



## wwPDB EM Validation Summary Report ⓘ

Jul 3, 2024 – 12:05 pm BST

PDB ID : 6ZW5  
EMDB ID : EMD-11481  
Title : C15 symmetry: Bacterial Vipp1 and PspA are members of the ancient ESCRT-III membrane-remodeling superfamily.  
Authors : Liu, J.W.; Tassinari, M.; Souza, D.P.; Naskar, S.; Noel, J.K.; Bohuszewicz, O.; Buck, M.; Williams, T.A.; Baum, B.; Low, H.H.  
Deposited on : 2020-07-27  
Resolution : 7.00 Å(reported)

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

We welcome your comments at [validation@mail.wwpdb.org](mailto: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>.

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

EMDB validation analysis : 0.0.1.dev92  
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) : 2.37.1

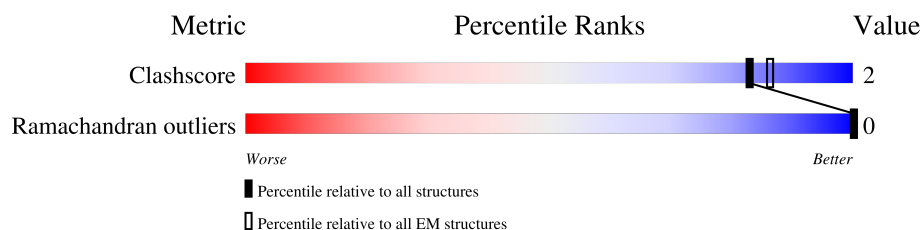
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*ELECTRON MICROSCOPY*

The reported resolution of this entry is 7.00 Å.

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



Metric	Whole archive (#Entries)	EM structures (#Entries)
Clashscore	158937	4297
Ramachandran outliers	154571	4023

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

Mol	Chain	Length	Quality of chain
1	A	258	<div> <div>7%</div> <div>52%</div> <div>48%</div> </div>
1	AA	258	<div> <div>12%</div> <div>83%</div> <div>15%</div> </div>
1	AB	258	<div> <div>9%</div> <div>83%</div> <div>15%</div> </div>
1	AC	258	<div> <div>36%</div> <div>61%</div> <div>39%</div> </div>
1	B	258	<div> <div>10%</div> <div>83%</div> <div>15%</div> </div>
1	BA	258	<div> <div>83%</div> <div>15%</div> </div>
1	BB	258	<div> <div>22%</div> <div>82%</div> <div>16%</div> </div>
1	BC	258	<div> <div>9%</div> <div>52%</div> <div>48%</div> </div>
1	C	258	<div> <div>83%</div> <div>15%</div> </div>

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Mol	Chain	Length	Quality of chain
1	CA	258	
1	CB	258	
1	CC	258	
1	D	258	
1	DA	258	
1	DB	258	
1	DC	258	
1	E	258	
1	EA	258	
1	EB	258	
1	EC	258	
1	F	258	
1	FA	258	
1	FB	258	
1	FC	258	
1	G	258	
1	GA	258	
1	GB	258	
1	GC	258	
1	H	258	
1	HA	258	
1	HB	258	
1	HC	258	
1	I	258	
1	IA	258	

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Mol	Chain	Length	Quality of chain
1	IB	258	
1	IC	258	
1	J	258	
1	JA	258	
1	JB	258	
1	JC	258	
1	K	258	
1	KA	258	
1	KB	258	
1	KC	258	
1	L	258	
1	LA	258	
1	LB	258	
1	LC	258	
1	M	258	
1	MA	258	
1	MB	258	
1	MC	258	
1	N	258	
1	NA	258	
1	NB	258	
1	O	258	
1	OA	258	
1	OB	258	
1	P	258	

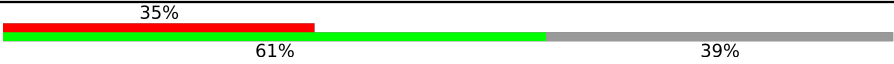

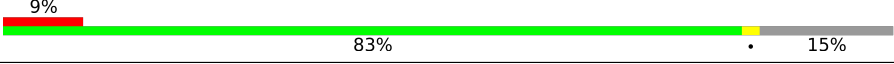
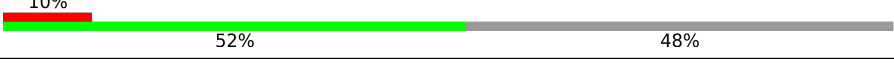

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Mol	Chain	Length	Quality of chain
1	PA	258	
1	PB	258	
1	Q	258	
1	QA	258	
1	QB	258	
1	R	258	
1	RA	258	
1	RB	258	
1	S	258	
1	SA	258	
1	SB	258	
1	T	258	
1	TA	258	
1	TB	258	
1	UA	258	
1	UB	258	
1	V	258	
1	VA	258	
1	VB	258	
1	W	258	
1	WA	258	
1	WB	258	
1	X	258	
1	XA	258	
1	XB	258	

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Mol	Chain	Length	Quality of chain
1	Y	258	
1	YA	258	
1	YB	258	
1	Z	258	
1	ZA	258	
1	ZB	258	

## 2 Entry composition [i](#)

There is only 1 type of molecule in this entry. The entry contains 86775 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 vipp1.

Mol	Chain	Residues	Atoms				AltConf	Trace
1	A	133	Total 661	C 395	N 133	O 133	0	0
1	B	219	Total 1086	C 648	N 219	O 219	0	0
1	C	219	Total 1086	C 648	N 219	O 219	0	0
1	D	219	Total 1086	C 648	N 219	O 219	0	0
1	E	218	Total 1081	C 645	N 218	O 218	0	0
1	F	158	Total 785	C 469	N 158	O 158	0	0
1	G	133	Total 661	C 395	N 133	O 133	0	0
1	H	219	Total 1086	C 648	N 219	O 219	0	0
1	I	219	Total 1086	C 648	N 219	O 219	0	0
1	J	219	Total 1086	C 648	N 219	O 219	0	0
1	K	218	Total 1081	C 645	N 218	O 218	0	0
1	L	158	Total 785	C 469	N 158	O 158	0	0
1	M	133	Total 661	C 395	N 133	O 133	0	0
1	N	219	Total 1086	C 648	N 219	O 219	0	0
1	O	219	Total 1086	C 648	N 219	O 219	0	0
1	P	219	Total 1086	C 648	N 219	O 219	0	0
1	Q	218	Total 1081	C 645	N 218	O 218	0	0

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Mol	Chain	Residues	Atoms				AltConf	Trace
1	R	158	Total 785	C 469	N 158	O 158	0	0
1	S	133	Total 661	C 395	N 133	O 133	0	0
1	T	219	Total 1086	C 648	N 219	O 219	0	0
1	V	219	Total 1086	C 648	N 219	O 219	0	0
1	W	219	Total 1086	C 648	N 219	O 219	0	0
1	X	218	Total 1081	C 645	N 218	O 218	0	0
1	Y	158	Total 785	C 469	N 158	O 158	0	0
1	Z	133	Total 661	C 395	N 133	O 133	0	0
1	AA	219	Total 1086	C 648	N 219	O 219	0	0
1	BA	219	Total 1086	C 648	N 219	O 219	0	0
1	CA	219	Total 1086	C 648	N 219	O 219	0	0
1	DA	218	Total 1081	C 645	N 218	O 218	0	0
1	EA	158	Total 785	C 469	N 158	O 158	0	0
1	FA	133	Total 661	C 395	N 133	O 133	0	0
1	GA	219	Total 1086	C 648	N 219	O 219	0	0
1	HA	219	Total 1086	C 648	N 219	O 219	0	0
1	IA	219	Total 1086	C 648	N 219	O 219	0	0
1	JA	218	Total 1081	C 645	N 218	O 218	0	0
1	KA	158	Total 785	C 469	N 158	O 158	0	0
1	LA	133	Total 661	C 395	N 133	O 133	0	0
1	MA	219	Total 1086	C 648	N 219	O 219	0	0

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Mol	Chain	Residues	Atoms				AltConf	Trace
1	NA	219	Total	C	N	O	0	0
			1086	648	219	219		
1	OA	219	Total	C	N	O	0	0
			1086	648	219	219		
1	PA	218	Total	C	N	O	0	0
			1081	645	218	218		
1	QA	158	Total	C	N	O	0	0
			785	469	158	158		
1	RA	133	Total	C	N	O	0	0
			661	395	133	133		
1	SA	219	Total	C	N	O	0	0
			1086	648	219	219		
1	TA	219	Total	C	N	O	0	0
			1086	648	219	219		
1	UA	219	Total	C	N	O	0	0
			1086	648	219	219		
1	VA	218	Total	C	N	O	0	0
			1081	645	218	218		
1	WA	158	Total	C	N	O	0	0
			785	469	158	158		
1	XA	133	Total	C	N	O	0	0
			661	395	133	133		
1	YA	219	Total	C	N	O	0	0
			1086	648	219	219		
1	ZA	219	Total	C	N	O	0	0
			1086	648	219	219		
1	AB	219	Total	C	N	O	0	0
			1086	648	219	219		
1	BB	218	Total	C	N	O	0	0
			1081	645	218	218		
1	CB	158	Total	C	N	O	0	0
			785	469	158	158		
1	DB	133	Total	C	N	O	0	0
			661	395	133	133		
1	EB	219	Total	C	N	O	0	0
			1086	648	219	219		
1	FB	219	Total	C	N	O	0	0
			1086	648	219	219		
1	GB	219	Total	C	N	O	0	0
			1086	648	219	219		
1	HB	218	Total	C	N	O	0	0
			1081	645	218	218		

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Mol	Chain	Residues	Atoms				AltConf	Trace
1	IB	158	Total 785	C 469	N 158	O 158	0	0
1	JB	133	Total 661	C 395	N 133	O 133	0	0
1	KB	219	Total 1086	C 648	N 219	O 219	0	0
1	LB	219	Total 1086	C 648	N 219	O 219	0	0
1	MB	219	Total 1086	C 648	N 219	O 219	0	0
1	NB	218	Total 1081	C 645	N 218	O 218	0	0
1	OB	158	Total 785	C 469	N 158	O 158	0	0
1	PB	133	Total 661	C 395	N 133	O 133	0	0
1	QB	219	Total 1086	C 648	N 219	O 219	0	0
1	RB	219	Total 1086	C 648	N 219	O 219	0	0
1	SB	219	Total 1086	C 648	N 219	O 219	0	0
1	TB	218	Total 1081	C 645	N 218	O 218	0	0
1	UB	158	Total 785	C 469	N 158	O 158	0	0
1	VB	133	Total 661	C 395	N 133	O 133	0	0
1	WB	219	Total 1086	C 648	N 219	O 219	0	0
1	XB	219	Total 1086	C 648	N 219	O 219	0	0
1	YB	219	Total 1086	C 648	N 219	O 219	0	0
1	ZB	218	Total 1081	C 645	N 218	O 218	0	0
1	AC	158	Total 785	C 469	N 158	O 158	0	0
1	BC	133	Total 661	C 395	N 133	O 133	0	0
1	CC	219	Total 1086	C 648	N 219	O 219	0	0

