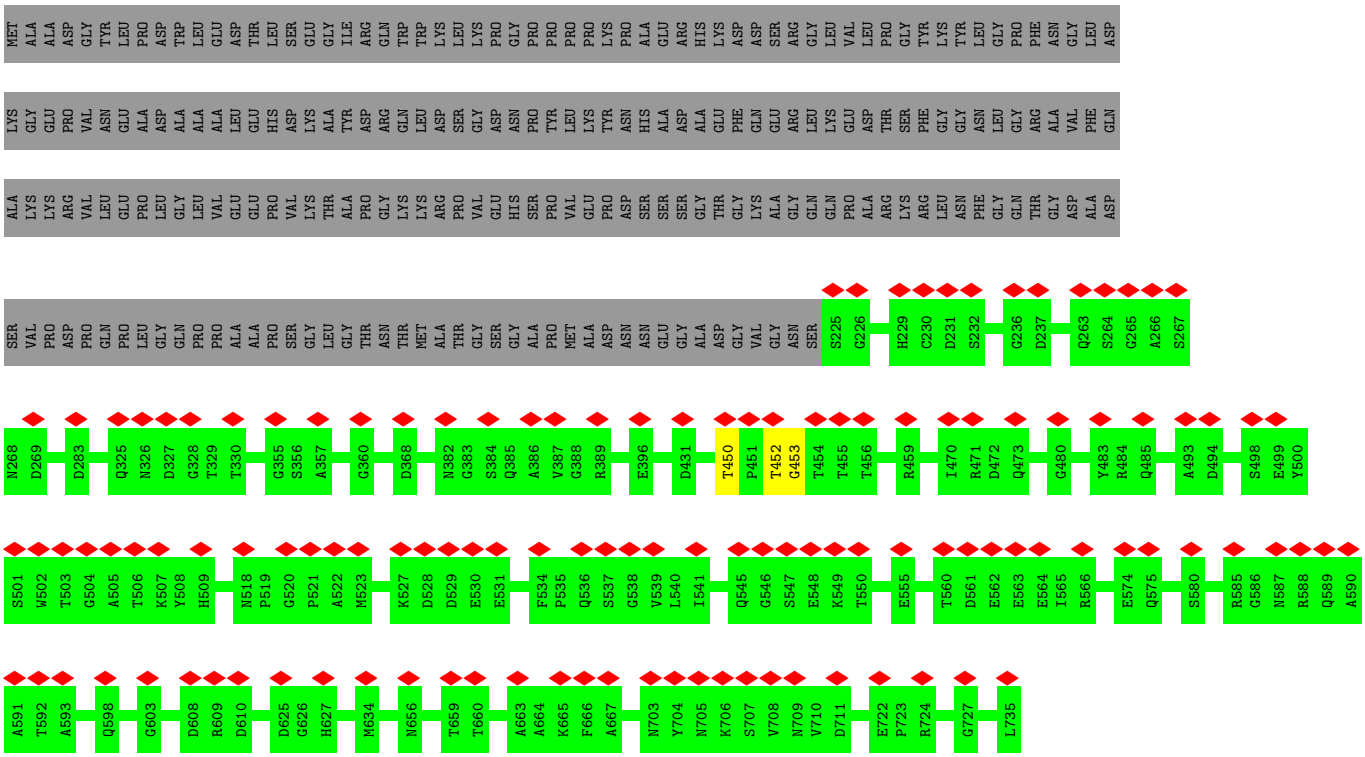
Chain n:



• Molecule 1: Capsid protein VP1

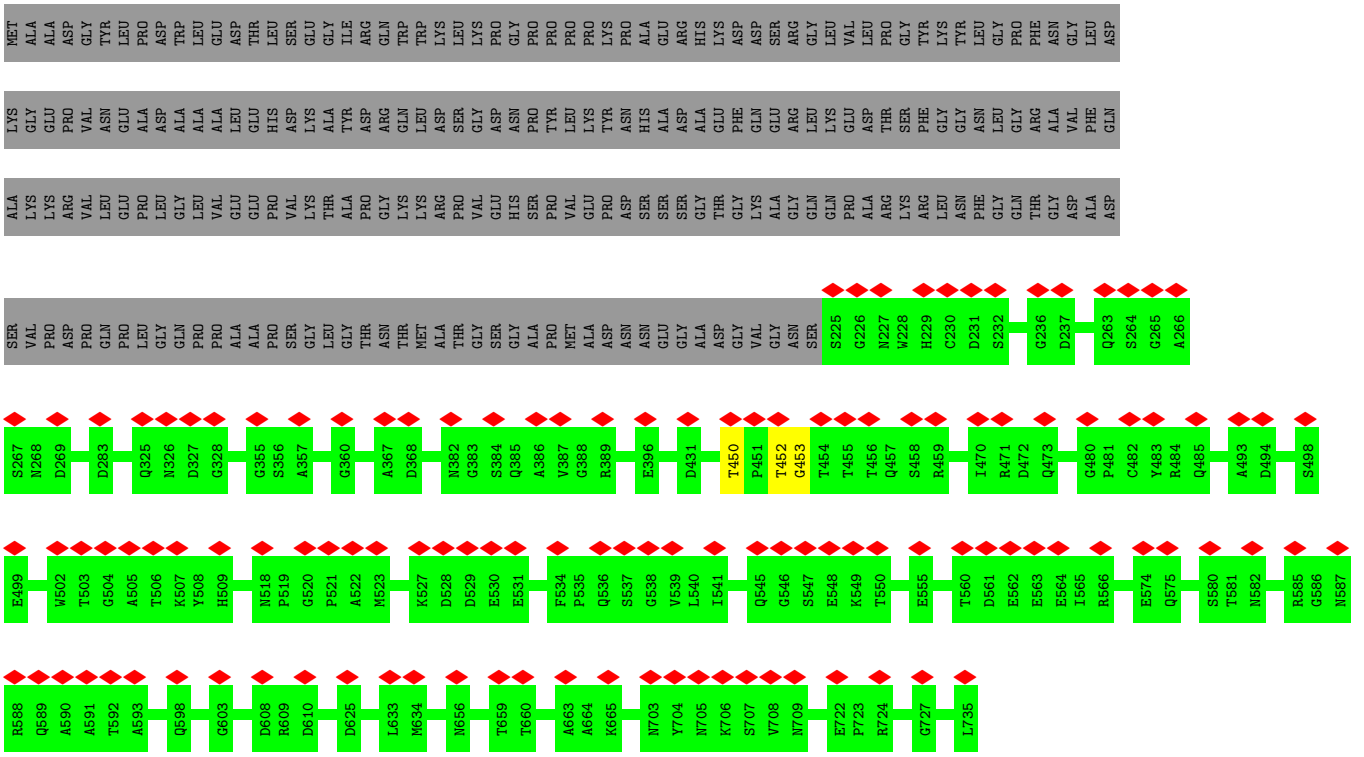


• Molecule 1: Capsid protein VP1



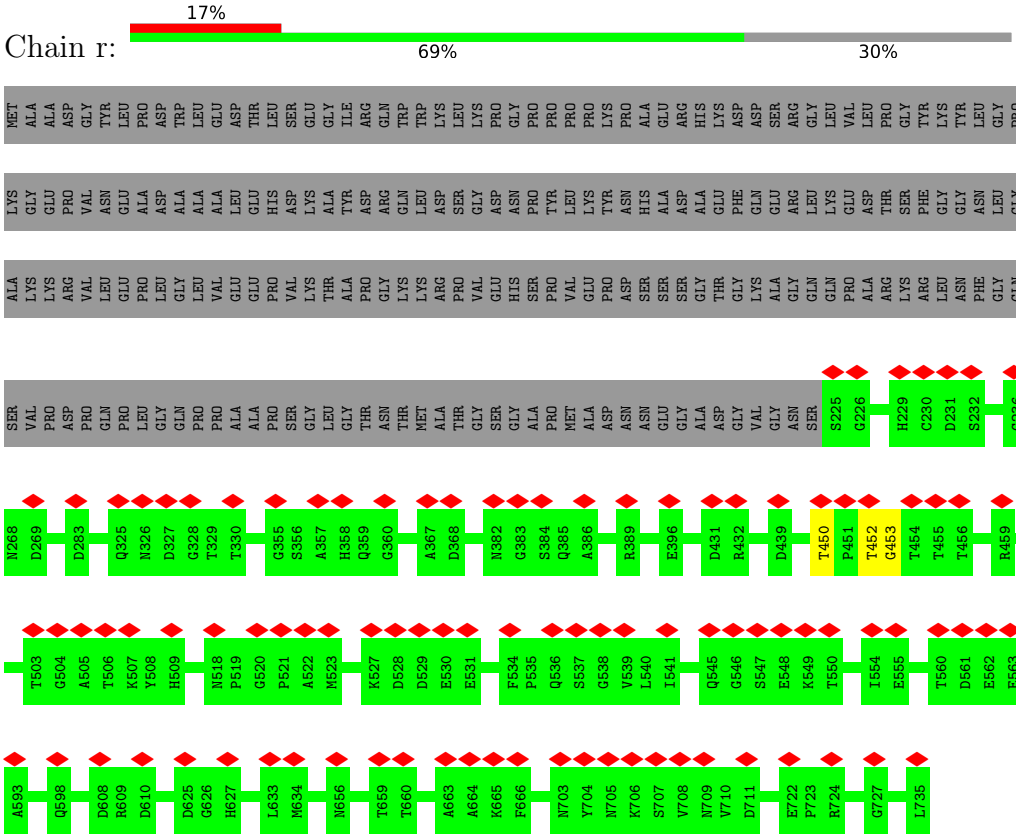
• Molecule 1: Capsid protein VP1

Chain q:



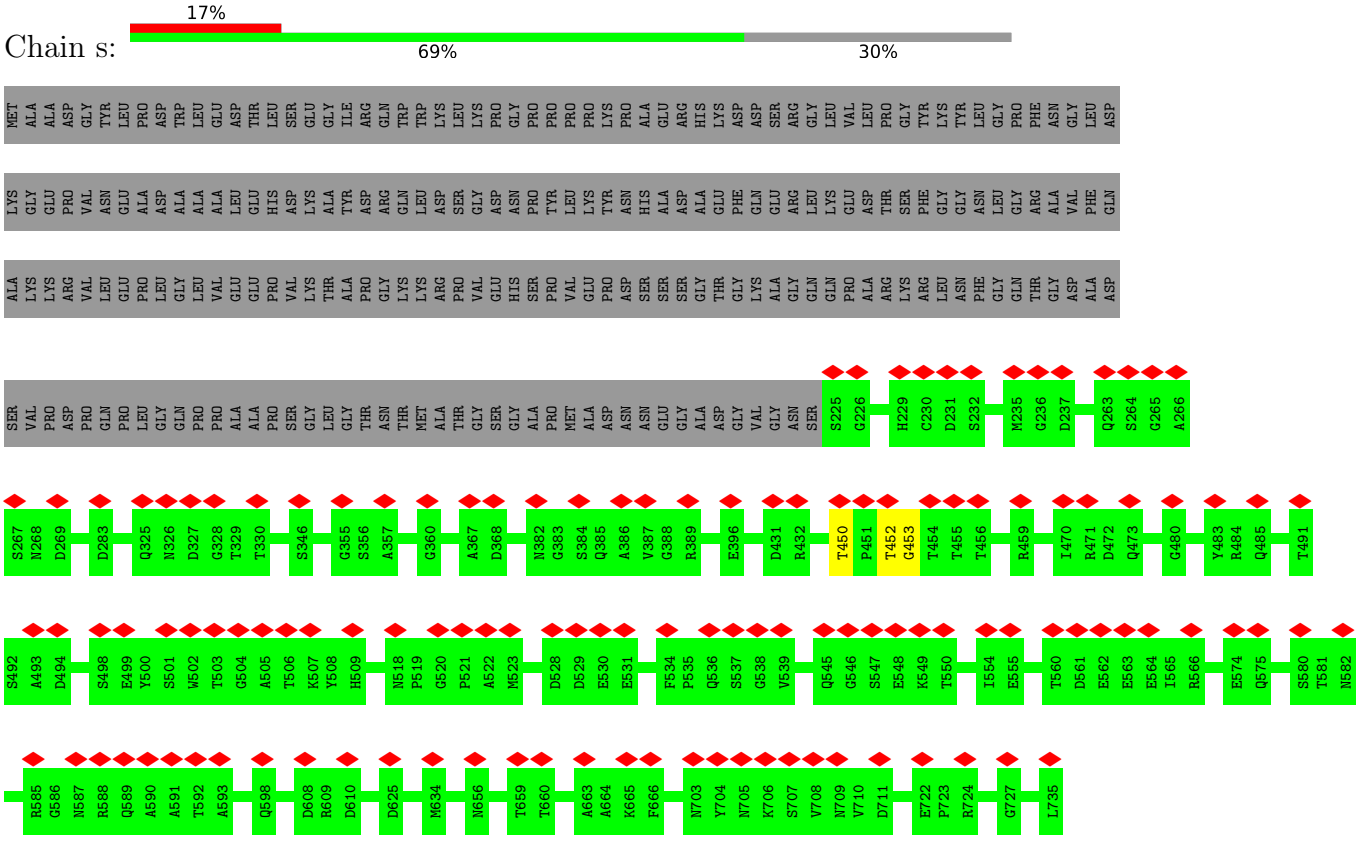
• Molecule 1: Capsid protein VP1

Chain r:



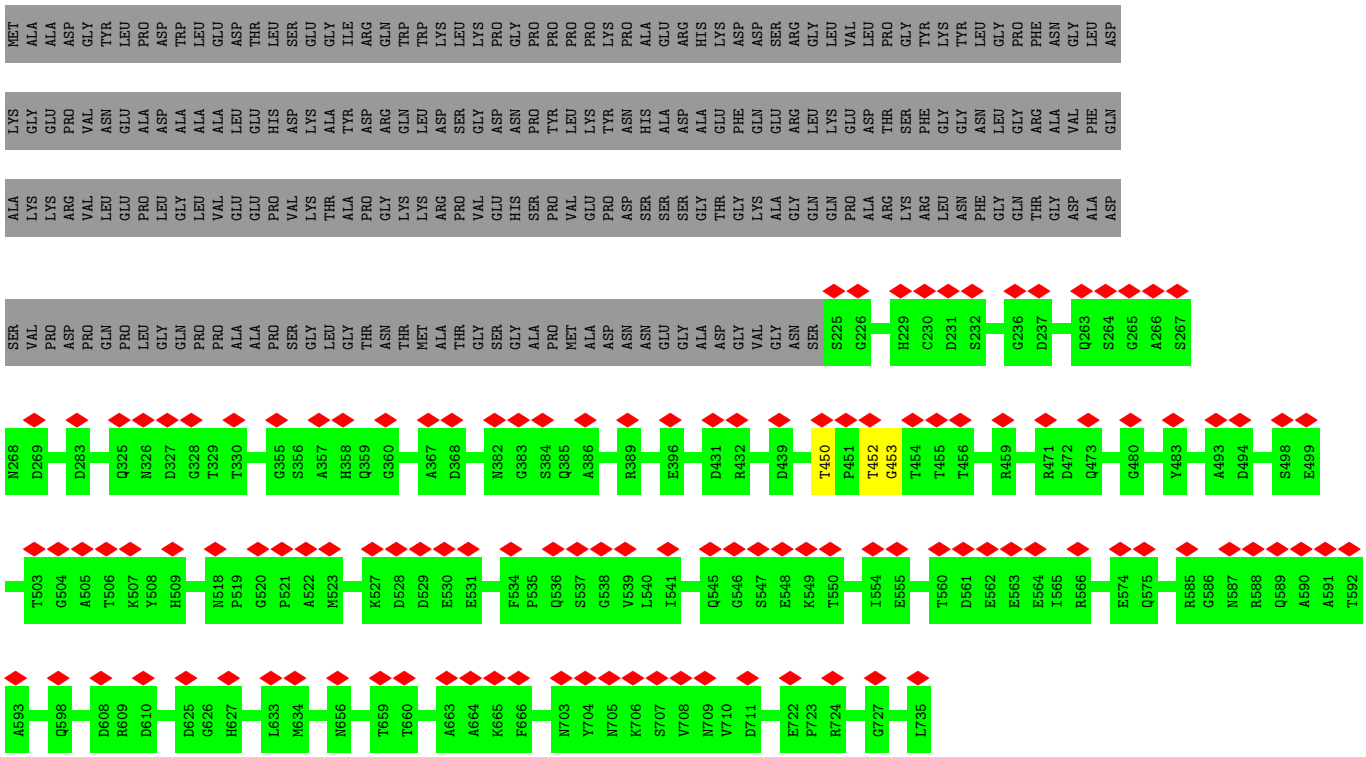
• Molecule 1: Capsid protein VP1

Chain s:



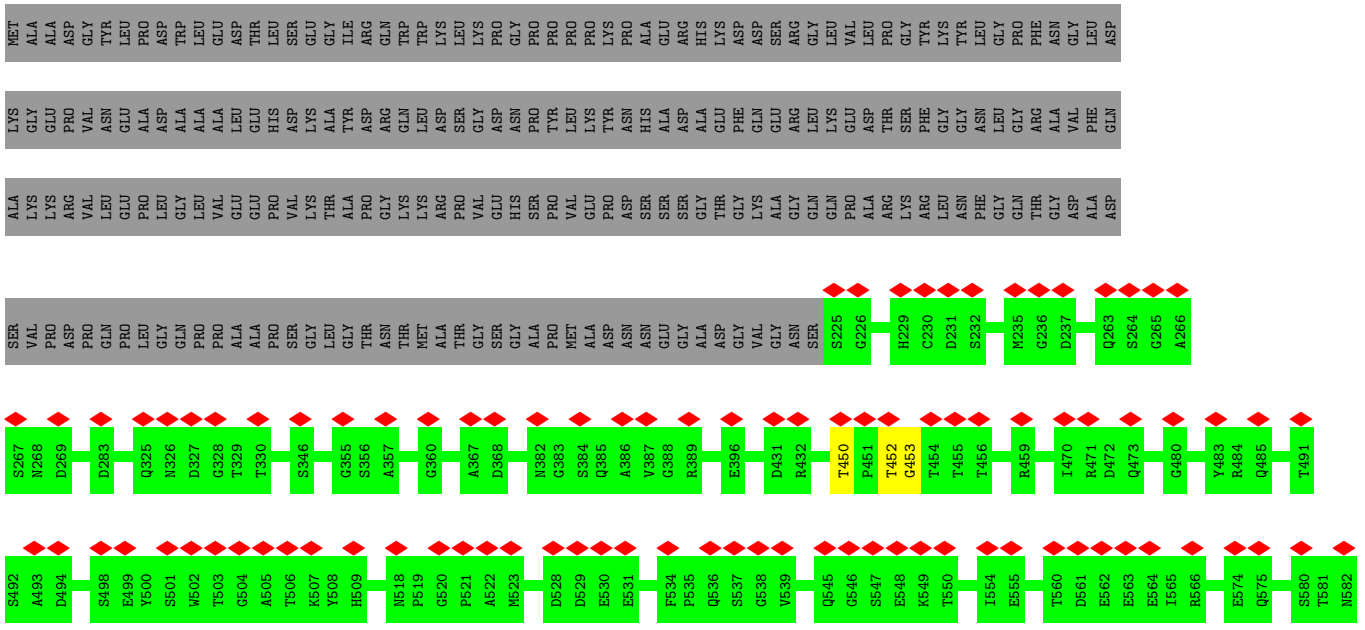
• Molecule 1: Capsid protein VP1

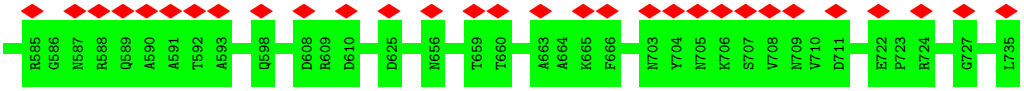
Chain t:



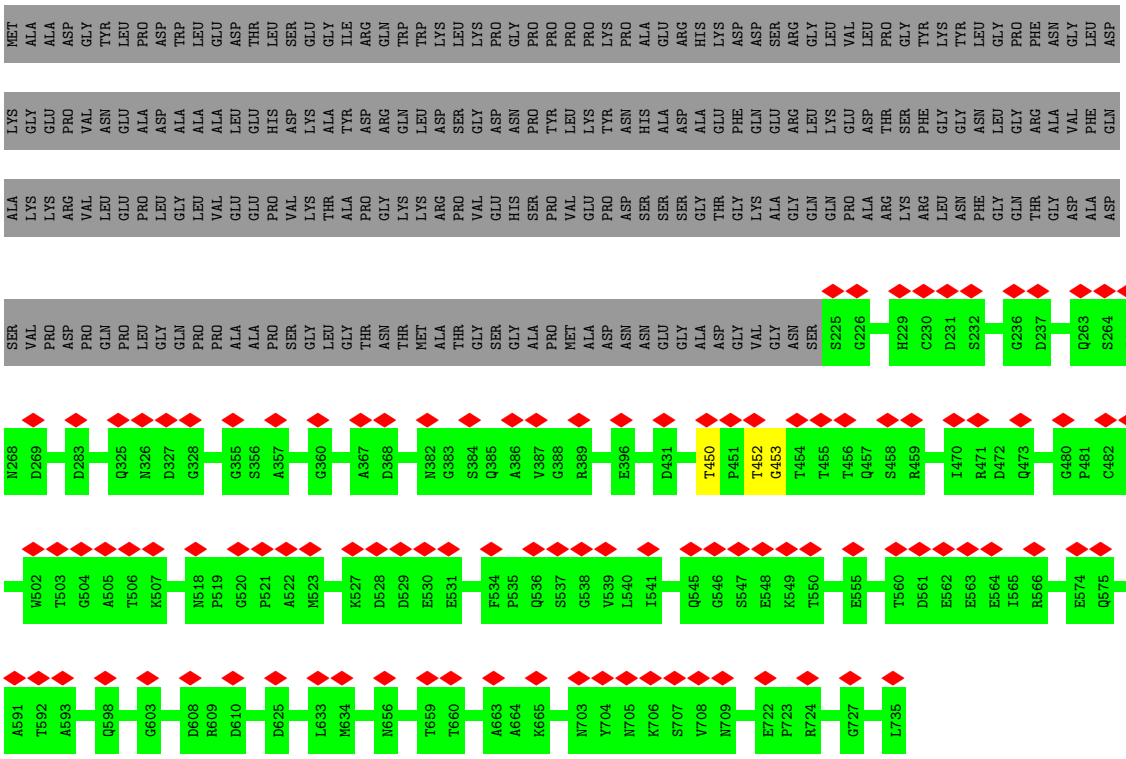
• Molecule 1: Capsid protein VP1

Chain u:

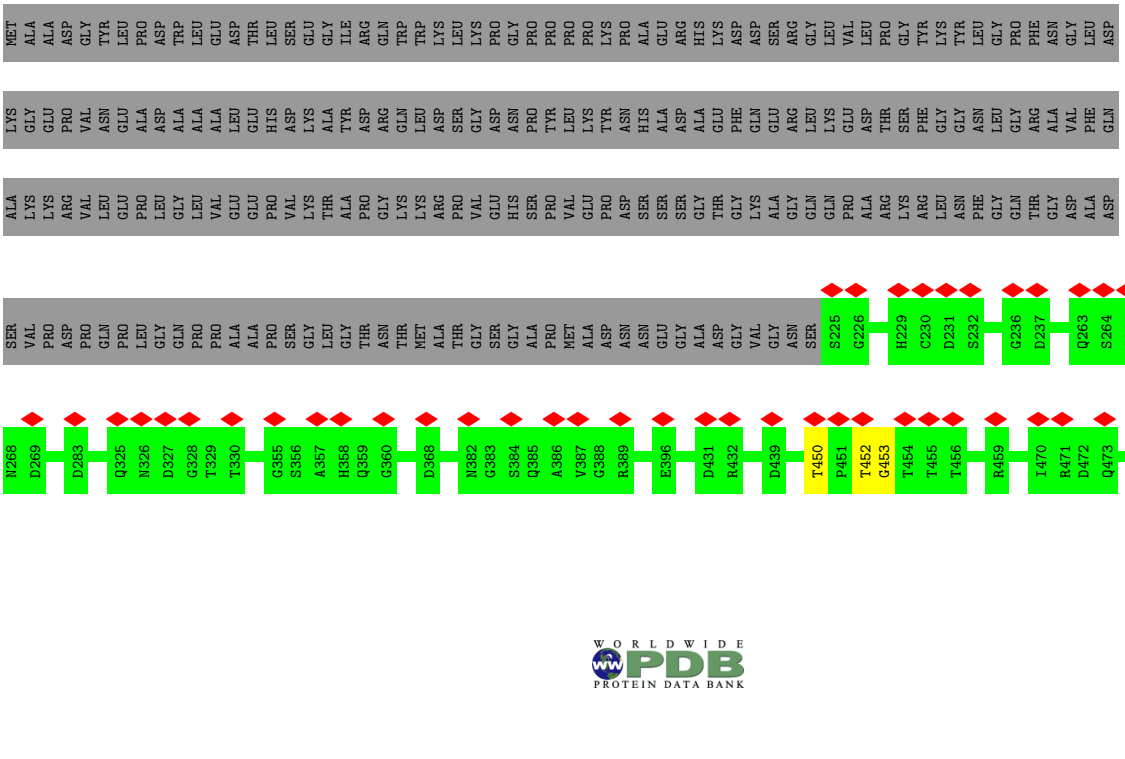


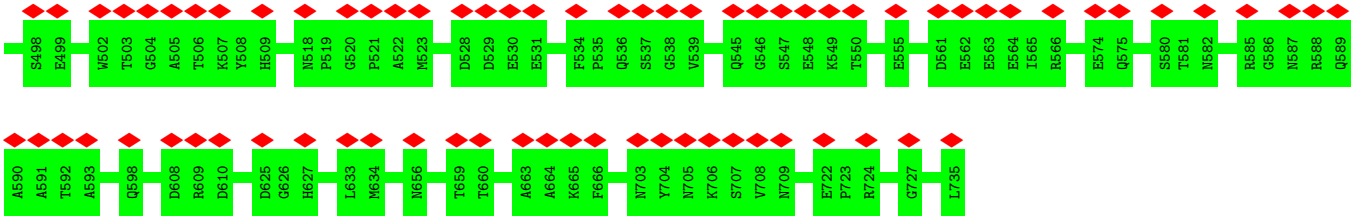


• Molecule 1: Capsid protein VP1

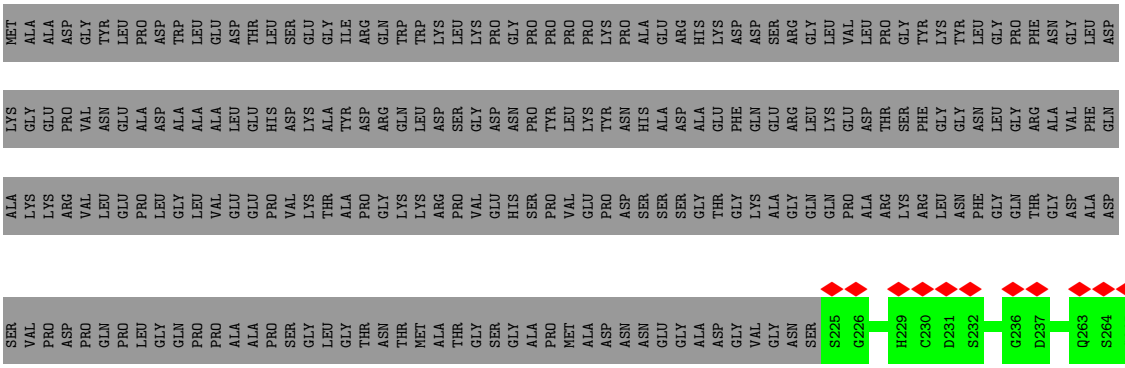


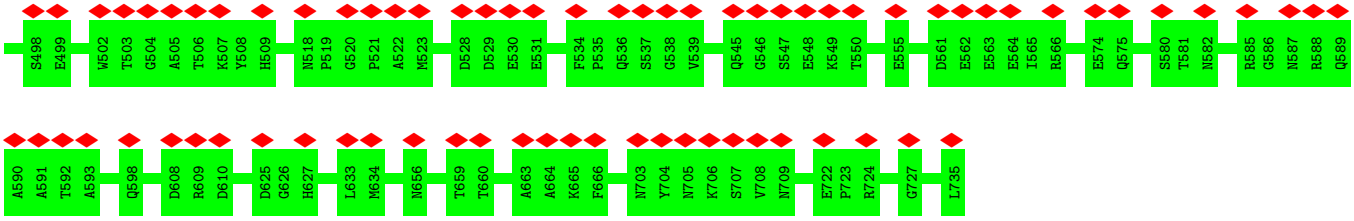
• Molecule 1: Capsid protein VP1



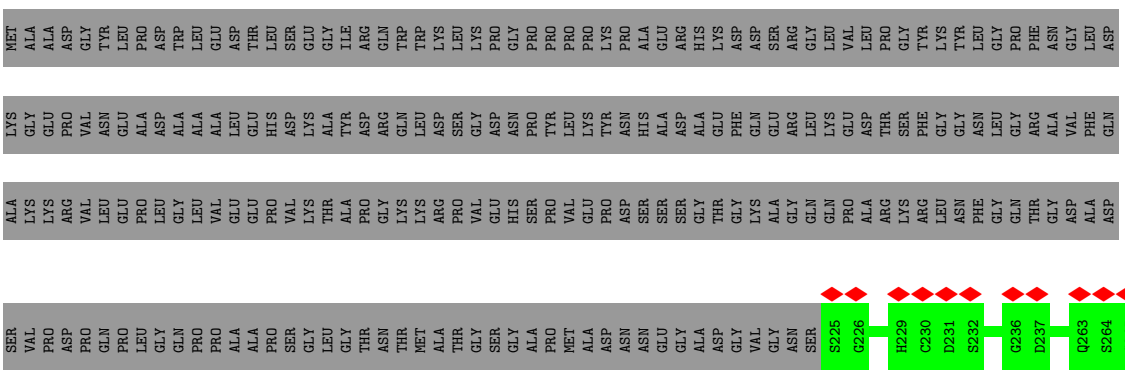


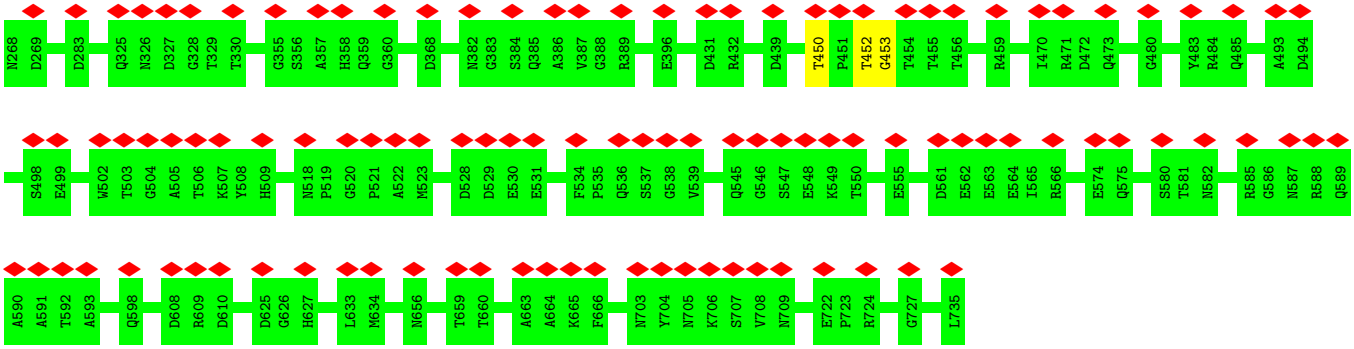
• Molecule 1: Capsid protein VP1





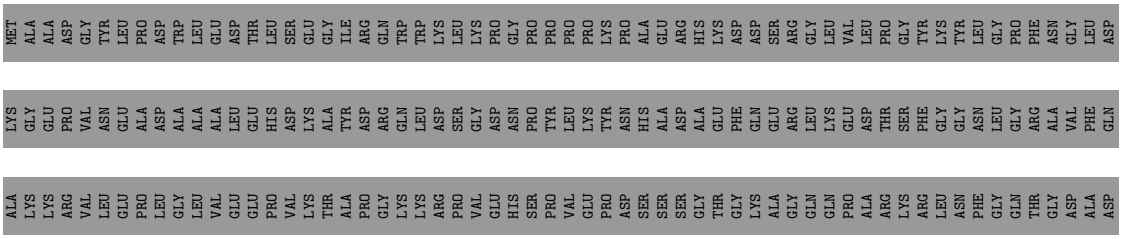
• Molecule 1: Capsid protein VP1

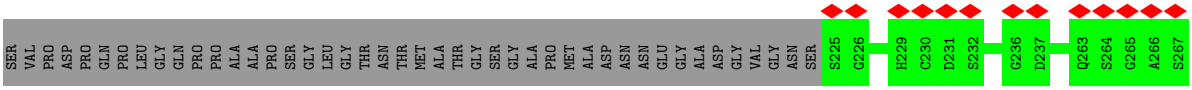


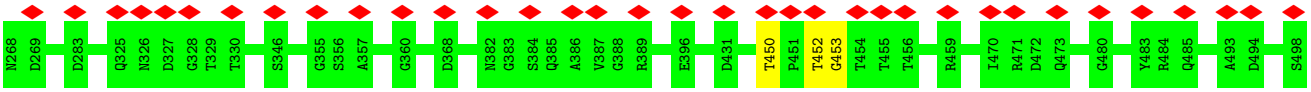


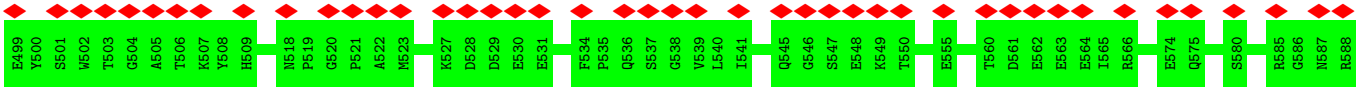
• Molecule 1: Capsid protein VP1

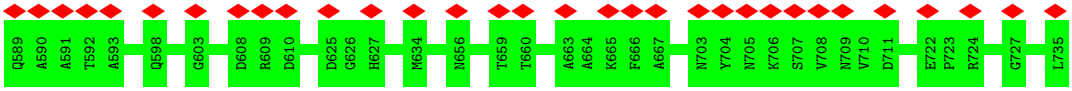
Chain z:





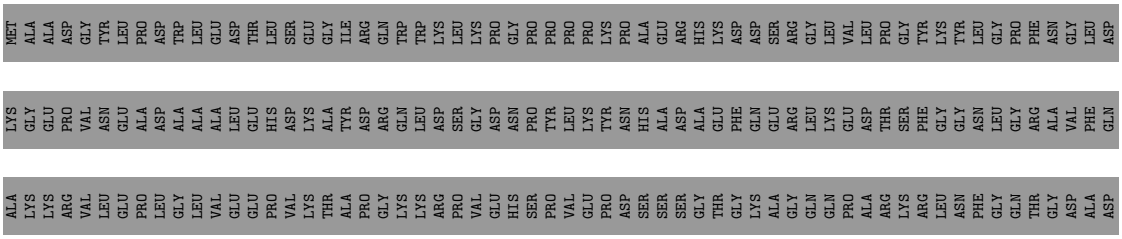


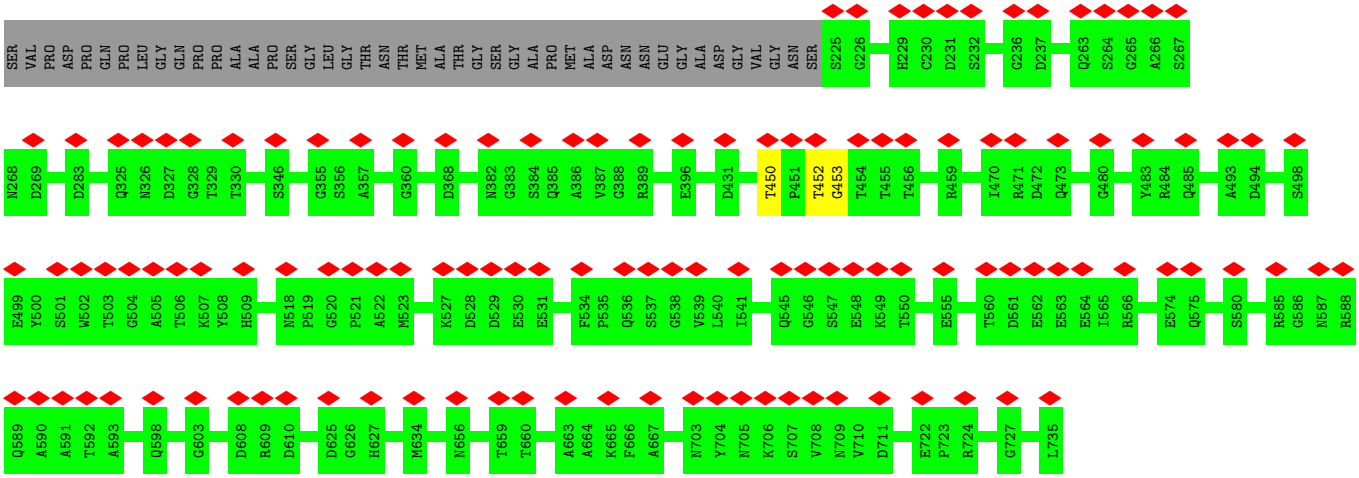




• Molecule 1: Capsid protein VP1

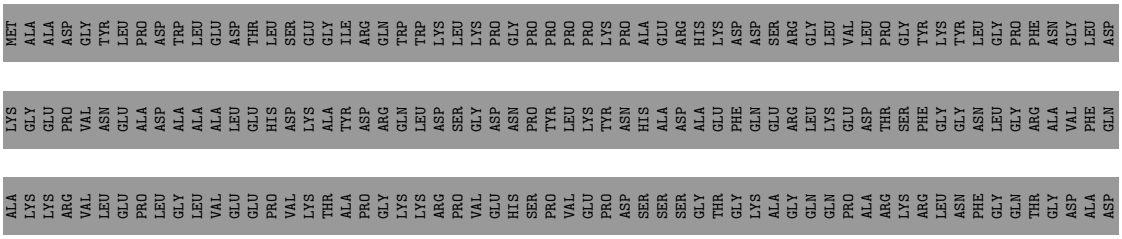
Chain 1:

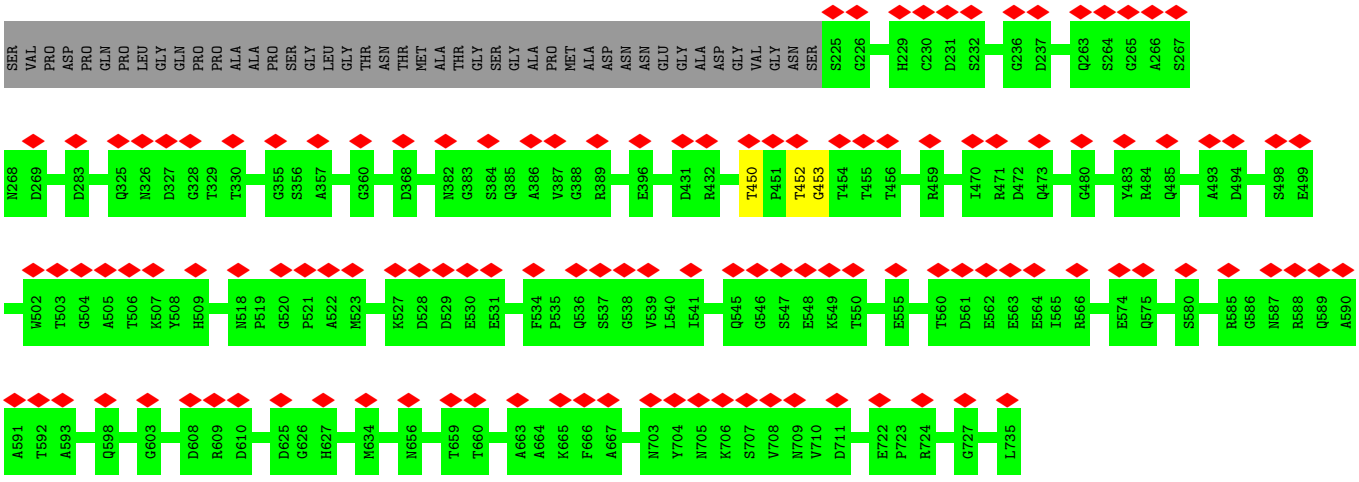




• Molecule 1: Capsid protein VP1

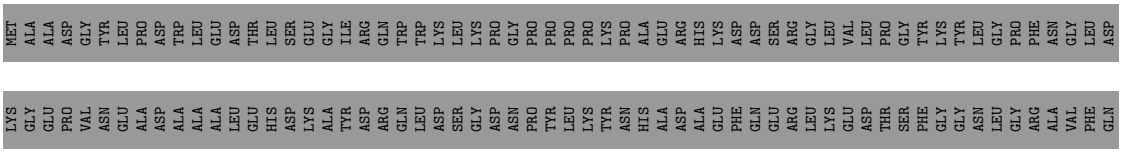
Chain 2:

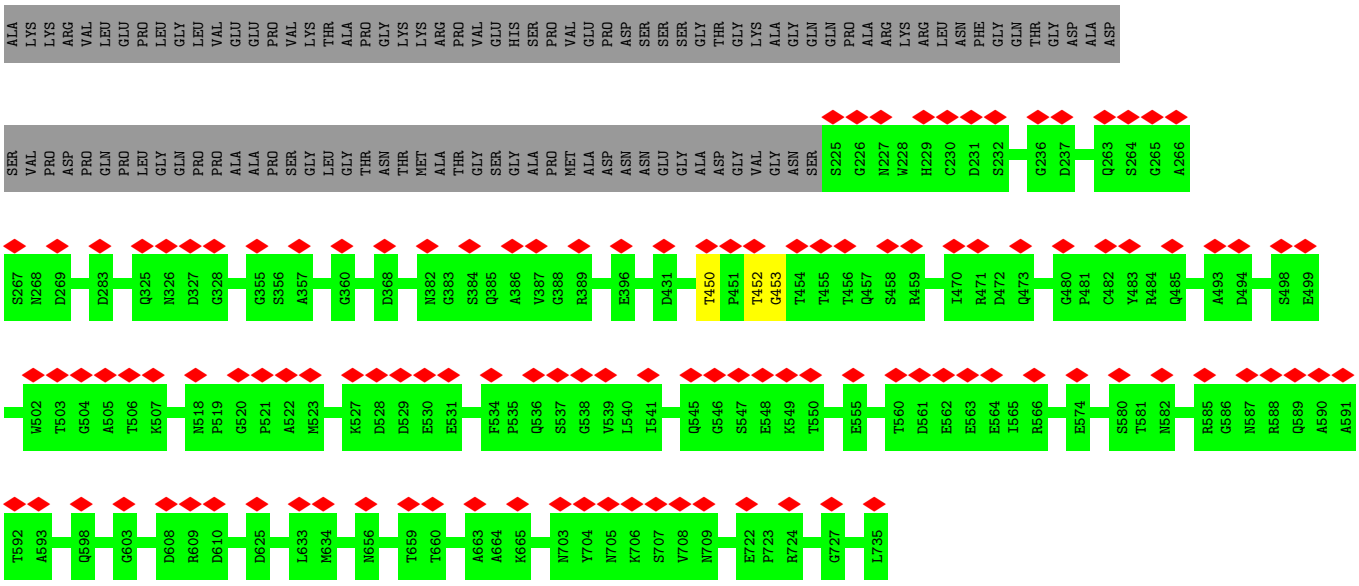




• Molecule 1: Capsid protein VP1

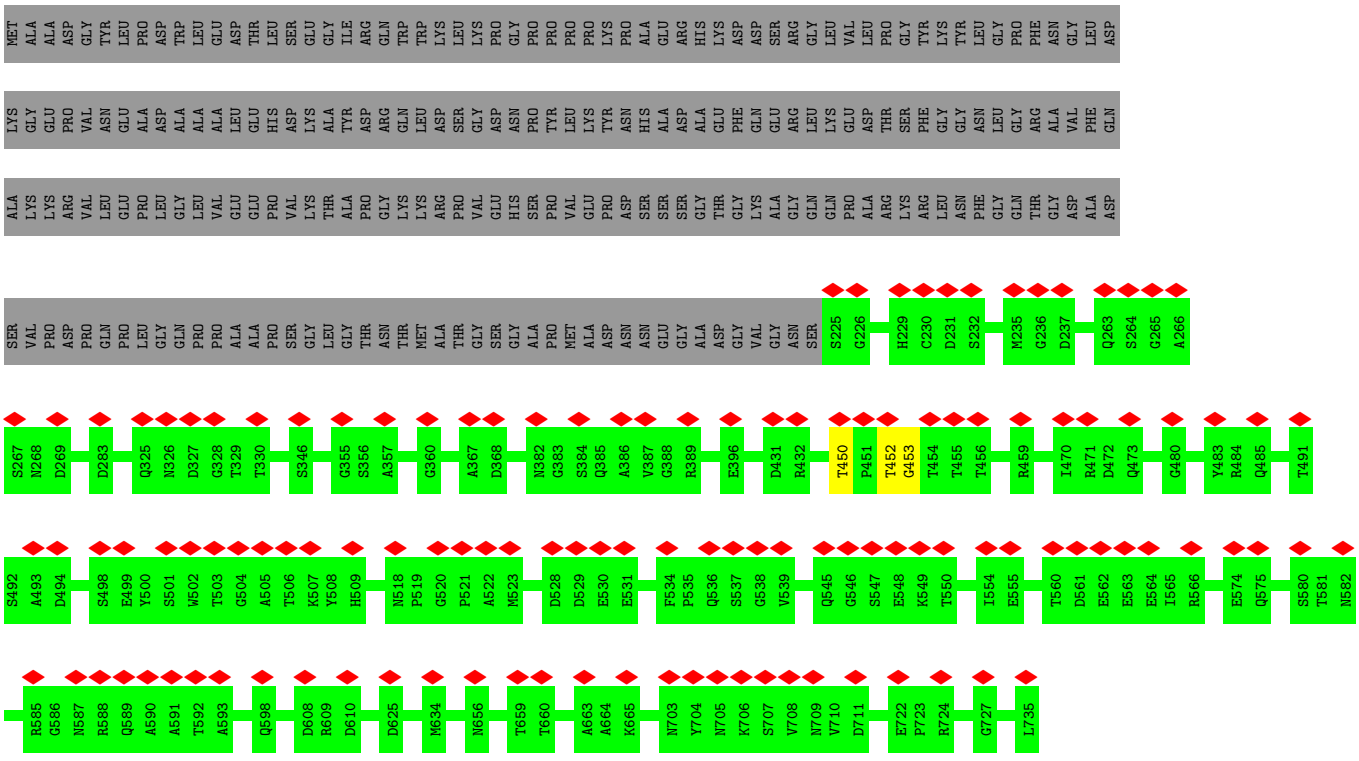
Chain 3:





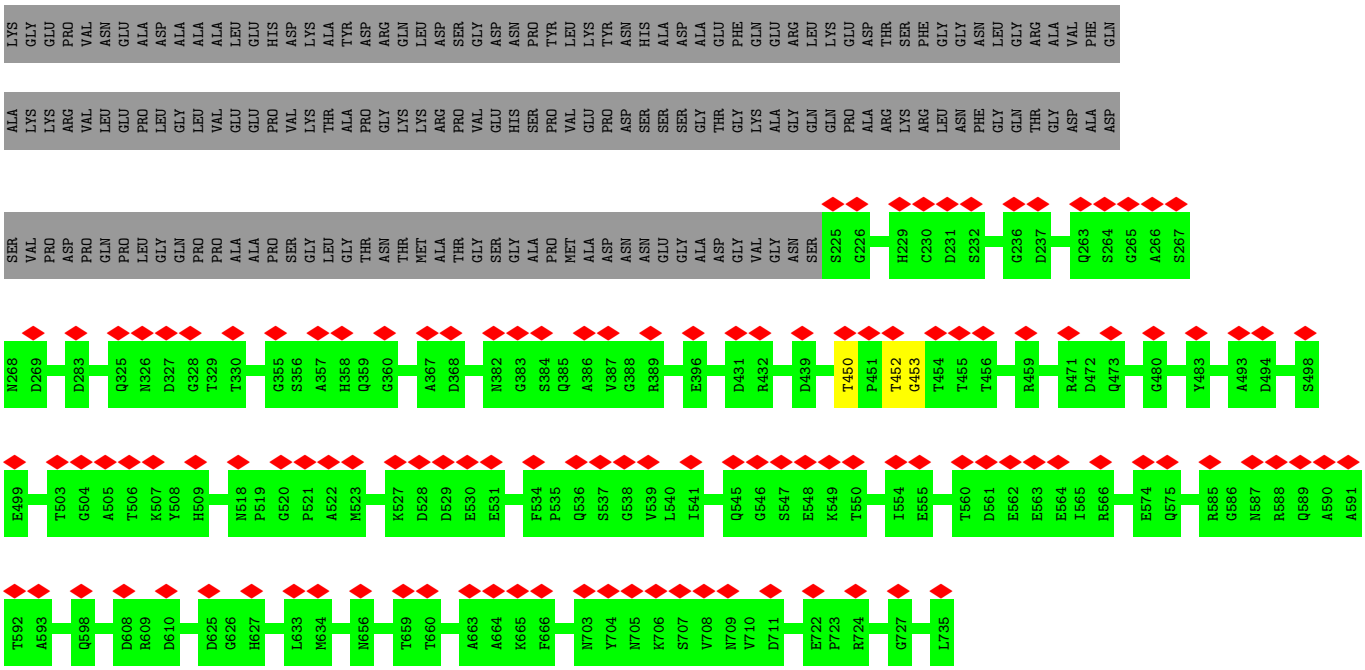
• Molecule 1: Capsid protein VP1

Chain 4:



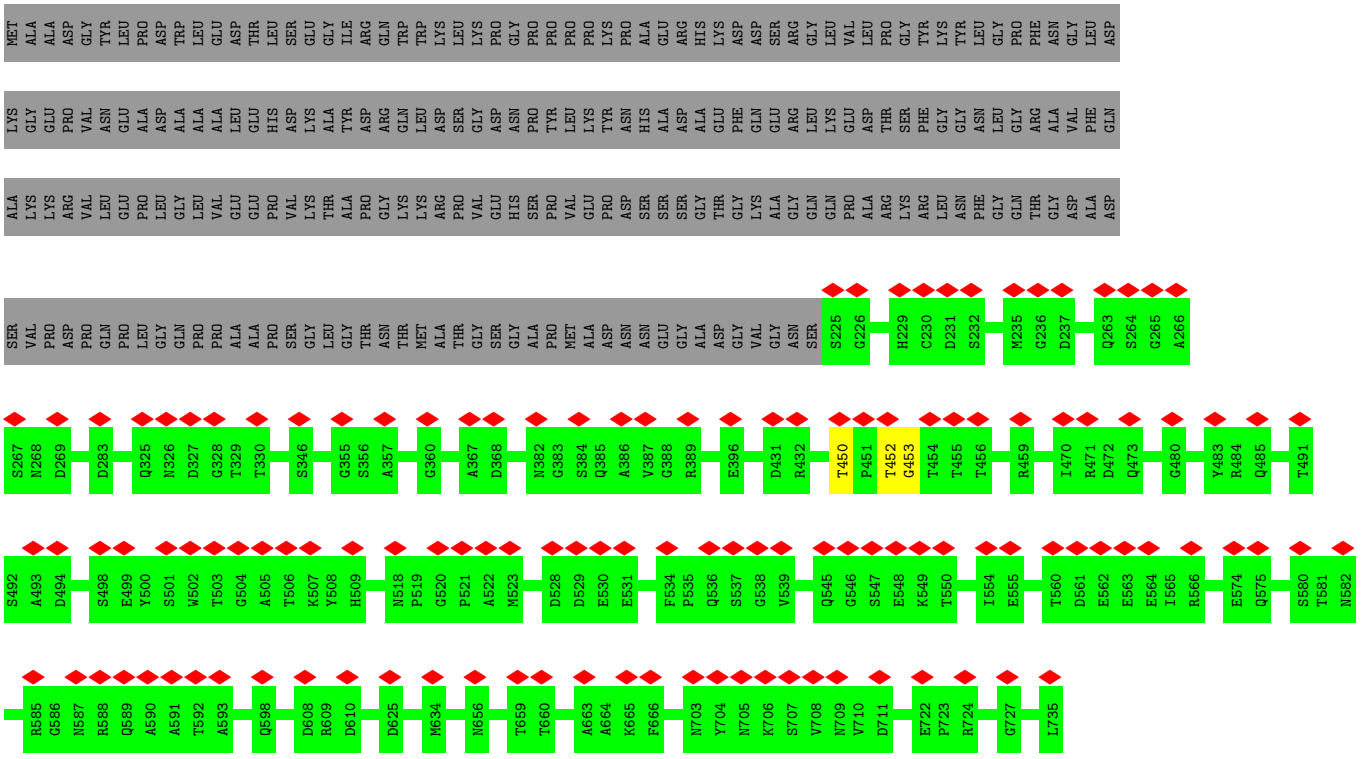
• Molecule 1: Capsid protein VP1

Chain 5:



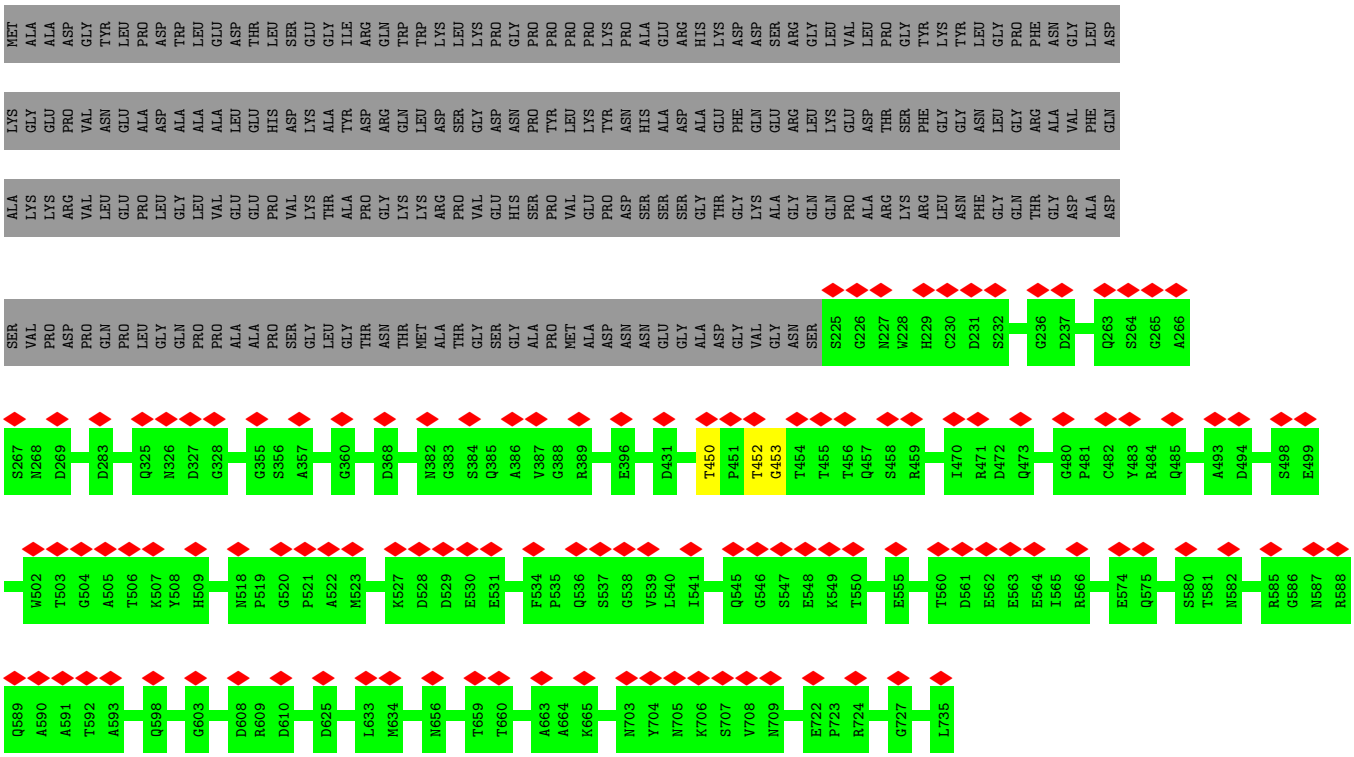
• Molecule 1: Capsid protein VP1

Chain 6:



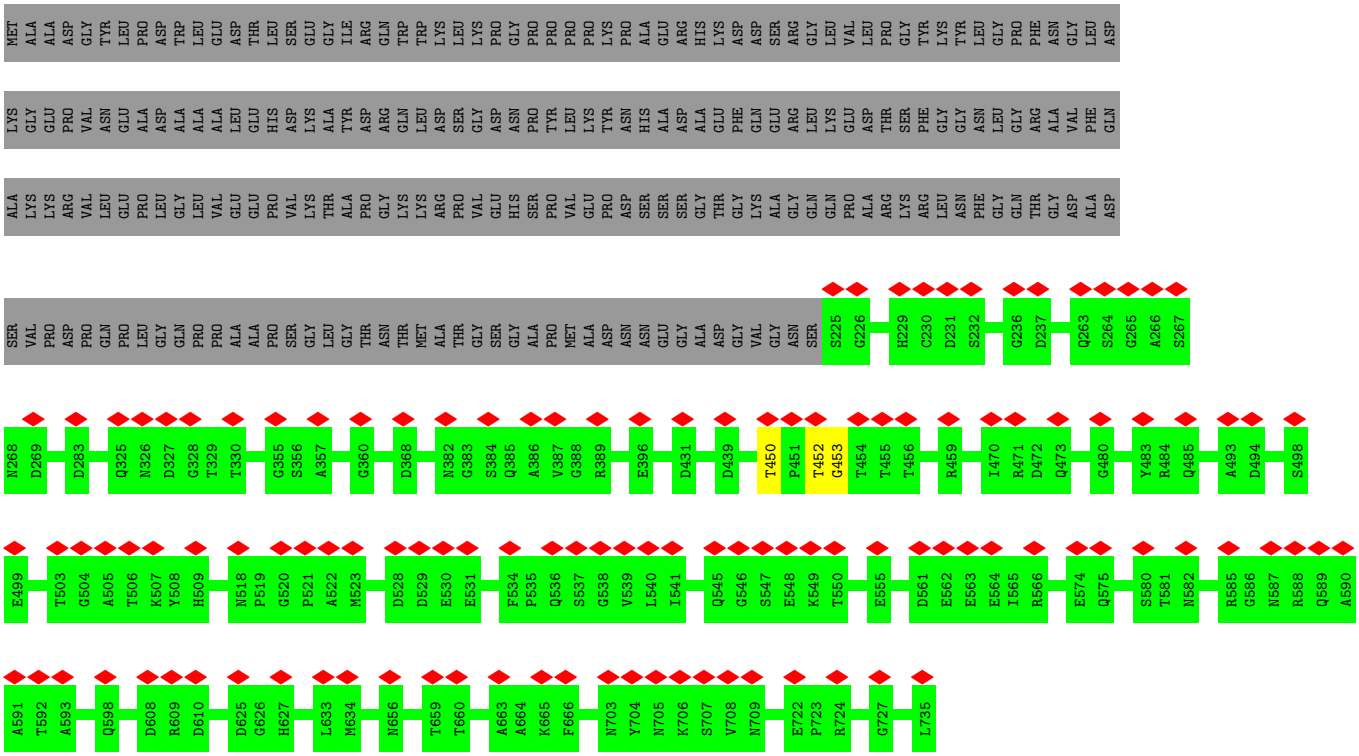
• Molecule 1: Capsid protein VP1

Chain 7:



• Molecule 1: Capsid protein VP1

Chain 8:



4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	23039	Depositor
Resolution determination method	FSC 0.5 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING ONLY	Depositor
Microscope	FEI POLARA 300	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/Å^2$)	20	Depositor
Minimum defocus (nm)	1100	Depositor
Maximum defocus (nm)	3700	Depositor
Magnification	56924	Depositor
Image detector	KODAK SO-163 FILM	Depositor
Maximum map value	14.791	Depositor
Minimum map value	-15.532	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	1.000	Depositor
Recommended contour level	2.9	Depositor
Map size (Å)	361.9, 361.9, 361.9	wwPDB
Map dimensions	329, 329, 329	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.1, 1.1, 1.1	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	1	0.54	0/4229	0.59	3/5767 (0.1%)
1	2	0.54	0/4229	0.59	3/5767 (0.1%)
1	3	0.54	0/4229	0.59	3/5767 (0.1%)
1	4	0.54	0/4229	0.59	3/5767 (0.1%)
1	5	0.54	0/4229	0.59	3/5767 (0.1%)
1	6	0.54	0/4229	0.59	3/5767 (0.1%)
1	7	0.54	0/4229	0.59	3/5767 (0.1%)
1	8	0.54	0/4229	0.59	3/5767 (0.1%)
1	A	0.54	0/4229	0.59	3/5767 (0.1%)
1	B	0.54	0/4229	0.59	3/5767 (0.1%)
1	C	0.54	0/4229	0.59	3/5767 (0.1%)
1	D	0.54	0/4229	0.59	3/5767 (0.1%)
1	E	0.54	0/4229	0.59	3/5767 (0.1%)
1	F	0.54	0/4229	0.59	3/5767 (0.1%)
1	G	0.54	0/4229	0.59	3/5767 (0.1%)
1	H	0.54	0/4229	0.59	3/5767 (0.1%)
1	I	0.54	0/4229	0.59	3/5767 (0.1%)
1	J	0.54	0/4229	0.59	3/5767 (0.1%)
1	K	0.54	0/4229	0.59	3/5767 (0.1%)
1	L	0.54	0/4229	0.59	3/5767 (0.1%)
1	M	0.54	0/4229	0.59	3/5767 (0.1%)
1	N	0.54	0/4229	0.59	3/5767 (0.1%)
1	O	0.54	0/4229	0.59	3/5767 (0.1%)
1	P	0.54	0/4229	0.59	3/5767 (0.1%)
1	Q	0.54	0/4229	0.59	3/5767 (0.1%)
1	R	0.54	0/4229	0.59	3/5767 (0.1%)
1	S	0.54	0/4229	0.59	3/5767 (0.1%)
1	T	0.54	0/4229	0.59	3/5767 (0.1%)
1	U	0.54	0/4229	0.59	3/5767 (0.1%)
1	V	0.54	0/4229	0.59	3/5767 (0.1%)
1	W	0.54	0/4229	0.59	3/5767 (0.1%)
1	X	0.54	0/4229	0.59	3/5767 (0.1%)
1	Y	0.54	0/4229	0.59	3/5767 (0.1%)
1	Z	0.54	0/4229	0.59	3/5767 (0.1%)

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	#|Z| >5	RMSZ	#|Z| >5
1	a	0.54	0/4229	0.59	3/5767 (0.1%)
1	b	0.54	0/4229	0.59	3/5767 (0.1%)
1	c	0.54	0/4229	0.59	3/5767 (0.1%)
1	d	0.54	0/4229	0.59	3/5767 (0.1%)
1	e	0.54	0/4229	0.59	3/5767 (0.1%)
1	f	0.54	0/4229	0.59	3/5767 (0.1%)
1	g	0.54	0/4229	0.59	3/5767 (0.1%)
1	h	0.54	0/4229	0.59	3/5767 (0.1%)
1	i	0.54	0/4229	0.59	3/5767 (0.1%)
1	j	0.54	0/4229	0.59	3/5767 (0.1%)
1	k	0.54	0/4229	0.59	3/5767 (0.1%)
1	l	0.54	0/4229	0.59	3/5767 (0.1%)
1	m	0.54	0/4229	0.59	3/5767 (0.1%)
1	n	0.54	0/4229	0.59	3/5767 (0.1%)
1	o	0.54	0/4229	0.59	3/5767 (0.1%)
1	p	0.54	0/4229	0.59	3/5767 (0.1%)
1	q	0.54	0/4229	0.59	3/5767 (0.1%)
1	r	0.54	0/4229	0.59	3/5767 (0.1%)
1	s	0.54	0/4229	0.59	3/5767 (0.1%)
1	t	0.54	0/4229	0.59	3/5767 (0.1%)
1	u	0.54	0/4229	0.59	3/5767 (0.1%)
1	v	0.54	0/4229	0.59	3/5767 (0.1%)
1	w	0.54	0/4229	0.59	3/5767 (0.1%)
1	x	0.54	0/4229	0.59	3/5767 (0.1%)
1	y	0.54	0/4229	0.59	3/5767 (0.1%)
1	z	0.54	0/4229	0.59	3/5767 (0.1%)
All	All	0.54	0/253740	0.59	180/346020 (0.1%)

There are no bond length outliers.

All (180) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	U	450	THR	C-N-CD	-8.07	102.84	120.60
1	w	450	THR	C-N-CD	-8.07	102.84	120.60
1	I	450	THR	C-N-CD	-8.07	102.85	120.60
1	l	450	THR	C-N-CD	-8.07	102.85	120.60
1	L	450	THR	C-N-CD	-8.07	102.85	120.60
1	V	450	THR	C-N-CD	-8.07	102.85	120.60
1	Z	450	THR	C-N-CD	-8.07	102.85	120.60
1	h	450	THR	C-N-CD	-8.07	102.85	120.60
1	q	450	THR	C-N-CD	-8.07	102.85	120.60
1	6	450	THR	C-N-CD	-8.07	102.85	120.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	450	THR	C-N-CD	-8.06	102.86	120.60
1	K	450	THR	C-N-CD	-8.06	102.86	120.60
1	D	450	THR	C-N-CD	-8.06	102.86	120.60
1	2	450	THR	C-N-CD	-8.06	102.86	120.60
1	c	450	THR	C-N-CD	-8.06	102.86	120.60
1	n	450	THR	C-N-CD	-8.06	102.86	120.60
1	r	450	THR	C-N-CD	-8.06	102.86	120.60
1	z	450	THR	C-N-CD	-8.06	102.86	120.60
1	1	450	THR	C-N-CD	-8.06	102.86	120.60
1	A	450	THR	C-N-CD	-8.06	102.87	120.60
1	j	450	THR	C-N-CD	-8.06	102.87	120.60
1	o	450	THR	C-N-CD	-8.06	102.87	120.60
1	t	450	THR	C-N-CD	-8.06	102.87	120.60
1	3	450	THR	C-N-CD	-8.06	102.87	120.60
1	7	450	THR	C-N-CD	-8.06	102.87	120.60
1	N	450	THR	C-N-CD	-8.05	102.88	120.60
1	T	450	THR	C-N-CD	-8.06	102.88	120.60
1	C	450	THR	C-N-CD	-8.05	102.89	120.60
1	a	450	THR	C-N-CD	-8.05	102.89	120.60
1	k	450	THR	C-N-CD	-8.05	102.89	120.60
1	g	450	THR	C-N-CD	-8.05	102.89	120.60
1	i	450	THR	C-N-CD	-8.05	102.89	120.60
1	s	450	THR	C-N-CD	-8.05	102.89	120.60
1	J	450	THR	C-N-CD	-8.05	102.89	120.60
1	m	450	THR	C-N-CD	-8.05	102.89	120.60
1	4	450	THR	C-N-CD	-8.05	102.90	120.60
1	E	450	THR	C-N-CD	-8.04	102.90	120.60
1	H	450	THR	C-N-CD	-8.04	102.90	120.60
1	M	450	THR	C-N-CD	-8.04	102.90	120.60
1	O	450	THR	C-N-CD	-8.04	102.90	120.60
1	P	450	THR	C-N-CD	-8.04	102.90	120.60
1	S	450	THR	C-N-CD	-8.04	102.90	120.60
1	b	450	THR	C-N-CD	-8.05	102.90	120.60
1	e	450	THR	C-N-CD	-8.04	102.90	120.60
1	f	450	THR	C-N-CD	-8.04	102.90	120.60
1	G	450	THR	C-N-CD	-8.04	102.91	120.60
1	R	450	THR	C-N-CD	-8.04	102.91	120.60
1	Y	450	THR	C-N-CD	-8.04	102.91	120.60
1	5	450	THR	C-N-CD	-8.04	102.91	120.60
1	8	450	THR	C-N-CD	-8.04	102.91	120.60
1	F	450	THR	C-N-CD	-8.04	102.92	120.60
1	X	450	THR	C-N-CD	-8.04	102.91	120.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	u	450	THR	C-N-CD	-8.04	102.92	120.60
1	x	450	THR	C-N-CD	-8.04	102.92	120.60
1	y	450	THR	C-N-CD	-8.04	102.92	120.60
1	W	450	THR	C-N-CD	-8.04	102.92	120.60
1	Q	450	THR	C-N-CD	-8.04	102.92	120.60
1	d	450	THR	C-N-CD	-8.04	102.92	120.60
1	p	450	THR	C-N-CD	-8.03	102.94	120.60
1	v	450	THR	C-N-CD	-8.03	102.94	120.60
1	c	450	THR	C-N-CA	5.96	147.03	122.00
1	V	450	THR	C-N-CA	5.96	147.03	122.00
1	C	450	THR	C-N-CA	5.95	147.00	122.00
1	T	450	THR	C-N-CA	5.95	147.00	122.00
1	h	450	THR	C-N-CA	5.95	147.00	122.00
1	2	450	THR	C-N-CA	5.95	147.01	122.00
1	z	450	THR	C-N-CA	5.95	147.00	122.00
1	J	450	THR	C-N-CA	5.95	147.00	122.00
1	K	450	THR	C-N-CA	5.95	146.99	122.00
1	M	450	THR	C-N-CA	5.95	147.00	122.00
1	q	450	THR	C-N-CA	5.95	146.99	122.00
1	L	450	THR	C-N-CA	5.95	146.99	122.00
1	Z	450	THR	C-N-CA	5.95	146.99	122.00
1	i	450	THR	C-N-CA	5.95	146.98	122.00
1	k	450	THR	C-N-CA	5.95	146.99	122.00
1	w	450	THR	C-N-CA	5.95	146.98	122.00
1	N	450	THR	C-N-CA	5.95	146.98	122.00
1	p	450	THR	C-N-CA	5.95	146.98	122.00
1	B	450	THR	C-N-CA	5.95	146.97	122.00
1	X	450	THR	C-N-CA	5.95	146.97	122.00
1	b	450	THR	C-N-CA	5.95	146.97	122.00
1	s	450	THR	C-N-CA	5.95	146.97	122.00
1	3	450	THR	C-N-CA	5.95	146.97	122.00
1	4	450	THR	C-N-CA	5.95	146.97	122.00
1	E	450	THR	C-N-CA	5.94	146.96	122.00
1	O	450	THR	C-N-CA	5.94	146.96	122.00
1	f	450	THR	C-N-CA	5.94	146.96	122.00
1	j	450	THR	C-N-CA	5.94	146.96	122.00
1	t	450	THR	C-N-CA	5.94	146.96	122.00
1	7	450	THR	C-N-CA	5.94	146.96	122.00
1	D	450	THR	C-N-CA	5.94	146.96	122.00
1	I	450	THR	C-N-CA	5.94	146.96	122.00
1	Q	450	THR	C-N-CA	5.94	146.96	122.00
1	R	450	THR	C-N-CA	5.94	146.96	122.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	S	450	THR	C-N-CA	5.94	146.96	122.00
1	U	450	THR	C-N-CA	5.94	146.96	122.00
1	Y	450	THR	C-N-CA	5.94	146.96	122.00
1	a	450	THR	C-N-CA	5.94	146.96	122.00
1	l	450	THR	C-N-CA	5.94	146.96	122.00
1	m	450	THR	C-N-CA	5.94	146.96	122.00
1	n	450	THR	C-N-CA	5.94	146.96	122.00
1	u	450	THR	C-N-CA	5.94	146.96	122.00
1	1	450	THR	C-N-CA	5.94	146.96	122.00
1	8	450	THR	C-N-CA	5.94	146.96	122.00
1	P	450	THR	C-N-CA	5.94	146.95	122.00
1	e	450	THR	C-N-CA	5.94	146.95	122.00
1	r	450	THR	C-N-CA	5.94	146.94	122.00
1	A	450	THR	C-N-CA	5.94	146.94	122.00
1	G	450	THR	C-N-CA	5.94	146.94	122.00
1	g	450	THR	C-N-CA	5.94	146.94	122.00
1	o	450	THR	C-N-CA	5.94	146.94	122.00
1	6	450	THR	C-N-CA	5.94	146.94	122.00
1	d	450	THR	C-N-CA	5.94	146.93	122.00
1	F	450	THR	C-N-CA	5.93	146.92	122.00
1	H	450	THR	C-N-CA	5.93	146.93	122.00
1	v	450	THR	C-N-CA	5.93	146.92	122.00
1	W	450	THR	C-N-CA	5.93	146.91	122.00
1	5	450	THR	C-N-CA	5.93	146.92	122.00
1	x	450	THR	C-N-CA	5.93	146.91	122.00
1	y	450	THR	C-N-CA	5.93	146.89	122.00
1	f	453	GLY	N-CA-C	5.21	126.12	113.10
1	8	453	GLY	N-CA-C	5.20	126.11	113.10
1	v	453	GLY	N-CA-C	5.20	126.10	113.10
1	r	453	GLY	N-CA-C	5.20	126.10	113.10
1	5	453	GLY	N-CA-C	5.20	126.10	113.10
1	K	453	GLY	N-CA-C	5.20	126.09	113.10
1	M	453	GLY	N-CA-C	5.20	126.10	113.10
1	R	453	GLY	N-CA-C	5.20	126.10	113.10
1	i	453	GLY	N-CA-C	5.20	126.09	113.10
1	s	453	GLY	N-CA-C	5.20	126.09	113.10
1	W	453	GLY	N-CA-C	5.20	126.09	113.10
1	w	453	GLY	N-CA-C	5.20	126.09	113.10
1	7	453	GLY	N-CA-C	5.20	126.09	113.10
1	g	453	GLY	N-CA-C	5.19	126.08	113.10
1	F	453	GLY	N-CA-C	5.19	126.08	113.10
1	Q	453	GLY	N-CA-C	5.19	126.08	113.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	Z	453	GLY	N-CA-C	5.19	126.08	113.10
1	p	453	GLY	N-CA-C	5.19	126.08	113.10
1	1	453	GLY	N-CA-C	5.19	126.08	113.10
1	X	453	GLY	N-CA-C	5.19	126.07	113.10
1	m	453	GLY	N-CA-C	5.19	126.07	113.10
1	x	453	GLY	N-CA-C	5.19	126.07	113.10
1	O	453	GLY	N-CA-C	5.19	126.07	113.10
1	U	453	GLY	N-CA-C	5.19	126.07	113.10
1	q	453	GLY	N-CA-C	5.19	126.07	113.10
1	B	453	GLY	N-CA-C	5.18	126.06	113.10
1	E	453	GLY	N-CA-C	5.18	126.06	113.10
1	D	453	GLY	N-CA-C	5.18	126.06	113.10
1	G	453	GLY	N-CA-C	5.18	126.06	113.10
1	I	453	GLY	N-CA-C	5.18	126.06	113.10
1	J	453	GLY	N-CA-C	5.18	126.06	113.10
1	N	453	GLY	N-CA-C	5.18	126.06	113.10
1	P	453	GLY	N-CA-C	5.18	126.06	113.10
1	d	453	GLY	N-CA-C	5.18	126.06	113.10
1	e	453	GLY	N-CA-C	5.18	126.06	113.10
1	j	453	GLY	N-CA-C	5.18	126.06	113.10
1	k	453	GLY	N-CA-C	5.18	126.05	113.10
1	y	453	GLY	N-CA-C	5.18	126.06	113.10
1	b	453	GLY	N-CA-C	5.18	126.05	113.10
1	n	453	GLY	N-CA-C	5.18	126.05	113.10
1	u	453	GLY	N-CA-C	5.18	126.05	113.10
1	H	453	GLY	N-CA-C	5.18	126.05	113.10
1	T	453	GLY	N-CA-C	5.18	126.05	113.10
1	Y	453	GLY	N-CA-C	5.18	126.05	113.10
1	a	453	GLY	N-CA-C	5.18	126.05	113.10
1	c	453	GLY	N-CA-C	5.18	126.05	113.10
1	t	453	GLY	N-CA-C	5.18	126.05	113.10
1	2	453	GLY	N-CA-C	5.18	126.05	113.10
1	z	453	GLY	N-CA-C	5.18	126.05	113.10
1	L	453	GLY	N-CA-C	5.17	126.03	113.10
1	V	453	GLY	N-CA-C	5.17	126.03	113.10
1	3	453	GLY	N-CA-C	5.17	126.04	113.10
1	6	453	GLY	N-CA-C	5.17	126.03	113.10
1	h	453	GLY	N-CA-C	5.17	126.03	113.10
1	l	453	GLY	N-CA-C	5.17	126.03	113.10
1	A	453	GLY	N-CA-C	5.17	126.02	113.10
1	C	453	GLY	N-CA-C	5.17	126.02	113.10
1	o	453	GLY	N-CA-C	5.17	126.02	113.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	4	453	GLY	N-CA-C	5.16	126.00	113.10
1	S	453	GLY	N-CA-C	5.16	125.99	113.10

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

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

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	1	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	2	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	3	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	4	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	5	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	6	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	7	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	8	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	A	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	B	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	C	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	D	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	E	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	F	509/735 (69%)	469 (92%)	40 (8%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	G	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	H	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	I	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	J	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	K	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	L	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	M	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	N	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	O	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	P	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	Q	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	R	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	S	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	T	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	U	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	V	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	W	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	X	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	Y	509/735 (69%)	468 (92%)	41 (8%)	0	100	100
1	Z	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	a	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	b	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	c	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	d	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	e	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	f	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	g	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	h	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	i	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	j	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	k	509/735 (69%)	469 (92%)	40 (8%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	l	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	m	509/735 (69%)	468 (92%)	41 (8%)	0	100	100
1	n	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	o	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	p	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	q	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	r	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	s	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	t	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	u	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	v	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	w	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	x	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	y	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
1	z	509/735 (69%)	469 (92%)	40 (8%)	0	100	100
All	All	30540/44100 (69%)	28138 (92%)	2402 (8%)	0	100	100

There are no Ramachandran outliers to report.