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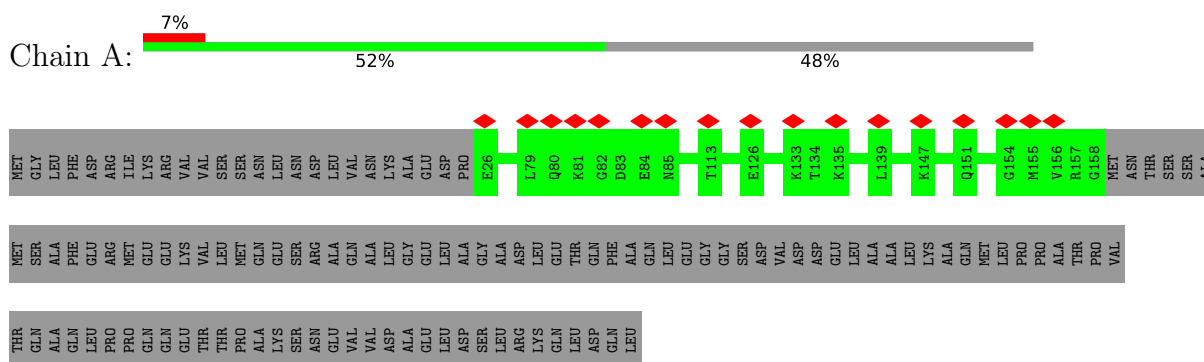
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Mol	Chain	Residues	Atoms				AltConf	Trace
1	DC	219	Total	C	N	O	0	0
			1086	648	219	219		
1	EC	219	Total	C	N	O	0	0
			1086	648	219	219		
1	FC	218	Total	C	N	O	0	0
			1081	645	218	218		
1	GC	158	Total	C	N	O	0	0
			785	469	158	158		
1	HC	133	Total	C	N	O	0	0
			661	395	133	133		
1	IC	219	Total	C	N	O	0	0
			1086	648	219	219		
1	JC	219	Total	C	N	O	0	0
			1086	648	219	219		
1	KC	219	Total	C	N	O	0	0
			1086	648	219	219		
1	LC	218	Total	C	N	O	0	0
			1081	645	218	218		
1	MC	158	Total	C	N	O	0	0
			785	469	158	158		

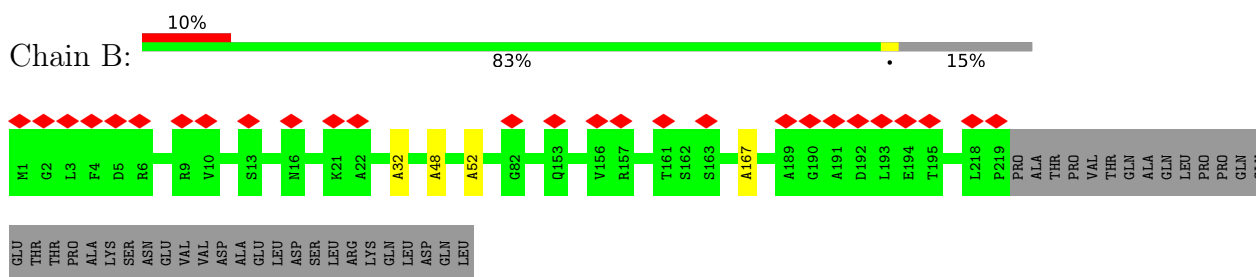
### 3 Residue-property plots [i](#)

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.

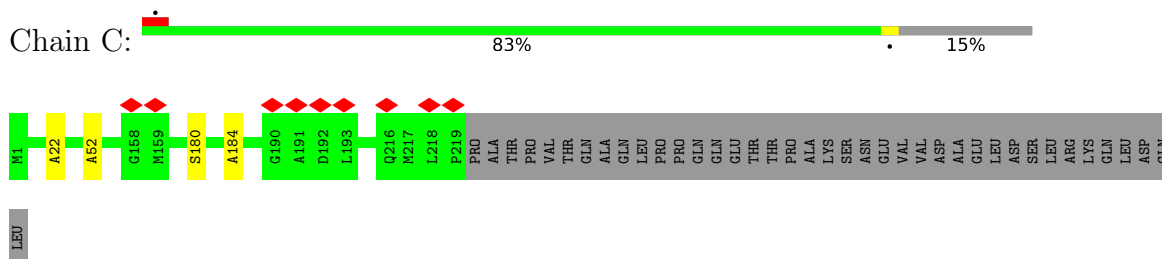
- Molecule 1: vipp1



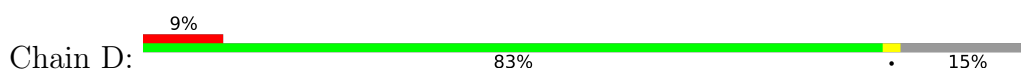
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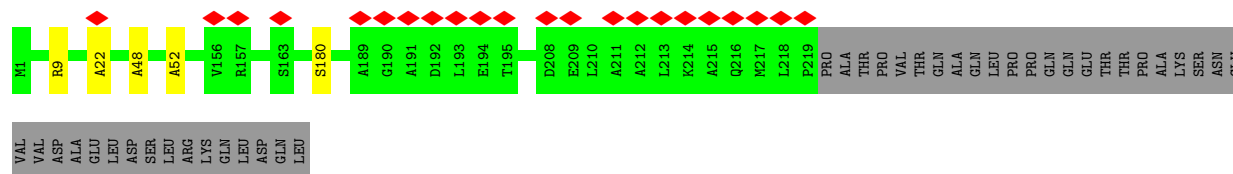


- Molecule 1: vipp1

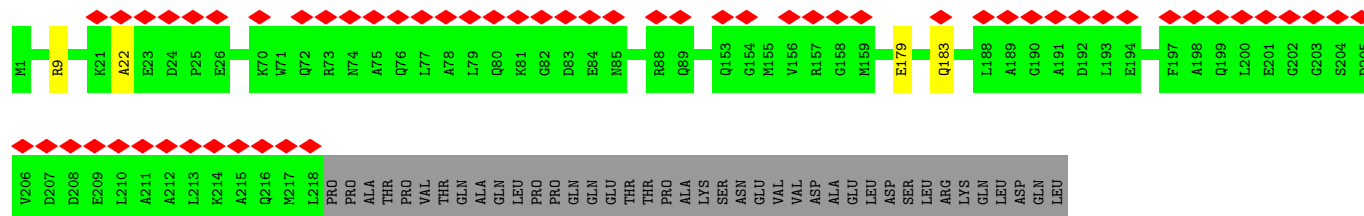
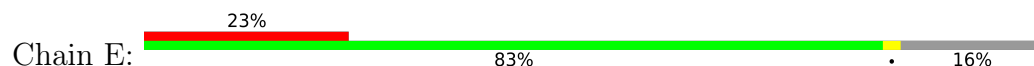


- Molecule 1: vipp1

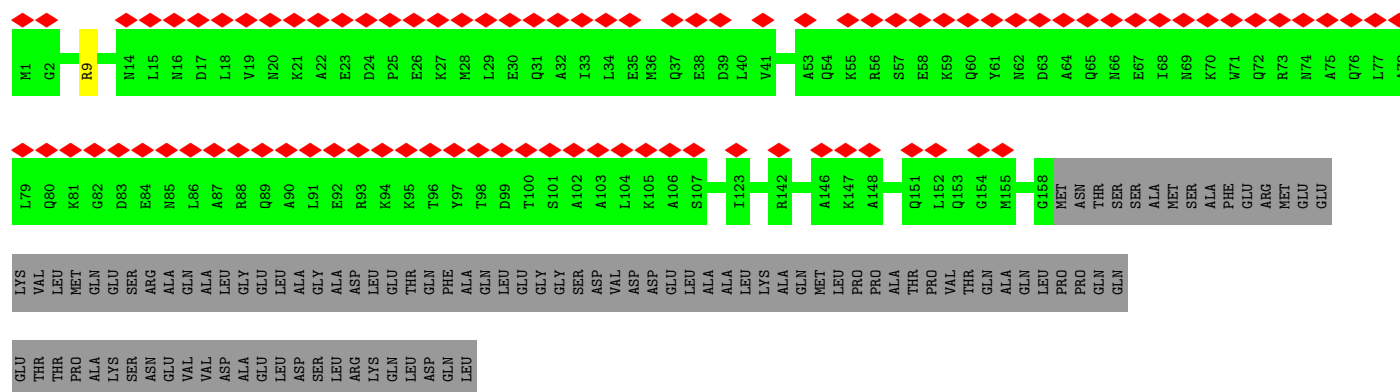




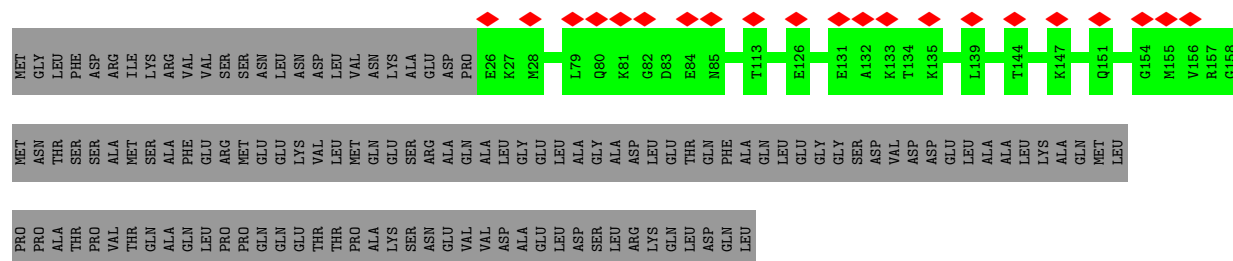
• Molecule 1: vipp1



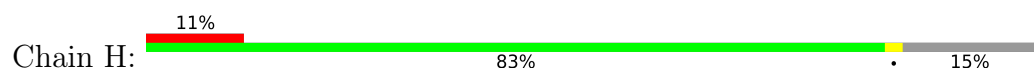
• Molecule 1: vipp1

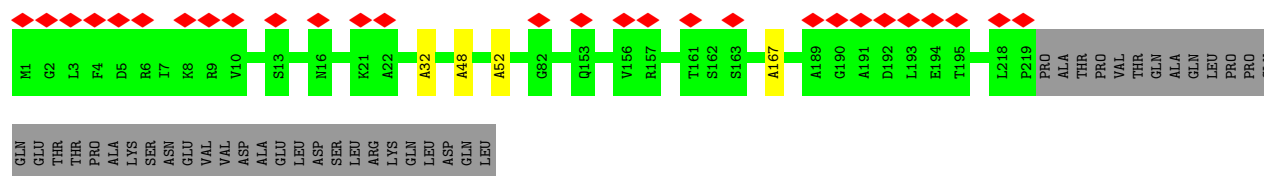


• Molecule 1: vipp1

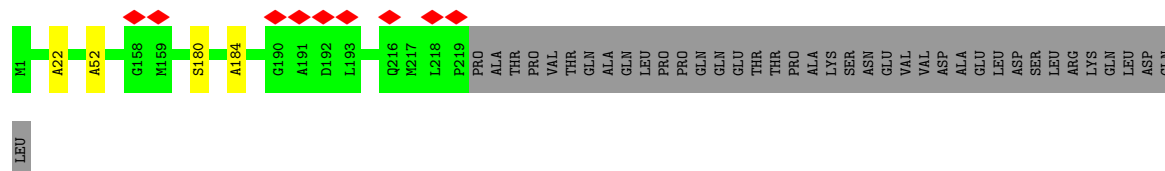
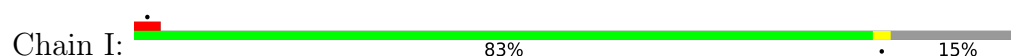


• Molecule 1: vipp1

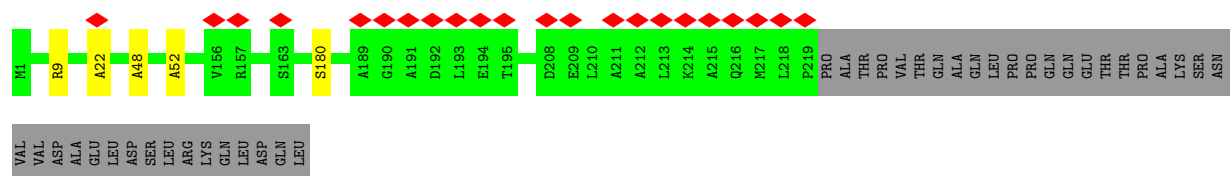
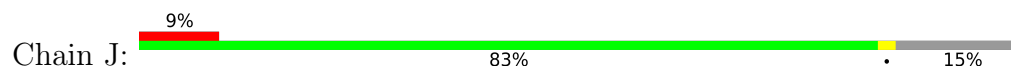




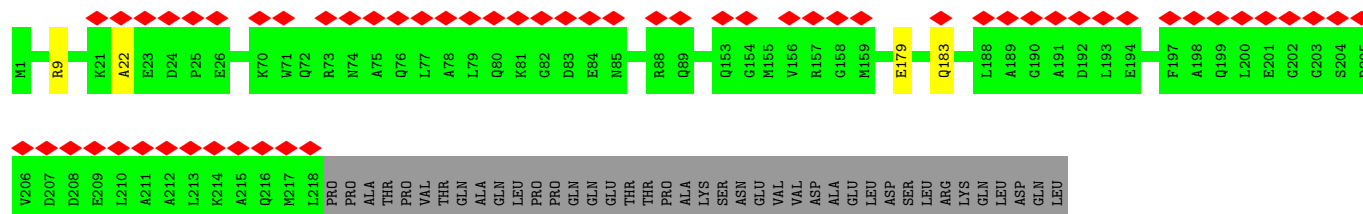
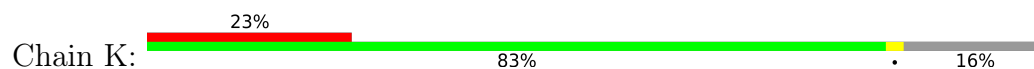
• Molecule 1: vipp1



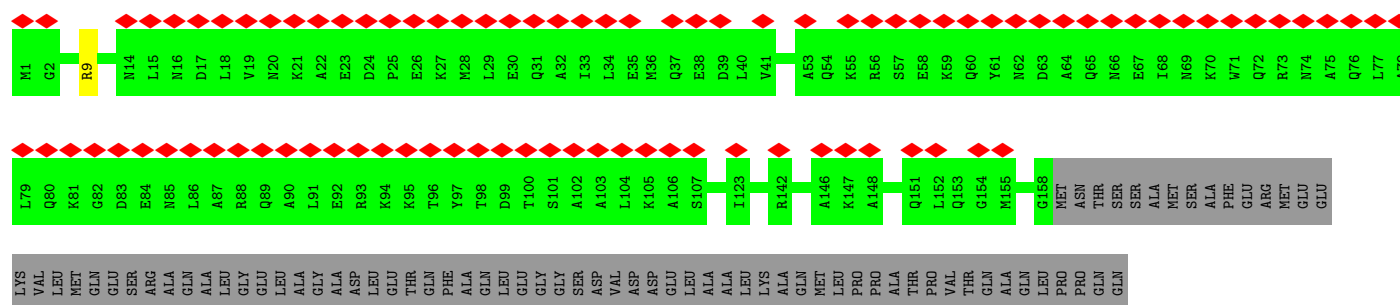
• Molecule 1: vipp1



• Molecule 1: vipp1



• Molecule 1: vipp1



GLU  
THR  
THR  
PRO  
PHE  
ALA  
LYS  
SER  
ASN  
GLU  
VAL  
VAL  
ASP  
ALA  
GLU  
LEU  
ASP  
SER  
SER  
ARG  
LYS  
GLN  
LEU  
ASP  
GLN  
LEU

• Molecule 1: vipp1


Chain M: 

MET  
GLY  
LEU  
PHE  
ASP  
MET  
ILE  
LYS  
ARG  
VAL  
VAL  
SER  
SER  
SER  
ASN  
LEU  
ASP  
SER  
LEU  
VAL  
ARG  
MET  
ASN  
PRO  
E26  
K27  
M28  
L79  
Q80  
K81  
G82  
D83  
E84  
N85  
L104  
T113  
E126  
T129  
S130  
E131  
A132  
K133  
T134  
K135  
K136  
E137  
M138  
L139  
T144  
K147  
Q151

G154  
M155  
V156  
R157  
G158  
MET  
THR  
THR  
SER  
SER  
ALA  
MET  
SER  
SER  
ALA  
GLN  
LEU  
PHE  
GLU  
MET  
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GLU  
LYS  
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PRO  
ALA  
THR  
PRO  
VAL  
THR  
GLN  
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THR  
ALA  
LYS  
SER  
ASN  
GLU  
VAL  
VAL  
ASP  
ALA  
GLU  
LEU  
LEU  
SER  
LEU  
GLN  
LEU


• Molecule 1: vipp1

Chain N: 

M1  
Q2  
L3  
F4  
D5  
R6  
I7  
K8  
R9  
V10  
V11  
S12  
S13  
M14  
L15  
N16  
K21  
A22  
A32  
A48  
A52  
G82  
Q153  
V156  
R157  
T161  
S162  
S163  
A167  
A189  
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A191  
D192  
L193  
E194  
T195  
L218  
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PRO  
ALA  
THR  
PRO  
VAL  
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ALA  
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• Molecule 1: vipp1

Chain O: 

M1  
A22  
A52  
G158  
M159  
S180  
A184  
G190  
A191  
D192  
L193  
Q216  
M217  
L218  
P219  
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THR  
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ALA  
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LEU  
ASP  
ASP  
ARG  
LYS  
GLN  
LEU  
ASP  
GLN

LEU


• Molecule 1: vipp1

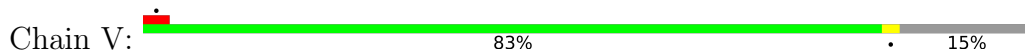
Chain P: 

M1  
R9  
A22  
A48  
A52  
V156  
R157  
S163  
S180  
A189  
G190  
A191  
D192  
L193  
E194  
T195  
D208  
E209  
L210  
A211  
A212  
L213  
K214  
A215  
Q216  
M217  
L218  
P219  
PRO  
ALA  
THR  
PRO  
VAL  
THR  
GLN  
ALA  
GLN  
LEU  
PRO  
PRO  
SER  
GLN  
LEU  
GLU  
THR  
THR  
PRO  
ALA  
LYS  
SER  
ASN  
GLU

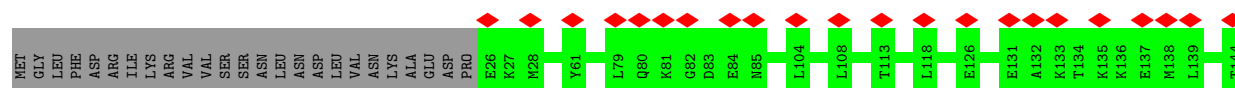
VAL  
VAL  
ASP  
GLU  
LEU  
ASP  
SER  
LEU  
ARG  
LYS  
GLN  
LEU  
ASP  
GLN  
LEU

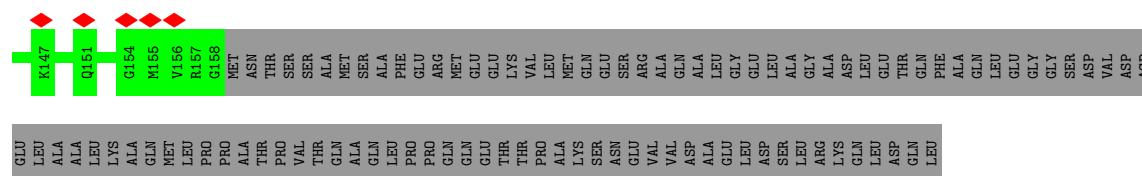
• Molecule 1: vipp1

Chain Q: 

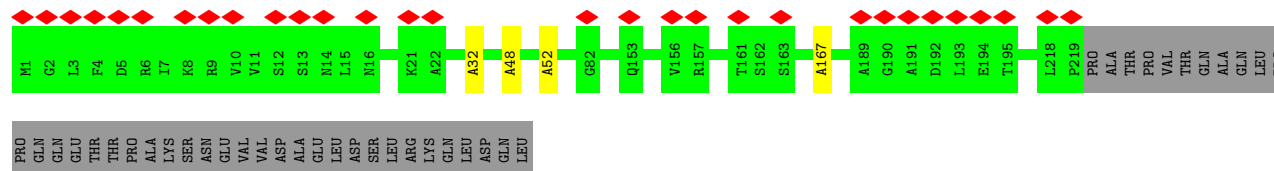
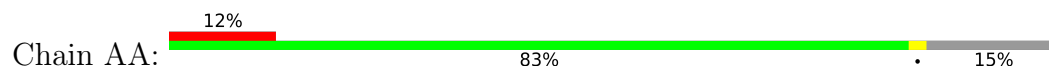