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	1	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	2	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	3	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	4	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	5	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	6	455/630 (72%)	454 (100%)	1 (0%)	93	97

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	7	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	8	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	A	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	B	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	C	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	D	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	E	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	F	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	G	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	H	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	I	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	J	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	K	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	L	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	M	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	N	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	O	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	P	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	Q	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	R	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	S	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	T	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	U	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	V	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	W	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	X	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	Y	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	Z	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	a	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	b	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	c	455/630 (72%)	454 (100%)	1 (0%)	93	97

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	d	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	e	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	f	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	g	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	h	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	i	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	j	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	k	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	l	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	m	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	n	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	o	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	p	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	q	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	r	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	s	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	t	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	u	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	v	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	w	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	x	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	y	455/630 (72%)	454 (100%)	1 (0%)	93	97
1	z	455/630 (72%)	454 (100%)	1 (0%)	93	97
All	All	27300/37800 (72%)	27240 (100%)	60 (0%)	93	97

All (60) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	452	THR
1	B	452	THR
1	C	452	THR
1	D	452	THR
1	E	452	THR
1	F	452	THR

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Mol	Chain	Res	Type
1	G	452	THR
1	H	452	THR
1	I	452	THR
1	J	452	THR
1	K	452	THR
1	L	452	THR
1	M	452	THR
1	N	452	THR
1	O	452	THR
1	P	452	THR
1	Q	452	THR
1	R	452	THR
1	S	452	THR
1	T	452	THR
1	U	452	THR
1	V	452	THR
1	W	452	THR
1	X	452	THR
1	Y	452	THR
1	Z	452	THR
1	a	452	THR
1	b	452	THR
1	c	452	THR
1	d	452	THR
1	e	452	THR
1	f	452	THR
1	g	452	THR
1	h	452	THR
1	i	452	THR
1	j	452	THR
1	k	452	THR
1	l	452	THR
1	m	452	THR
1	n	452	THR
1	o	452	THR
1	p	452	THR
1	q	452	THR
1	r	452	THR
1	s	452	THR
1	t	452	THR
1	u	452	THR
1	v	452	THR

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Mol	Chain	Res	Type
1	w	452	THR
1	x	452	THR
1	y	452	THR
1	z	452	THR
1	1	452	THR
1	2	452	THR
1	3	452	THR
1	4	452	THR
1	5	452	THR
1	6	452	THR
1	7	452	THR
1	8	452	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (424) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	325	GLN
1	A	457	GLN
1	A	485	GLN
1	A	497	ASN
1	A	518	ASN
1	A	598	GLN
1	A	650	ASN
1	B	325	GLN
1	B	457	GLN
1	B	485	GLN
1	B	497	ASN
1	B	518	ASN
1	B	598	GLN
1	B	650	ASN
1	C	325	GLN
1	C	457	GLN
1	C	485	GLN
1	C	497	ASN
1	C	518	ASN
1	C	598	GLN
1	C	650	ASN
1	D	325	GLN
1	D	457	GLN
1	D	485	GLN
1	D	497	ASN
1	D	518	ASN

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Mol	Chain	Res	Type
1	D	598	GLN
1	D	650	ASN
1	E	325	GLN
1	E	457	GLN
1	E	485	GLN
1	E	497	ASN
1	E	518	ASN
1	E	598	GLN
1	E	650	ASN
1	F	325	GLN
1	F	457	GLN
1	F	485	GLN
1	F	497	ASN
1	F	518	ASN
1	F	598	GLN
1	F	650	ASN
1	G	325	GLN
1	G	457	GLN
1	G	485	GLN
1	G	497	ASN
1	G	518	ASN
1	G	598	GLN
1	G	650	ASN
1	H	325	GLN
1	H	457	GLN
1	H	485	GLN
1	H	497	ASN
1	H	518	ASN
1	H	598	GLN
1	H	650	ASN
1	I	325	GLN
1	I	457	GLN
1	I	485	GLN
1	I	497	ASN
1	I	518	ASN
1	I	598	GLN
1	I	650	ASN
1	J	325	GLN
1	J	457	GLN
1	J	485	GLN
1	J	497	ASN
1	J	518	ASN

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Mol	Chain	Res	Type
1	J	598	GLN
1	J	650	ASN
1	K	325	GLN
1	K	457	GLN
1	K	485	GLN
1	K	497	ASN
1	K	518	ASN
1	K	598	GLN
1	K	650	ASN
1	L	325	GLN
1	L	457	GLN
1	L	485	GLN
1	L	497	ASN
1	L	518	ASN
1	L	598	GLN
1	L	650	ASN
1	M	325	GLN
1	M	457	GLN
1	M	485	GLN
1	M	497	ASN
1	M	518	ASN
1	M	598	GLN
1	M	650	ASN
1	N	325	GLN
1	N	457	GLN
1	N	485	GLN
1	N	497	ASN
1	N	518	ASN
1	N	598	GLN
1	N	650	ASN
1	O	325	GLN
1	O	457	GLN
1	O	485	GLN
1	O	497	ASN
1	O	518	ASN
1	O	598	GLN
1	O	650	ASN
1	P	325	GLN
1	P	457	GLN
1	P	485	GLN
1	P	497	ASN
1	P	518	ASN

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Mol	Chain	Res	Type
1	P	598	GLN
1	P	650	ASN
1	Q	325	GLN
1	Q	457	GLN
1	Q	485	GLN
1	Q	497	ASN
1	Q	518	ASN
1	Q	598	GLN
1	Q	650	ASN
1	R	325	GLN
1	R	457	GLN
1	R	485	GLN
1	R	497	ASN
1	R	518	ASN
1	R	598	GLN
1	R	650	ASN
1	S	325	GLN
1	S	457	GLN
1	S	485	GLN
1	S	497	ASN
1	S	518	ASN
1	S	598	GLN
1	S	650	ASN
1	T	325	GLN
1	T	457	GLN
1	T	485	GLN
1	T	497	ASN
1	T	518	ASN
1	T	598	GLN
1	T	650	ASN
1	U	325	GLN
1	U	457	GLN
1	U	485	GLN
1	U	497	ASN
1	U	518	ASN
1	U	598	GLN
1	U	650	ASN
1	V	325	GLN
1	V	457	GLN
1	V	485	GLN
1	V	497	ASN
1	V	518	ASN

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Mol	Chain	Res	Type
1	V	598	GLN
1	V	650	ASN
1	W	325	GLN
1	W	449	ASN
1	W	457	GLN
1	W	485	GLN
1	W	497	ASN
1	W	518	ASN
1	W	598	GLN
1	W	650	ASN
1	X	325	GLN
1	X	457	GLN
1	X	485	GLN
1	X	497	ASN
1	X	518	ASN
1	X	598	GLN
1	X	650	ASN
1	Y	325	GLN
1	Y	457	GLN
1	Y	485	GLN
1	Y	497	ASN
1	Y	518	ASN
1	Y	598	GLN
1	Y	650	ASN
1	Z	325	GLN
1	Z	457	GLN
1	Z	485	GLN
1	Z	497	ASN
1	Z	518	ASN
1	Z	598	GLN
1	Z	650	ASN
1	a	325	GLN
1	a	457	GLN
1	a	485	GLN
1	a	497	ASN
1	a	518	ASN
1	a	598	GLN
1	a	650	ASN
1	b	325	GLN
1	b	457	GLN
1	b	485	GLN
1	b	497	ASN

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Mol	Chain	Res	Type
1	b	518	ASN
1	b	598	GLN
1	b	650	ASN
1	c	325	GLN
1	c	457	GLN
1	c	485	GLN
1	c	497	ASN
1	c	518	ASN
1	c	598	GLN
1	c	650	ASN
1	d	325	GLN
1	d	457	GLN
1	d	485	GLN
1	d	497	ASN
1	d	518	ASN
1	d	598	GLN
1	d	650	ASN
1	e	325	GLN
1	e	449	ASN
1	e	457	GLN
1	e	485	GLN
1	e	497	ASN
1	e	518	ASN
1	e	598	GLN
1	e	650	ASN
1	f	325	GLN
1	f	457	GLN
1	f	485	GLN
1	f	497	ASN
1	f	518	ASN
1	f	598	GLN
1	f	650	ASN
1	g	325	GLN
1	g	457	GLN
1	g	485	GLN
1	g	497	ASN
1	g	518	ASN
1	g	598	GLN
1	g	650	ASN
1	h	325	GLN
1	h	457	GLN
1	h	485	GLN

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Mol	Chain	Res	Type
1	h	497	ASN
1	h	518	ASN
1	h	598	GLN
1	h	650	ASN
1	i	325	GLN
1	i	457	GLN
1	i	485	GLN
1	i	497	ASN
1	i	518	ASN
1	i	598	GLN
1	i	650	ASN
1	j	325	GLN
1	j	457	GLN
1	j	485	GLN
1	j	497	ASN
1	j	518	ASN
1	j	598	GLN
1	j	650	ASN
1	k	325	GLN
1	k	457	GLN
1	k	485	GLN
1	k	497	ASN
1	k	518	ASN
1	k	598	GLN
1	k	650	ASN
1	l	325	GLN
1	l	457	GLN
1	l	485	GLN
1	l	497	ASN
1	l	518	ASN
1	l	598	GLN
1	l	650	ASN
1	m	325	GLN
1	m	457	GLN
1	m	485	GLN
1	m	497	ASN
1	m	518	ASN
1	m	598	GLN
1	m	650	ASN
1	n	325	GLN
1	n	457	GLN
1	n	485	GLN

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Mol	Chain	Res	Type
1	n	497	ASN
1	n	518	ASN
1	n	598	GLN
1	n	650	ASN
1	o	325	GLN
1	o	457	GLN
1	o	485	GLN
1	o	497	ASN
1	o	518	ASN
1	o	598	GLN
1	o	650	ASN
1	p	325	GLN
1	p	457	GLN
1	p	485	GLN
1	p	497	ASN
1	p	518	ASN
1	p	598	GLN
1	p	650	ASN
1	q	325	GLN
1	q	449	ASN
1	q	457	GLN
1	q	485	GLN
1	q	497	ASN
1	q	518	ASN
1	q	598	GLN
1	q	650	ASN
1	r	325	GLN
1	r	457	GLN
1	r	485	GLN
1	r	497	ASN
1	r	518	ASN
1	r	598	GLN
1	r	650	ASN
1	s	325	GLN
1	s	457	GLN
1	s	485	GLN
1	s	497	ASN
1	s	518	ASN
1	s	598	GLN
1	s	650	ASN
1	t	325	GLN
1	t	457	GLN

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Mol	Chain	Res	Type
1	t	485	GLN
1	t	497	ASN
1	t	518	ASN
1	t	598	GLN
1	t	650	ASN
1	u	325	GLN
1	u	457	GLN
1	u	485	GLN
1	u	497	ASN
1	u	518	ASN
1	u	598	GLN
1	u	650	ASN
1	v	325	GLN
1	v	457	GLN
1	v	485	GLN
1	v	497	ASN
1	v	518	ASN
1	v	598	GLN
1	v	650	ASN
1	w	325	GLN
1	w	457	GLN
1	w	485	GLN
1	w	497	ASN
1	w	518	ASN
1	w	598	GLN
1	w	650	ASN
1	x	325	GLN
1	x	457	GLN
1	x	485	GLN
1	x	497	ASN
1	x	518	ASN
1	x	598	GLN
1	x	650	ASN
1	y	325	GLN
1	y	457	GLN
1	y	485	GLN
1	y	497	ASN
1	y	518	ASN
1	y	598	GLN
1	y	650	ASN
1	z	325	GLN
1	z	457	GLN

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Mol	Chain	Res	Type
1	z	485	GLN
1	z	497	ASN
1	z	518	ASN
1	z	598	GLN
1	z	650	ASN
1	z	690	ASN
1	1	325	GLN
1	1	457	GLN
1	1	485	GLN
1	1	497	ASN
1	1	518	ASN
1	1	598	GLN
1	1	650	ASN
1	2	325	GLN
1	2	457	GLN
1	2	485	GLN
1	2	497	ASN
1	2	518	ASN
1	2	598	GLN
1	2	650	ASN
1	3	325	GLN
1	3	457	GLN
1	3	485	GLN
1	3	497	ASN
1	3	518	ASN
1	3	598	GLN
1	3	650	ASN
1	4	325	GLN
1	4	457	GLN
1	4	485	GLN
1	4	497	ASN
1	4	518	ASN
1	4	598	GLN
1	4	650	ASN
1	5	325	GLN
1	5	457	GLN
1	5	485	GLN
1	5	497	ASN
1	5	518	ASN
1	5	598	GLN
1	5	650	ASN
1	6	325	GLN

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Mol	Chain	Res	Type
1	6	457	GLN
1	6	485	GLN
1	6	497	ASN
1	6	518	ASN
1	6	598	GLN
1	6	650	ASN
1	7	325	GLN
1	7	457	GLN
1	7	485	GLN
1	7	497	ASN
1	7	518	ASN
1	7	598	GLN
1	7	650	ASN
1	8	325	GLN
1	8	457	GLN
1	8	485	GLN
1	8	497	ASN
1	8	518	ASN
1	8	598	GLN
1	8	650	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 monosaccharides in this entry.

5.6 Ligand geometry

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.