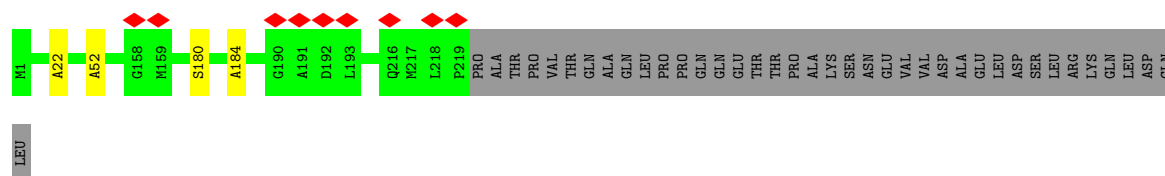
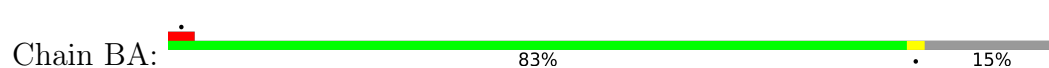




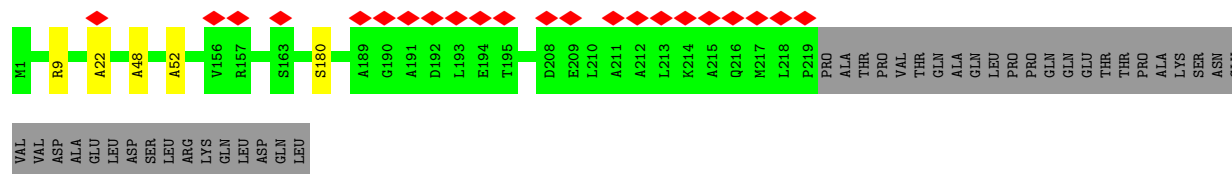
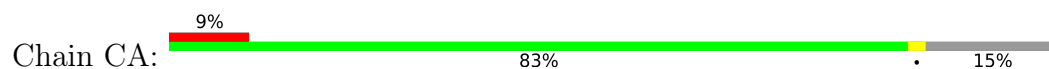
## • Molecule 1: vipp1



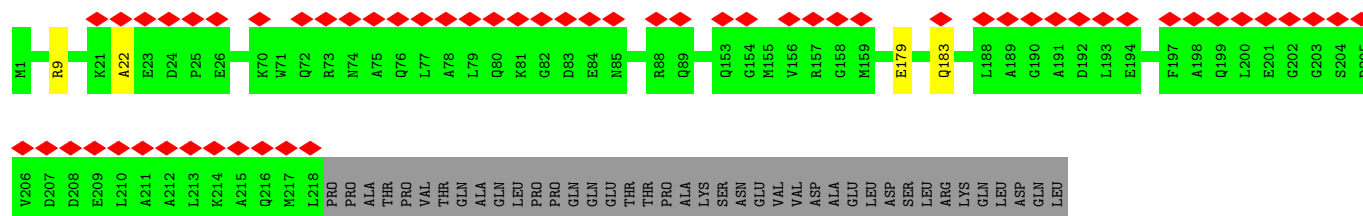
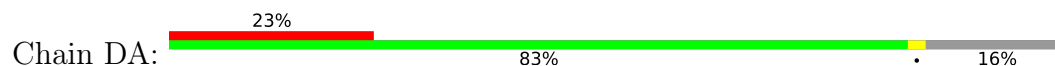
## • Molecule 1: vipp1



## • Molecule 1: vipp1

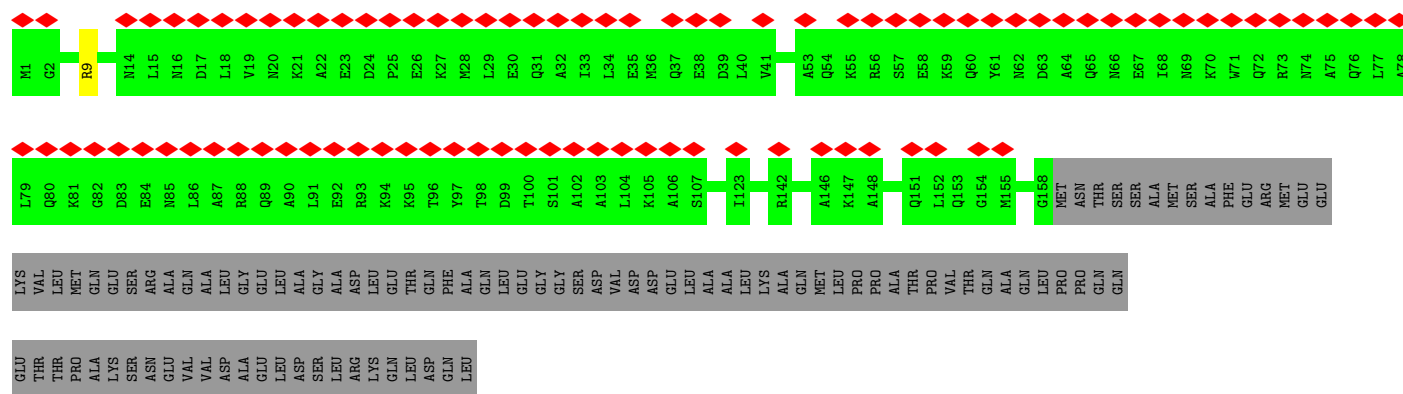


## • Molecule 1: vipp1



## • Molecule 1: vipp1

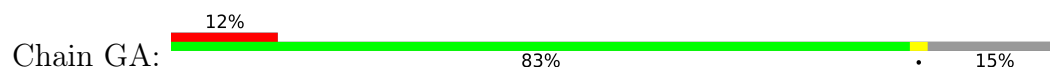




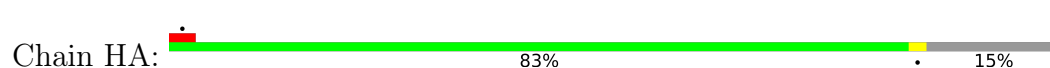
• Molecule 1: vipp1



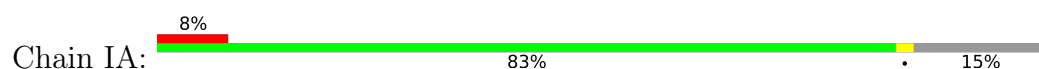
• Molecule 1: vipp1

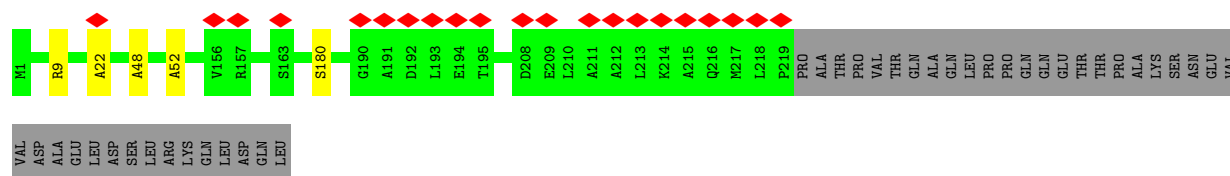


• Molecule 1: vipp1

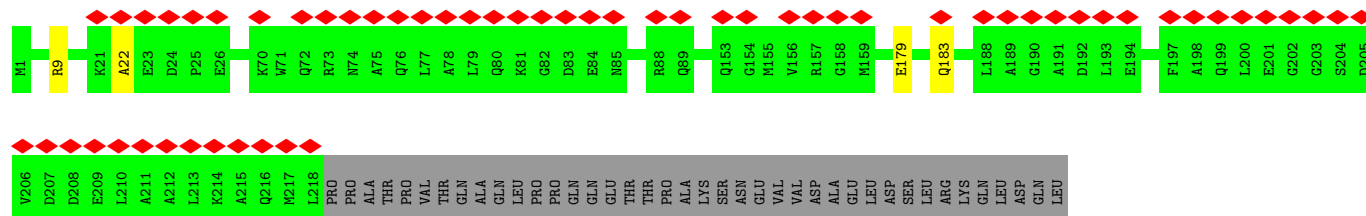
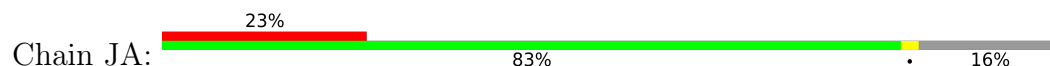


• Molecule 1: vipp1

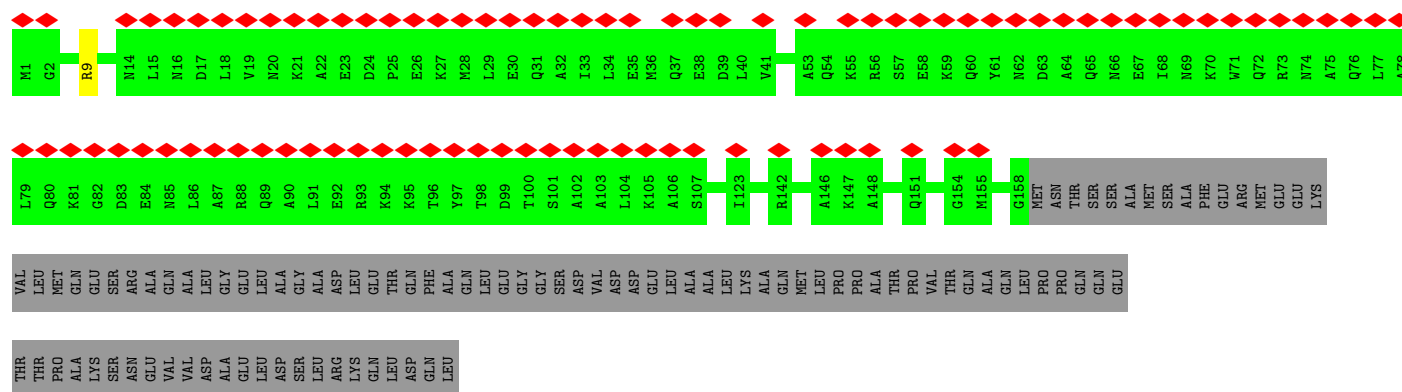




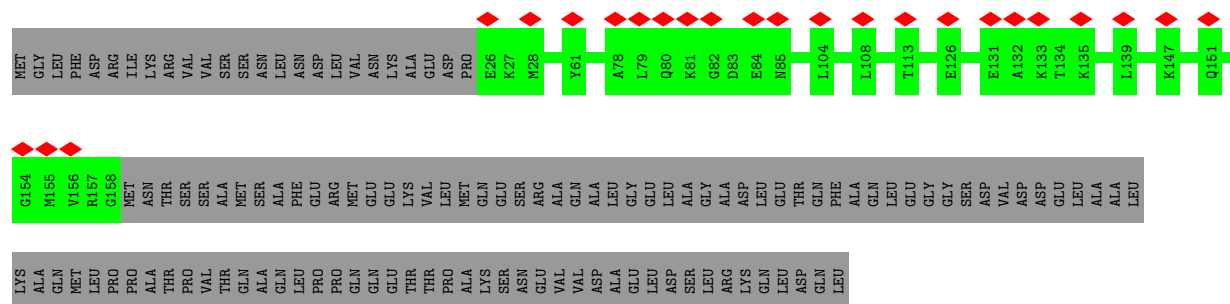
- Molecule 1: vipp1



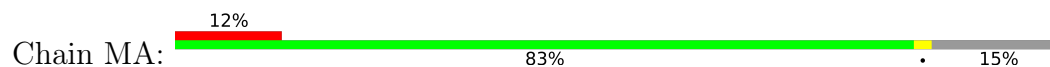
- Molecule 1: vipp1

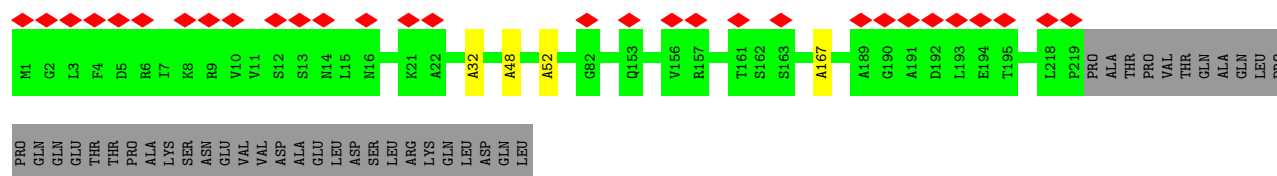


- Molecule 1: vipp1



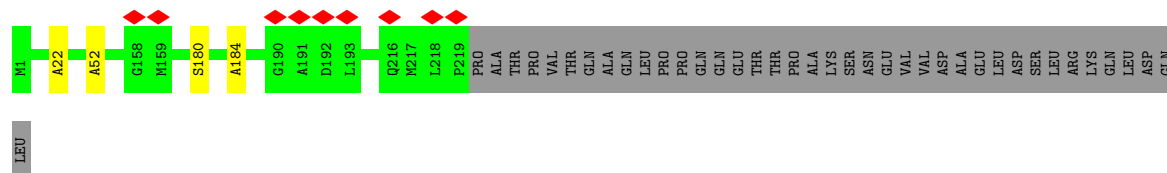
- Molecule 1: vipp1





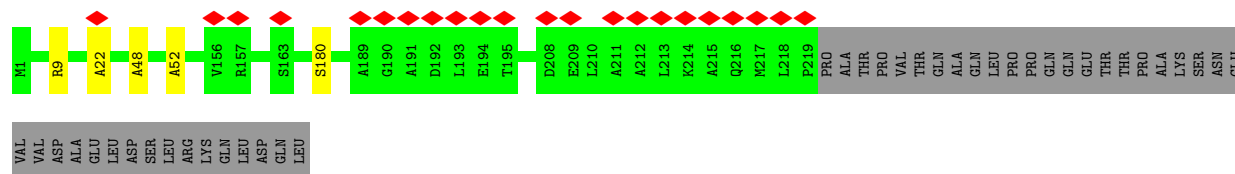
• Molecule 1: vipp1

Chain NA: 83% 15%



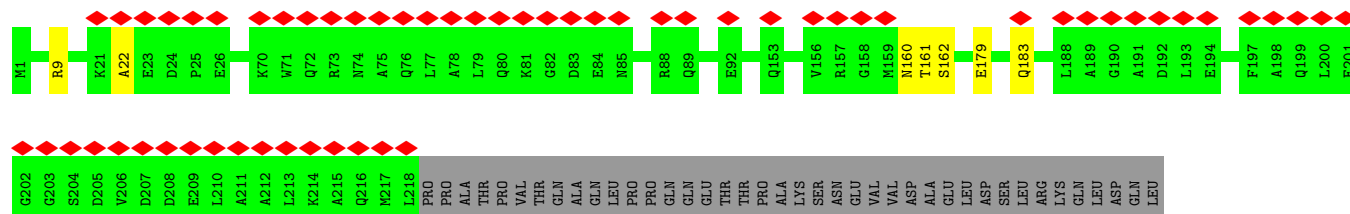
• Molecule 1: vipp1

Chain OA: 9% 83% 15%



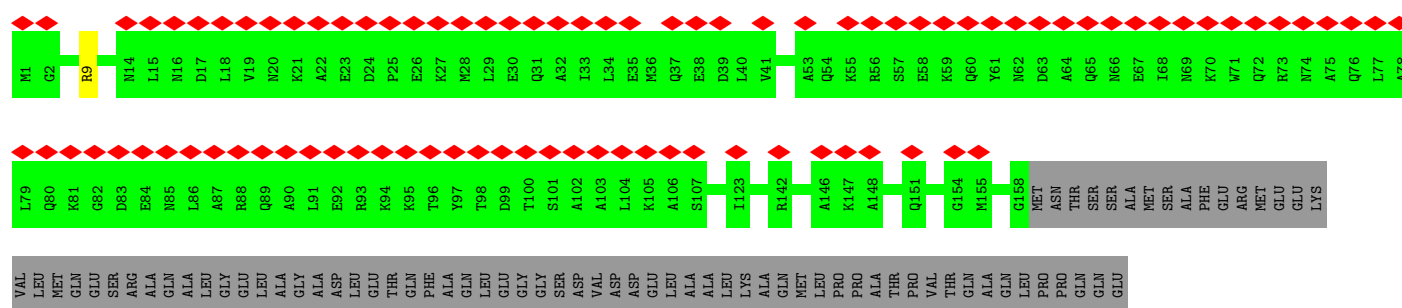
• Molecule 1: vipp1

Chain PA: 23% 82% 16%



• Molecule 1: vipp1

Chain QA: 35% 61% 39%



THR  
THR  
PRO  
ALA  
LYS  
SER  
ASN  
GLU  
VAL  
VAL  
ASP  
ALA  
GLU  
LEU  
ASP  
SER  
LEU  
ARG  
LYS  
GLN  
LEU  
ASP  
GLN  
LEU

• Molecule 1: vipp1

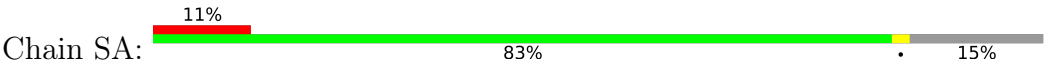


MET  
GLY  
PHE  
ASP  
MET  
ILE  
LYS  
ARG  
VAL  
MET  
GLU  
SER  
SER  
ASN  
LEU  
ASN  
ASP  
MET  
LEU  
VAL  
GLN  
SER  
LYS  
ASN  
GLN  
LEU  
ASP  
PRO  
E26  
L79  
Q80  
K91  
G82  
D83  
E84  
N85  
L104  
T113  
E126  
E131  
A132  
K133  
T134  
K135  
L139  
K147  
Q151  
G154  
M155  
V156  
R157  
G158  
MET  
ASN

THR  
SER  
SER  
ALA  
MET  
SER  
MET  
PHE  
GLN  
ARG  
MET  
GLU  
GLU  
GLN  
LYS  
VAL  
THR  
LEU  
THR  
PRO  
MET  
GLN  
GLU  
SER  
ARG  
GLN  
ALA  
GLN  
VAL  
LEU  
ASP  
GLY  
GLU  
LEU  
ALA  
GLY  
GLU  
GLY  
GLY  
SER  
ASP  
VAL  
ASP  
ASP  
GLU  
LEU  
ALA  
ALA  
LYS  
ALA  
GLN  
MET  
LEU  
PRO  
PRO

ALA  
THR  
PRO  
VAL  
THR  
GLN  
ALA  
GLN  
PHE  
GLU  
PRO  
PRO  
GLN  
GLN  
GLU  
THR  
THR  
PRO  
ALA  
GLN  
SER  
SER  
ASN  
GLN  
VAL  
VAL  
ASP  
ALA  
GLY  
GLU  
LEU  
SER  
SER  
LEU  
ARG  
LYS  
GLN  
LEU  
ASP  
GLN  
LEU

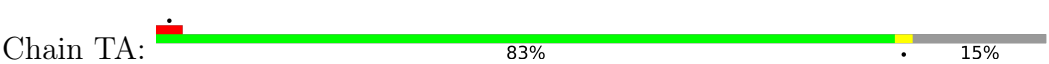
• Molecule 1: vipp1



M1  
Q2  
L3  
F4  
D5  
R6  
I7  
K8  
R9  
V10  
S13  
M16  
K21  
A22  
A32  
A48  
A52  
G82  
Q153  
V156  
R157  
T161  
S162  
S163  
A167  
A189  
G190  
A191  
D192  
L193  
E194  
T195  
L218  
P219  
PRO  
ALA  
THR  
PRO  
VAL  
THR  
GLN  
ALA  
GLN  
LEU  
PRO  
GLN

GLN  
GLU  
THR  
THR  
PRO  
PRO  
LYS  
ALA  
SER  
SER  
GLU  
VAL  
VAL  
ASP  
ALA  
GLU  
LEU  
ASP  
SER  
SER  
LEU  
ARG  
LYS  
GLN  
LEU  
ASP  
GLN  
LEU

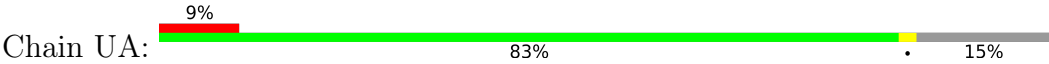
• Molecule 1: vipp1



M1  
A22  
A52  
G158  
M159  
S180  
A184  
G190  
A191  
D192  
L193  
Q216  
M217  
L218  
P219  
PRO  
THR  
THR  
PRO  
VAL  
THR  
GLN  
GLN  
GLN  
LEU  
PRO  
PRO  
GLN  
GLU  
THR  
THR  
PRO  
ALA  
ALA  
LYS  
SER  
ASN  
GLU  
VAL  
VAL  
ASP  
ASP  
GLU  
LEU  
ASP  
SER  
LEU  
ARG  
GLN  
GLN  
LEU  
GLN