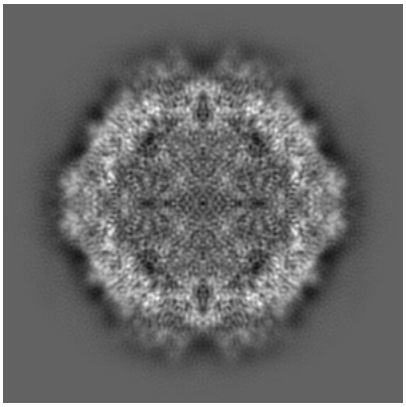
6 Map visualisation

This section contains visualisations of the EMDB entry EMD-8099. These allow visual inspection of the internal detail of the map and identification of artifacts.

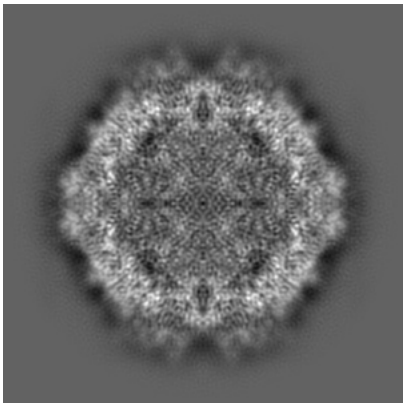
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections

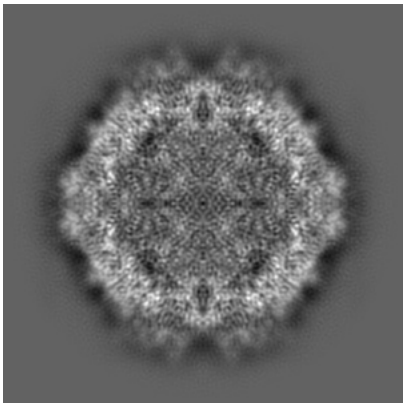
6.1.1 Primary map



X



Y

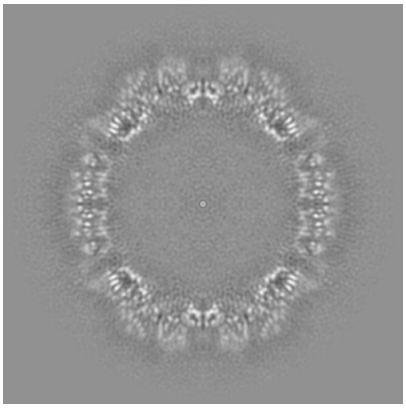


Z

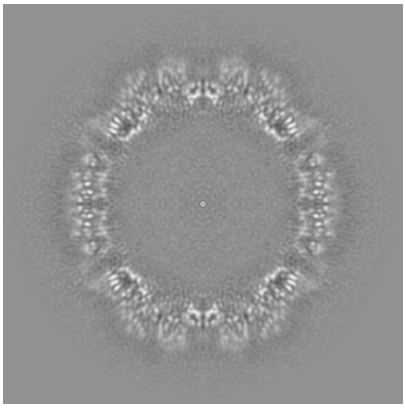
The images above show the map projected in three orthogonal directions.

6.2 Central slices

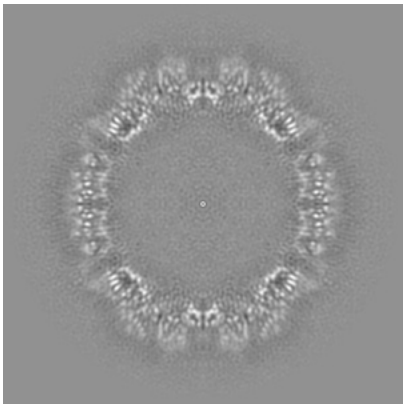
6.2.1 Primary map



X Index: 164



Y Index: 164

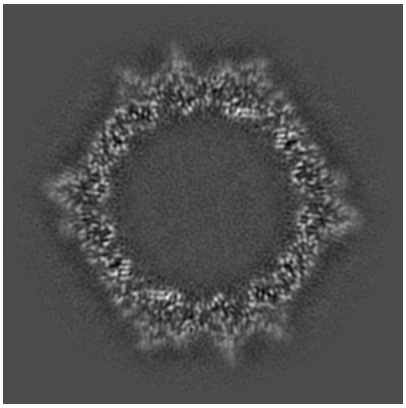


Z Index: 164

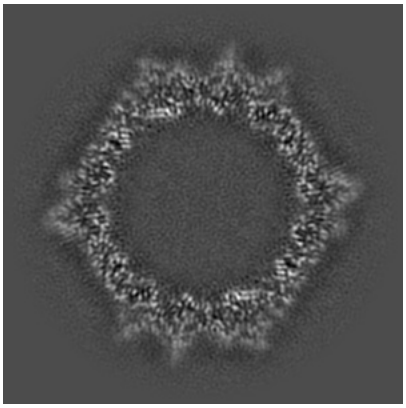
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices (i)

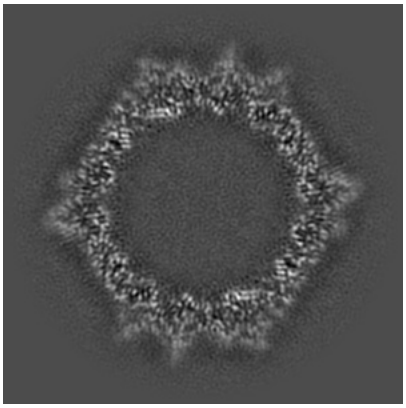
6.3.1 Primary map



X Index: 183



Y Index: 145

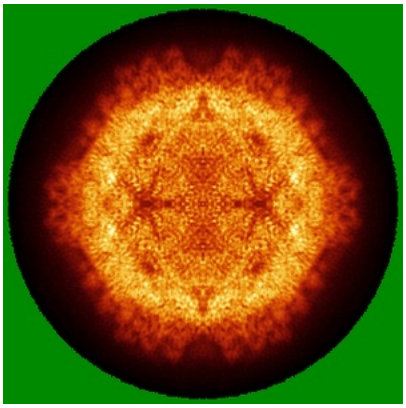


Z Index: 145

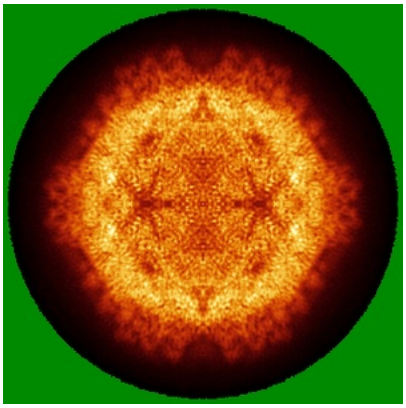
The images above show the largest variance slices of the map in three orthogonal directions.

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

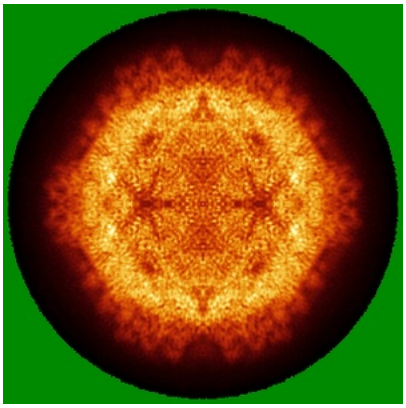
6.4.1 Primary map



X



Y

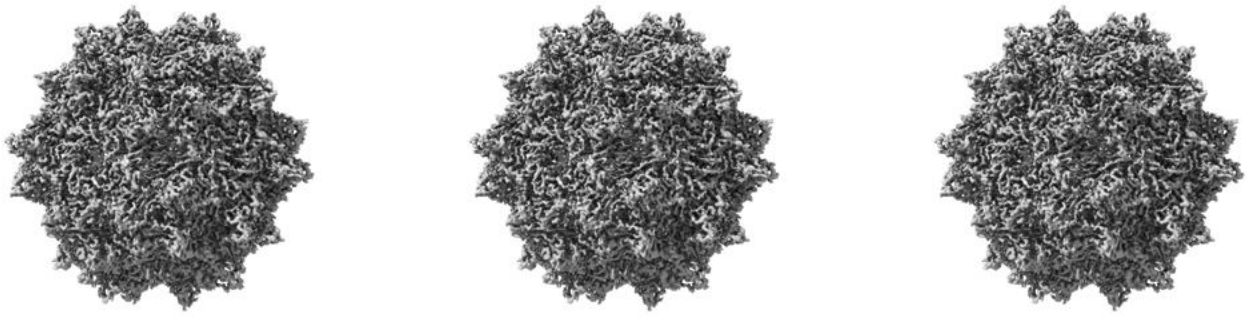


Z

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 2.9. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

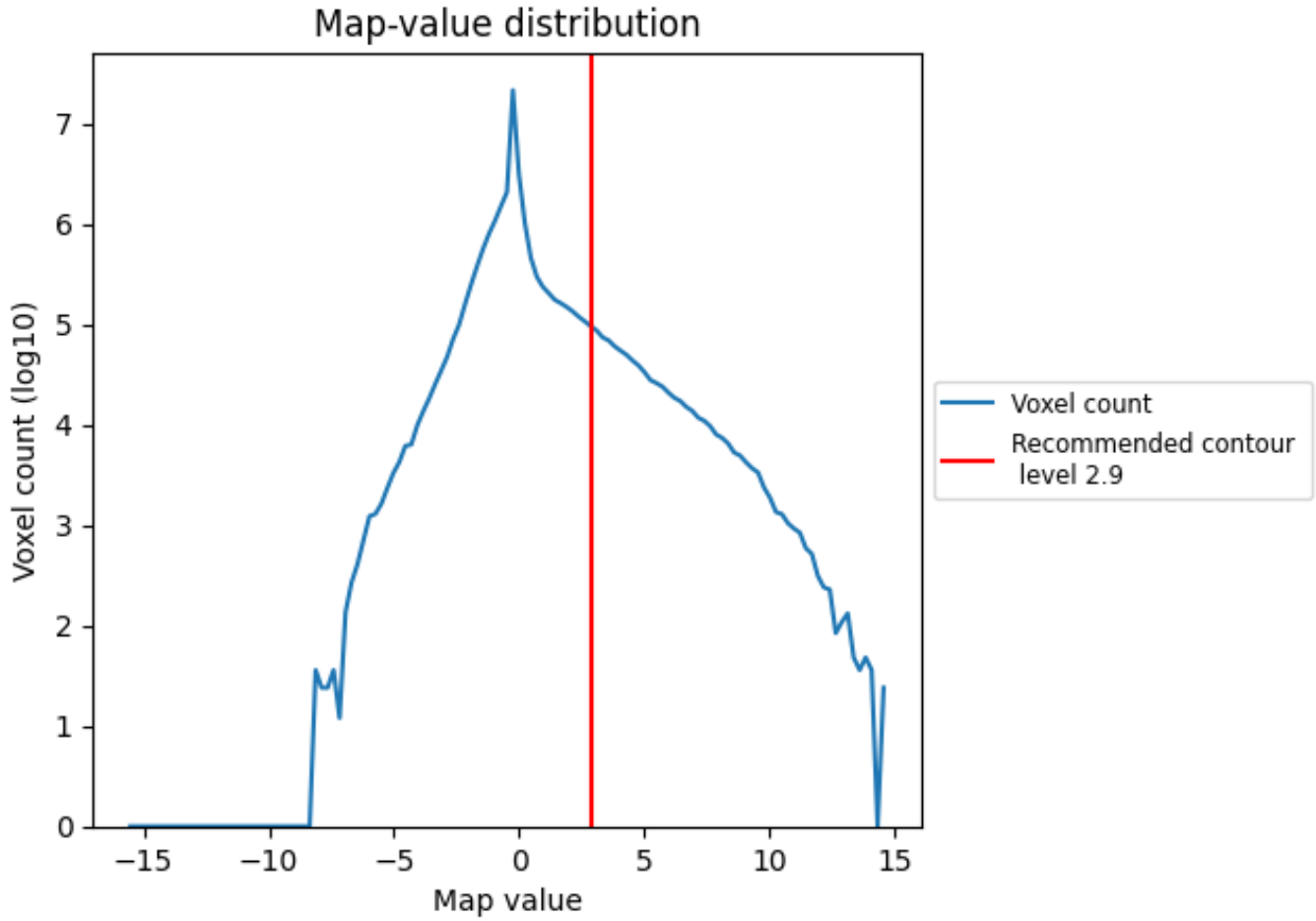
6.6 Mask visualisation (i)

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

7 Map analysis

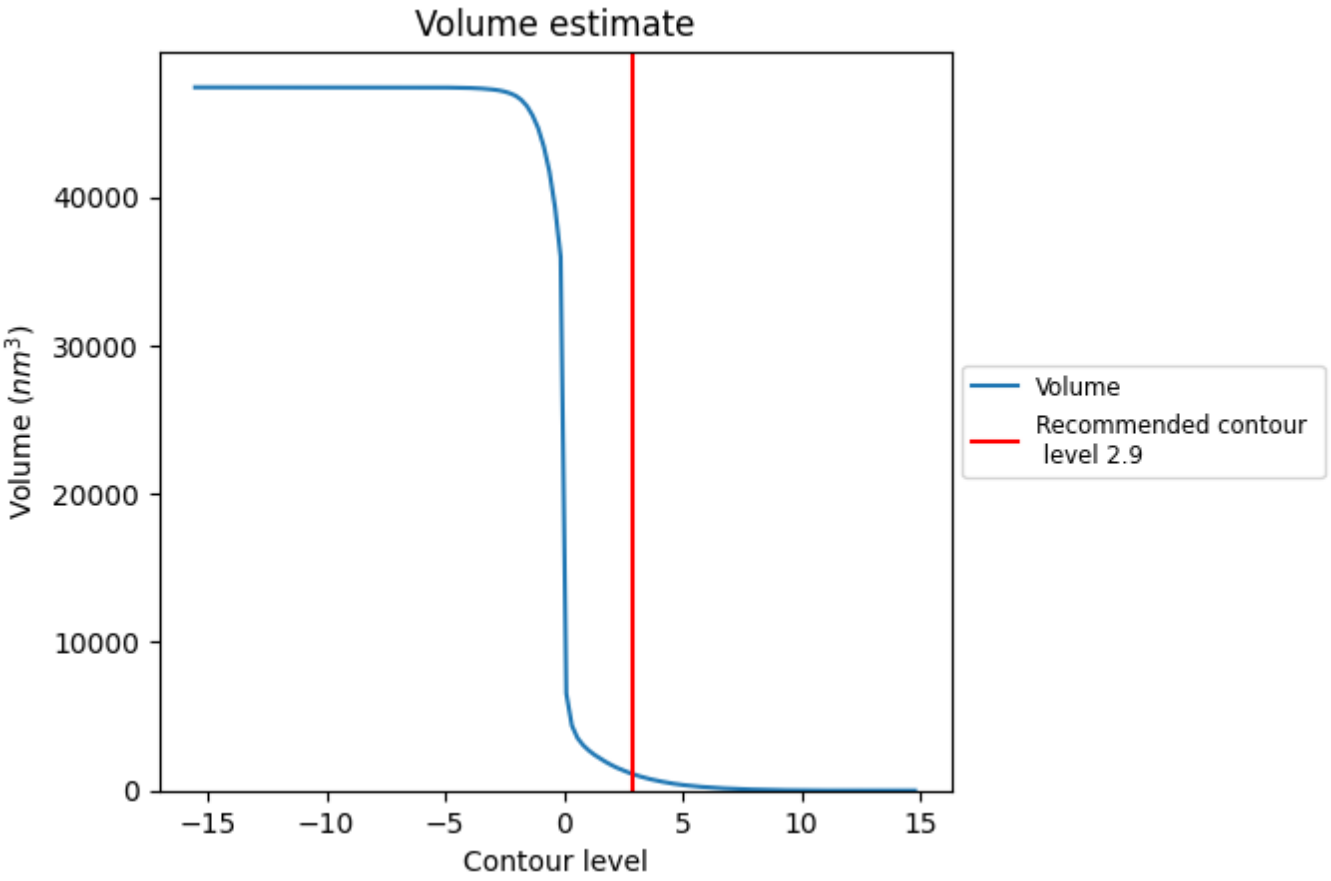
This section contains the results of statistical analysis of the map.