LEU

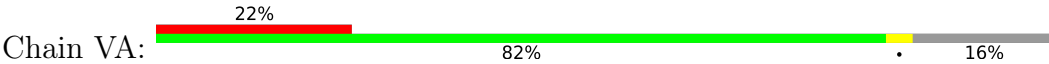
• Molecule 1: vipp1

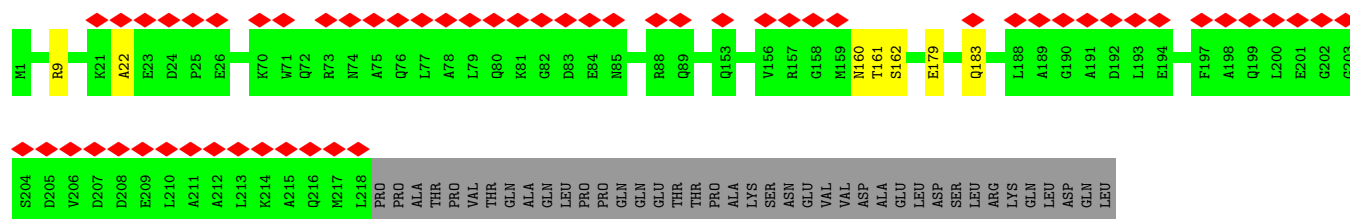


M1  
R9  
A22  
A48  
A52  
V156  
R157  
S163  
S180  
A189  
G190  
A191  
D192  
L193  
E194  
T195  
D208  
E209  
L210  
A211  
A212  
L213  
K214  
A215  
Q216  
M217  
L218  
P219  
PRO  
ALA  
THR  
PRO  
VAL  
THR  
GLN  
GLN  
GLN  
LEU  
PRO  
PRO  
GLN  
GLN  
GLU  
THR  
THR  
PRO  
ALA  
LYS  
SER  
ASN  
GLU

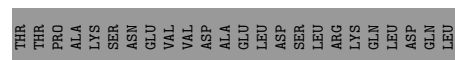
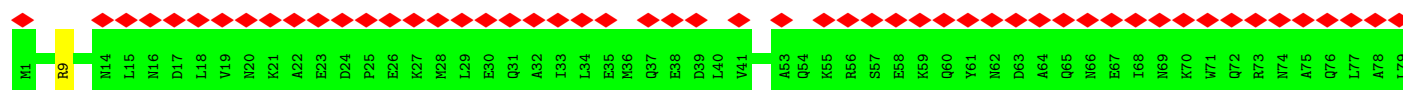
VAL  
VAL  
ASP  
ALA  
GLU  
LEU  
ASP  
SER  
LEU  
LYS  
ARG  
GLN  
GLN  
ASP  
GLN  
LEU

• Molecule 1: vipp1

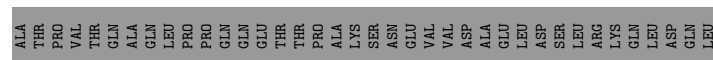
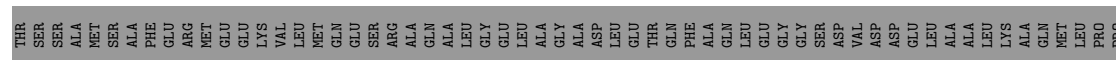
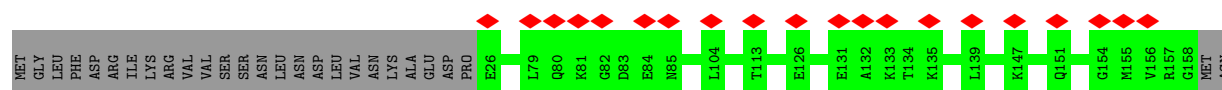




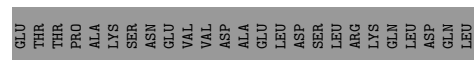
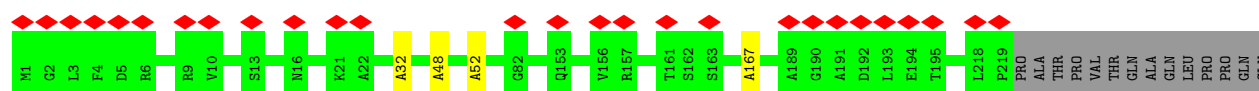
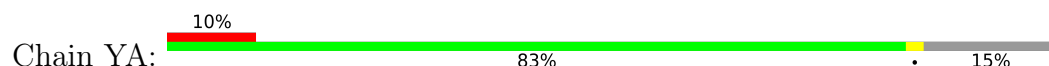
• Molecule 1: vipp1



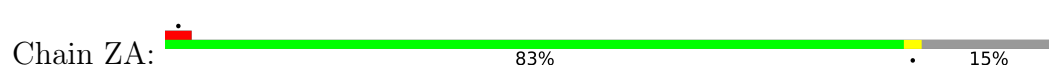
• Molecule 1: vipp1

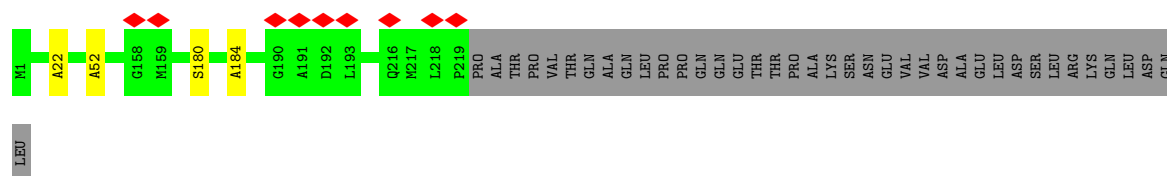


• Molecule 1: vipp1

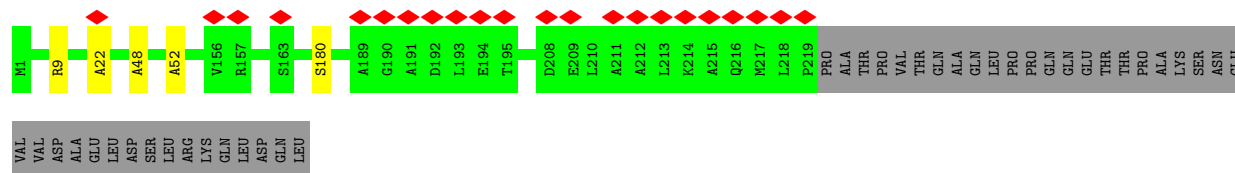
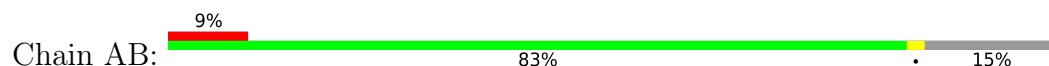


• Molecule 1: vipp1

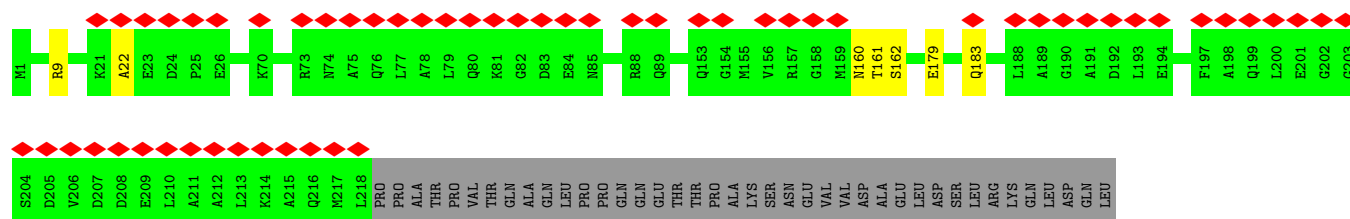
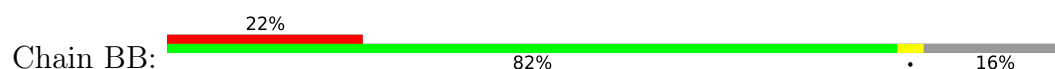




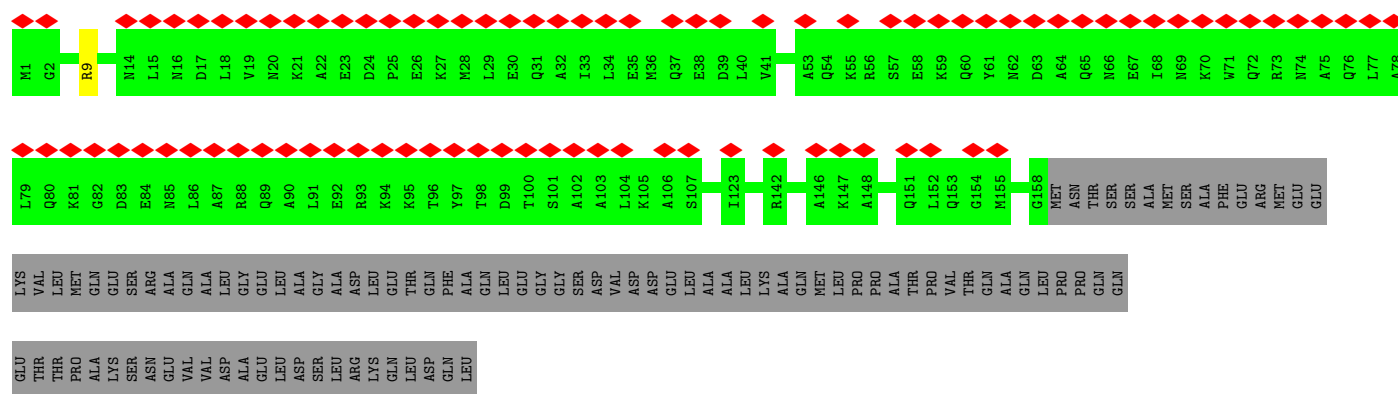
- Molecule 1: vipp1



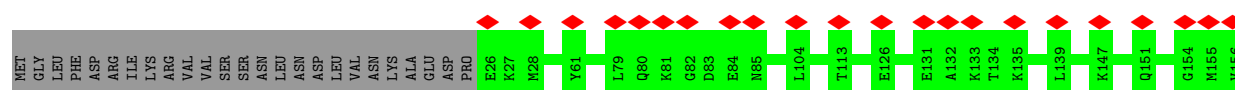
- Molecule 1: vipp1



- Molecule 1: vipp1



- Molecule 1: vipp1





R157  
G158  
MET  
ASN  
THR  
SER  
SER  
ALA  
MET  
SER  
ALA  
PHE  
GLU  
ARG  
MET  
GLU  
GLU  
LYS  
VAL  
MET  
GLN  
GLU  
SER  
ARG  
ALA  
GLN  
ALA  
LEU  
GLY  
GLU  
LEU  
ALA  
GLY  
LEU  
GLU  
THR  
GLN  
PHE  
ALA  
GLN  
LEU  
GLU  
GLY  
GLY  
SER  
SER  
VAL  
ASP  
ASP  
GLU  
LEU  
ALA  
ALA  
LEU  
LYS  
GLN

MET  
LEU  
PRO  
PRO  
ALA  
THR  
THR  
PRO  
VAL  
THR  
GLN  
ALA  
GLN  
LEU  
PRO  
PRO  
GLN  
GLN  
GLU  
LYS  
THR  
THR  
PRO  
PRO  
ALA  
LYS  
ASN  
SER  
ASN  
GLU  
VAL  
VAL  
ASP  
GLU  
ALA  
GLU  
LEU  
ASP  
SER  
LEU  
ARG  
LYS  
GLN  
LEU  
GLN  
LEU

• Molecule 1: vipp1

Chain EB: 11% 83% 15%

M1  
G2  
L3  
F4  
D5  
R6  
R9  
V10  
S13  
N14  
L15  
N16  
K21  
A22  
A32  
A48  
A52  
G82  
Q153  
V156  
R157  
T161  
S162  
S163  
A167  
A189  
G190  
A191  
D192  
L193  
E194  
T195  
L218  
P219  
PRO  
ALA  
THR  
PRO  
VAL  
THR  
GLN  
ALA  
LEU  
PRO  
GLN

GLN  
GLU  
THR  
PRO  
ALA  
LYS  
SER  
ASN  
VAL  
VAL  
ASP  
ALA  
GLU  
LEU  
ASP  
SER  
LEU  
ARG  
LYS  
GLN  
LEU  
GLN  
LEU

• Molecule 1: vipp1

Chain FB: 83% 15%

M1  
A22  
A52  
G158  
M159  
S180  
A184  
G190  
A191  
D192  
L193  
Q216  
M217  
L218  
P219  
PRO  
ALA  
THR  
PRO  
VAL  
THR  
GLN  
GLN  
ALA  
LEU  
PRO  
PRO  
GLN  
GLU  
THR  
THR  
PRO  
ALA  
LYS  
SER  
ASN  
GLU  
VAL  
VAL  
ASP  
ALA  
GLU  
LEU  
SER  
SER  
LEU  
ARG  
LYS  
GLN  
LEU  
GLN

LEU

• Molecule 1: vipp1

Chain GB: 9% 83% 15%

M1  
R9  
A22  
A48  
A52  
V156  
R157  
S163  
S180  
A189  
G190  
A191  
D192  
L193  
E194  
T195  
D208  
E209  
L210  
A211  
A212  
L213  
K214  
A215  
Q216  
M217  
L218  
P219  
PRO  
THR  
PRO  
VAL  
THR  
GLN  
GLN  
GLN  
LEU  
PRO  
PRO  
GLN  
GLU  
THR  
PRO  
ALA  
LYS  
SER  
ASN  
GLU

VAL  
VAL  
ASP  
ALA  
GLU  
LEU  
ASP  
SER  
ARG  
LYS  
GLN  
ASP  
GLN  
LEU

• Molecule 1: vipp1

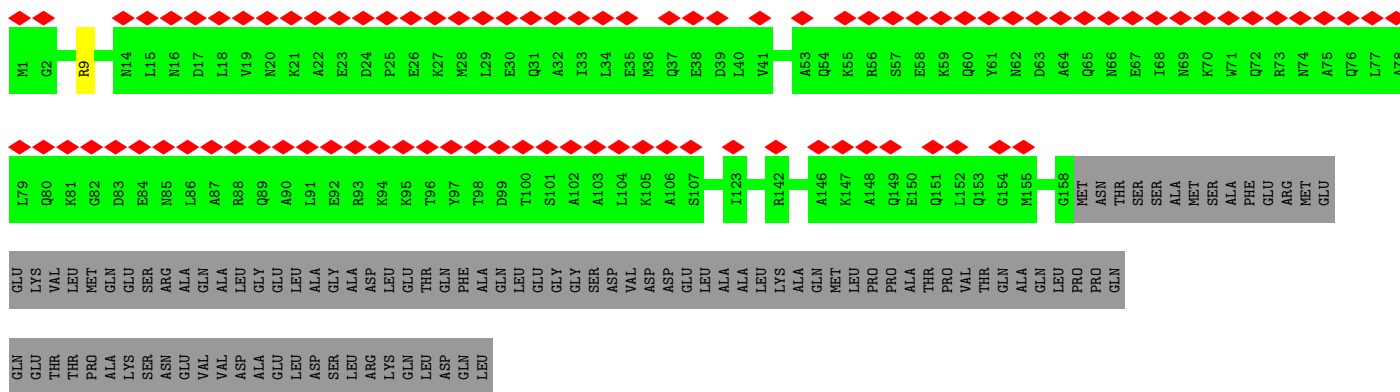
Chain HB: 22% 82% 16%

M1  
R9  
K21  
A22  
E23  
D24  
P25  
E26  
K70  
R73  
N74  
A75  
Q76  
L77  
A78  
L79  
Q80  
K81  
G82  
D83  
E84  
N85  
R88  
Q89  
Q153  
G154  
M155  
V156  
R157  
G158  
M159  
N160  
T161  
S162  
E179  
Q183  
L188  
A189  
G190  
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D192  
L193  
E194  
F197  
A198  
Q199  
L200  
E201  
G202  
G203

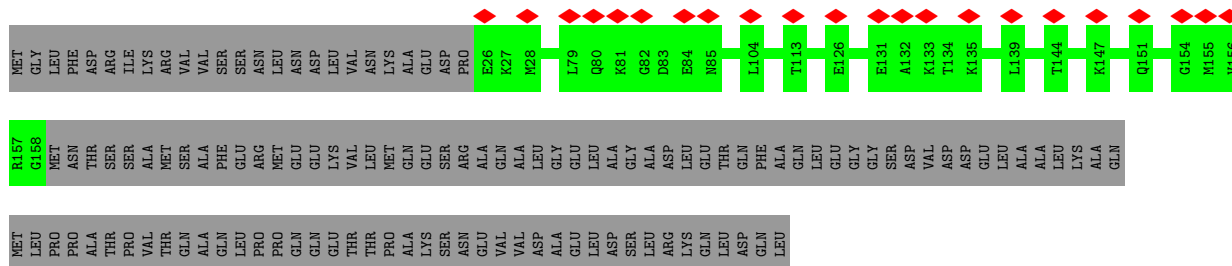
S204  
D205  
V206  
D207  
D208  
E209  
L210  
A211  
A212  
L213  
K214  
A215  
Q216  
M217  
L218  
PRO  
PRO  
ALA  
THR  
PRO  
VAL  
THR  
GLN  
GLN  
ALA  
LEU  
PRO  
PRO  
GLN  
GLN  
GLU  
THR  
THR  
PRO  
ALA  
LYS  
SER  
ASN  
GLU  
VAL  
VAL  
ASP  
ALA  
GLU  
LEU  
ASP  
SER  
LEU  
LEU  
ARG  
LYS  
GLN  
LEU  
ASP  
GLN  
LEU

• Molecule 1: vipp1

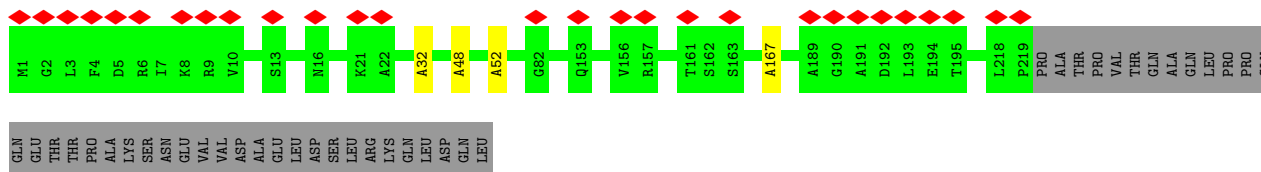
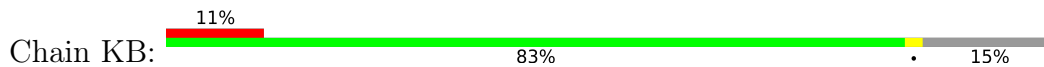
Chain IB: 36% 61% 39%



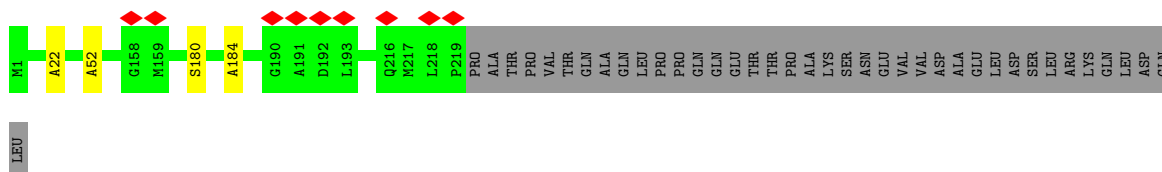
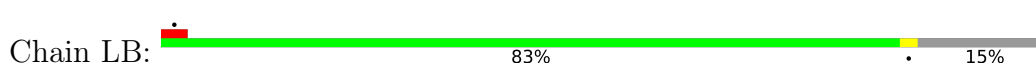
- Molecule 1: vippp1