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.

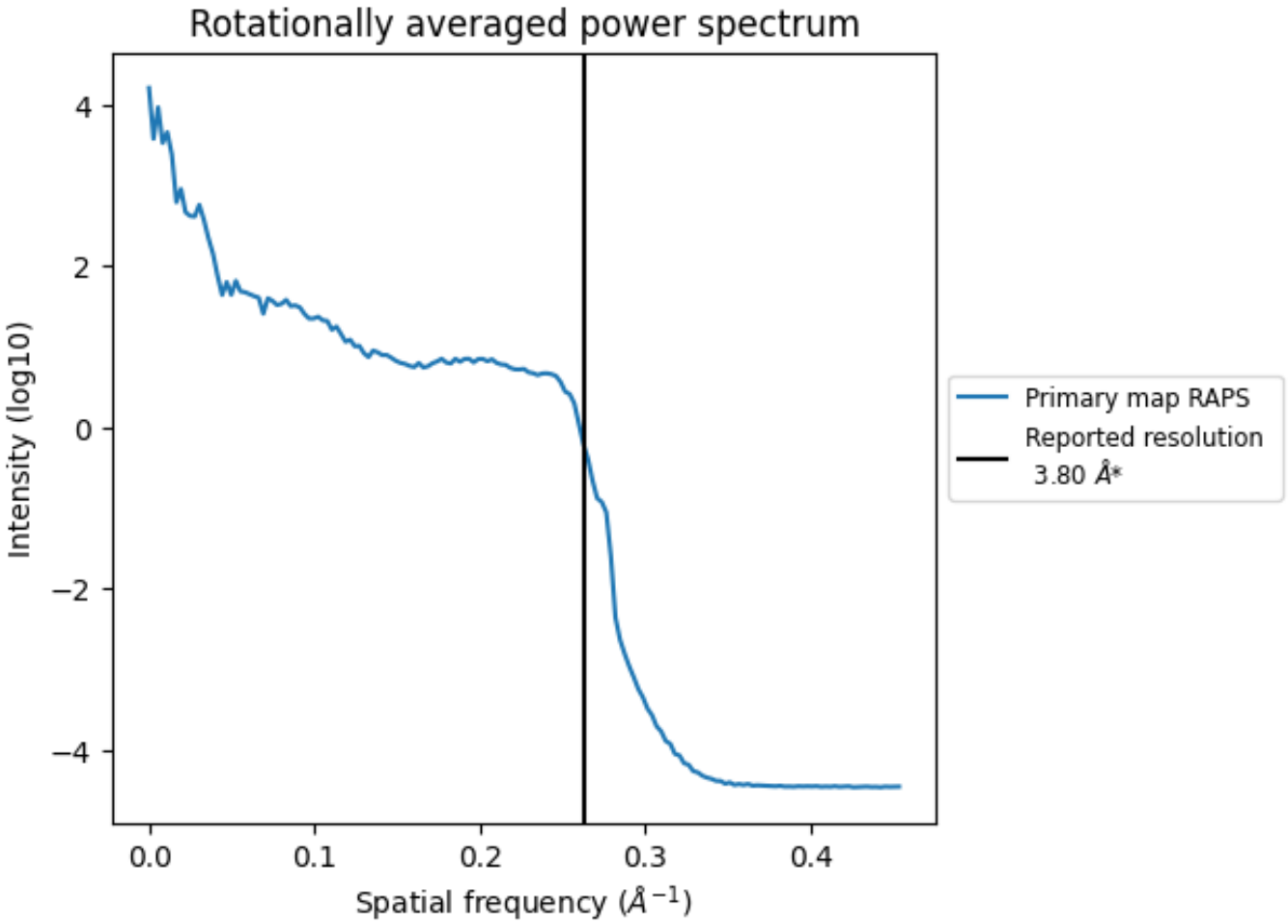
7.2 Volume estimate



The volume at the recommended contour level is 1115 nm^3; this corresponds to an approximate mass of 1007 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 0.263 Å^{-1}

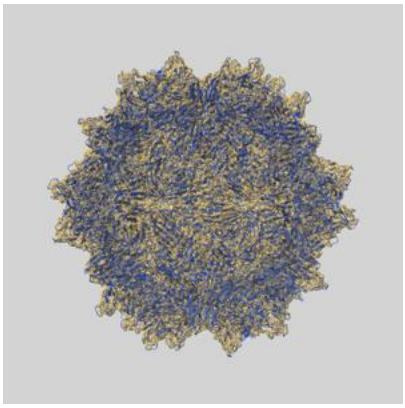
8 Fourier-Shell correlation (i)

This section was not generated. No FSC curve or half-maps provided.

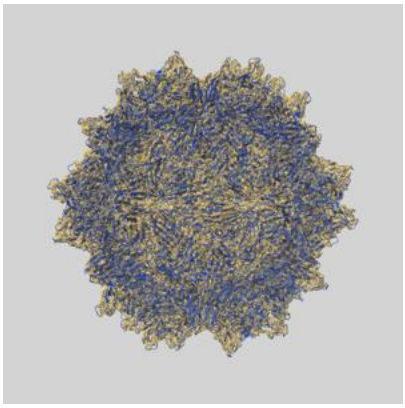
9 Map-model fit

This section contains information regarding the fit between EMDB map EMD-8099 and PDB model 5IPI. Per-residue inclusion information can be found in section 3 on page 11.

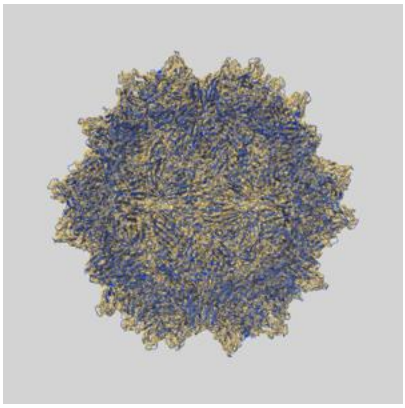
9.1 Map-model overlay



X



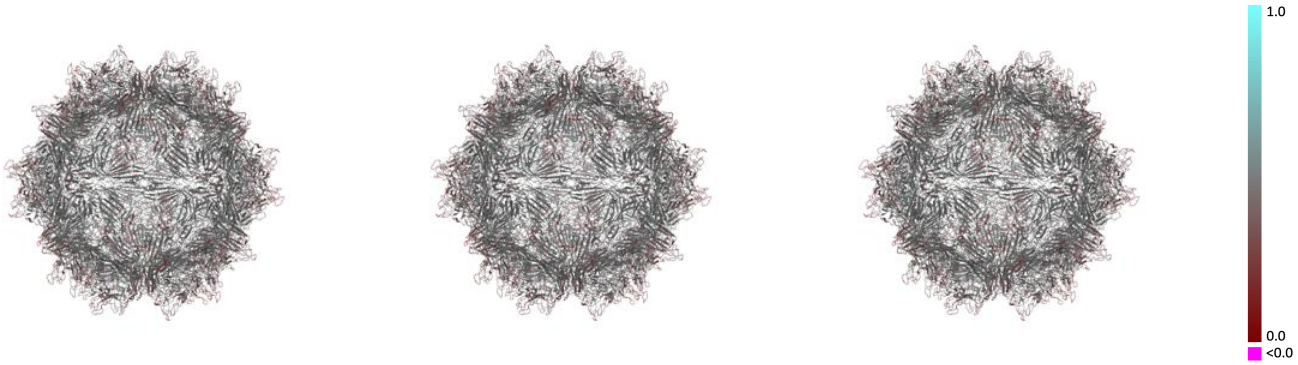
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 2.9 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.

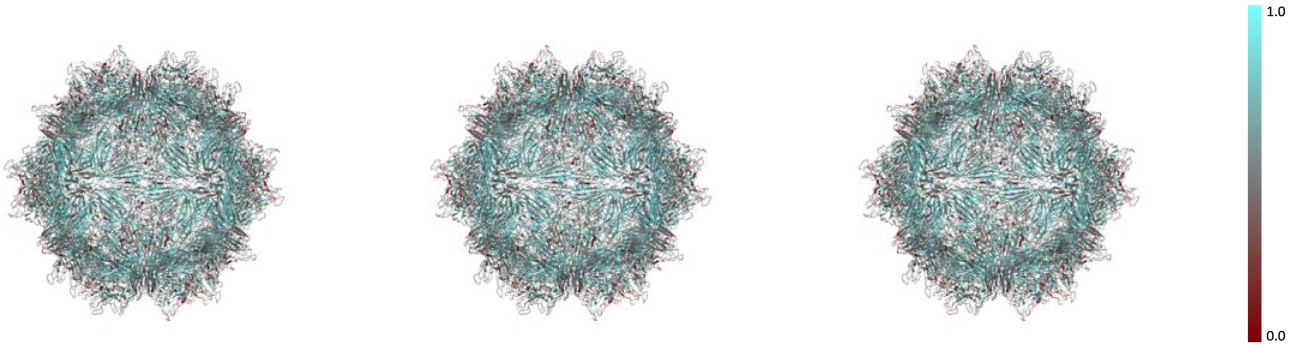
9.2 Q-score mapped to coordinate model (i)



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.

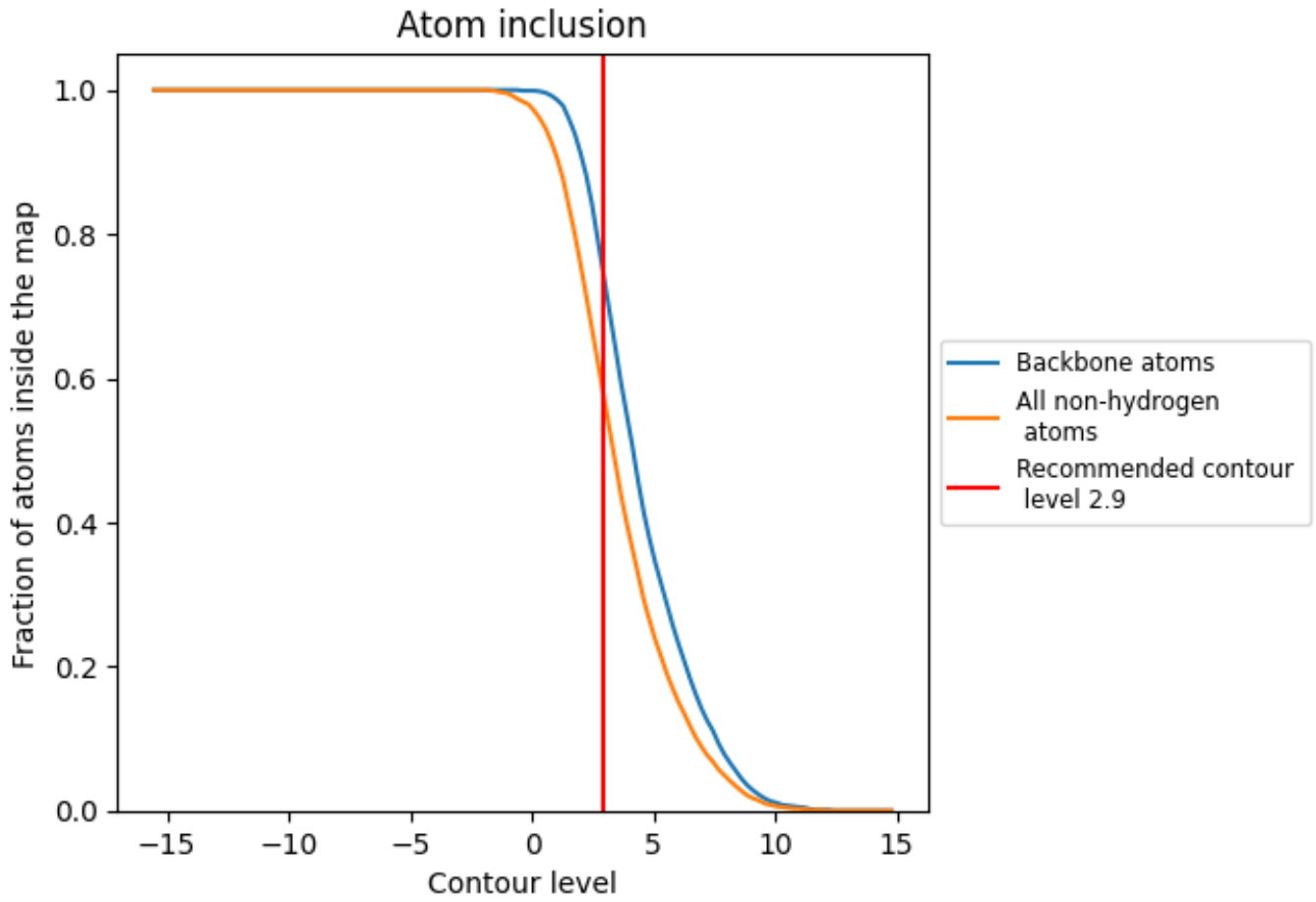
9.3 Atom inclusion mapped to coordinate model (i)



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 (2.9).

9.4 Atom inclusion (i)



At the recommended contour level, 75% of all backbone atoms, 58% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary

The table lists the average atom inclusion at the recommended contour level (2.9) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.5850	0.4410
1	0.5860	0.4430
2	0.5850	0.4410
3	0.5860	0.4400
4	0.5840	0.4400
5	0.5840	0.4420
6	0.5840	0.4420
7	0.5860	0.4420
8	0.5860	0.4410
A	0.5860	0.4420
B	0.5840	0.4410
C	0.5860	0.4410
D	0.5830	0.4410
E	0.5860	0.4420
F	0.5850	0.4410
G	0.5860	0.4420
H	0.5840	0.4410
I	0.5860	0.4410
J	0.5860	0.4410
K	0.5840	0.4420
L	0.5850	0.4410
M	0.5860	0.4410
N	0.5840	0.4400
O	0.5860	0.4400
P	0.5840	0.4410
Q	0.5860	0.4400
R	0.5860	0.4410
S	0.5850	0.4410
T	0.5860	0.4410
U	0.5850	0.4420
V	0.5840	0.4410
W	0.5860	0.4410
X	0.5840	0.4410
Y	0.5850	0.4400
Z	0.5850	0.4390

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Chain	Atom inclusion	Q-score
a	0.5840	0.4390
b	0.5840	0.4410
c	0.5860	0.4420
d	0.5860	0.4410
e	0.5840	0.4420
f	0.5860	0.4420
g	0.5840	0.4420
h	0.5850	0.4410
i	0.5840	0.4420
j	0.5860	0.4410
k	0.5860	0.4410
l	0.5860	0.4420
m	0.5850	0.4390
n	0.5830	0.4400
o	0.5860	0.4420
p	0.5850	0.4420
q	0.5860	0.4400
r	0.5850	0.4390
s	0.5830	0.4410
t	0.5860	0.4420
u	0.5840	0.4420
v	0.5870	0.4400
w	0.5840	0.4410
x	0.5840	0.4400
y	0.5830	0.4410
z	0.5870	0.4430