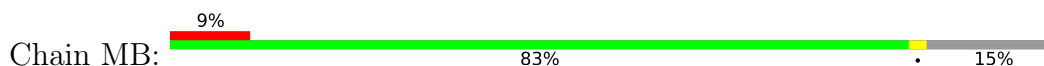
- Molecule 1: vippp1

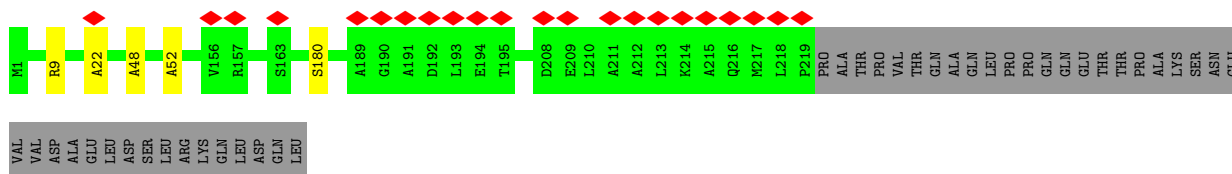


- Molecule 1: vippp1

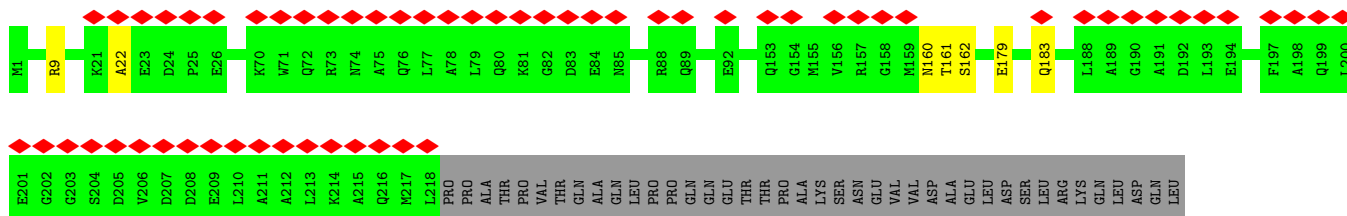
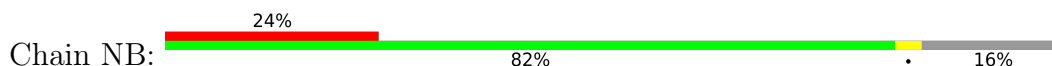


- Molecule 1: vippp1

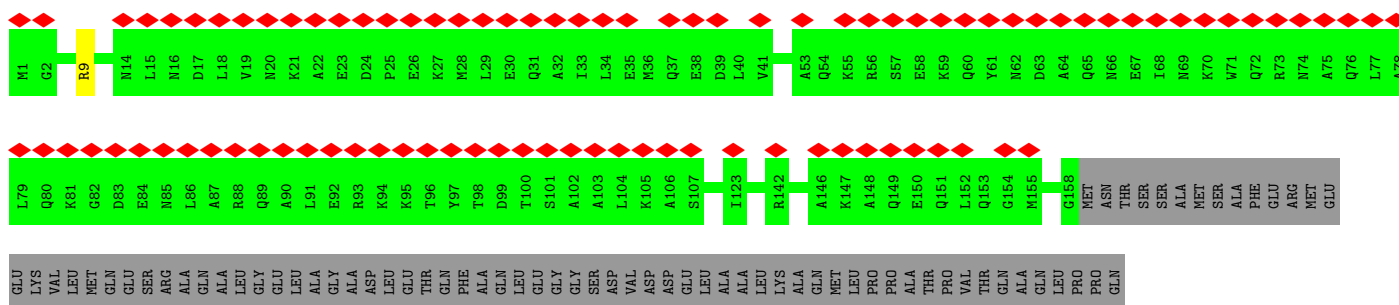




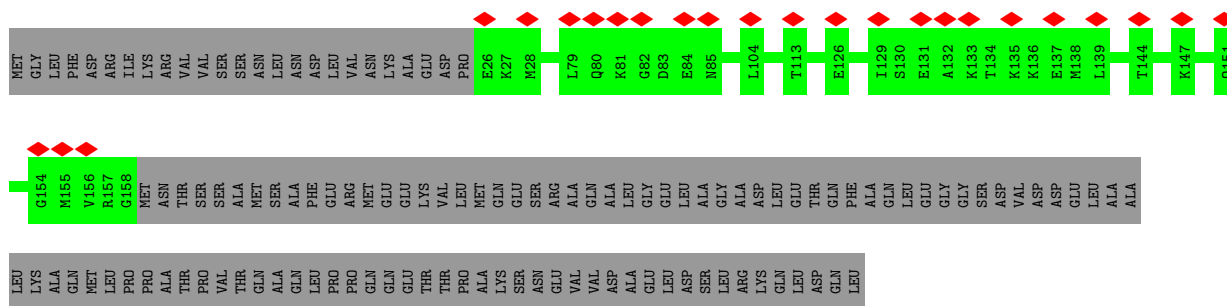
- Molecule 1: vipp1



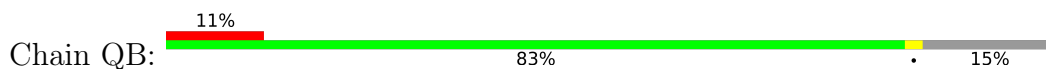
- Molecule 1: vipp1



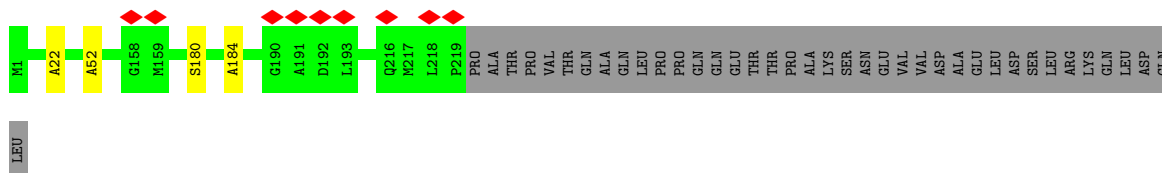
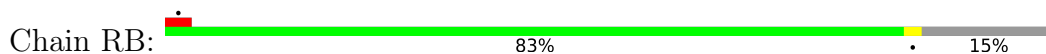
- Molecule 1: vipp1



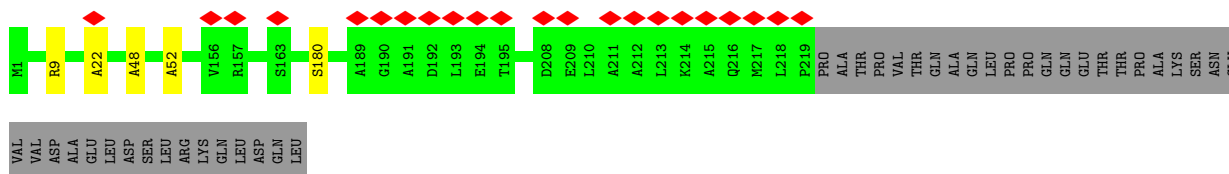
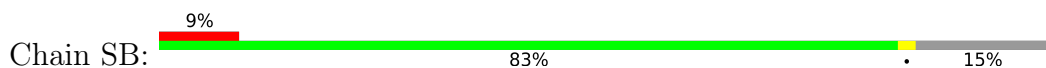
- Molecule 1: vipp1



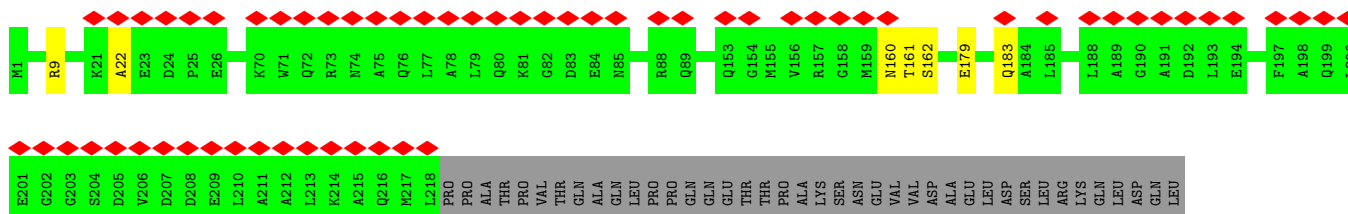
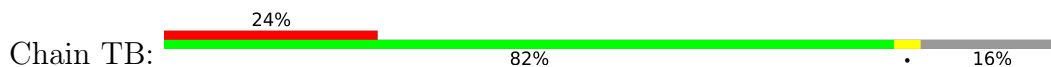
- Molecule 1: vippp1



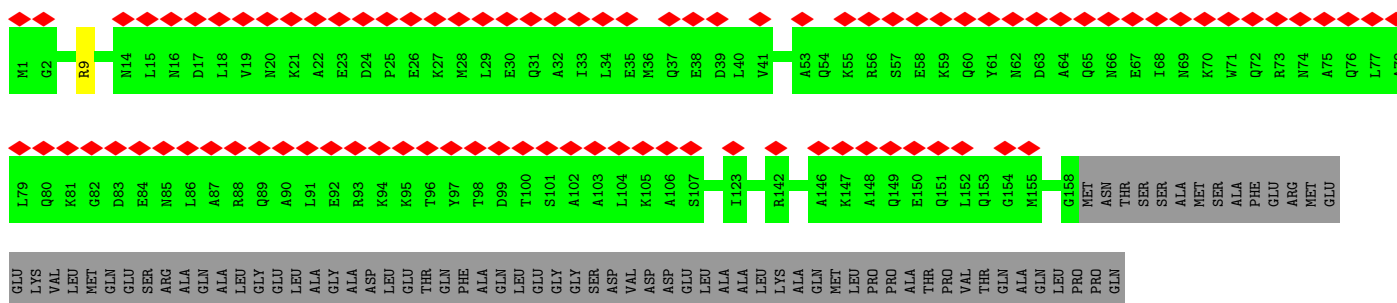
- Molecule 1: vippp1



- Molecule 1: vippp1



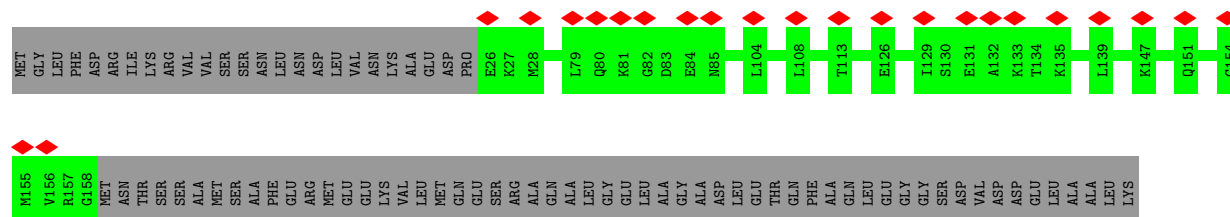
- Molecule 1: vippp1



GLN  
GLU  
THR  
THR  
PRO  
ALA  
LYS  
SER  
ASN  
GLU  
VAL  
VAL  
SER  
ASP  
ALA  
GLU  
LEU  
ASP  
SER  
LEU  
ARG  
LYS  
GLN  
LEU  
LEU  
ASP  
GLN  
LEU

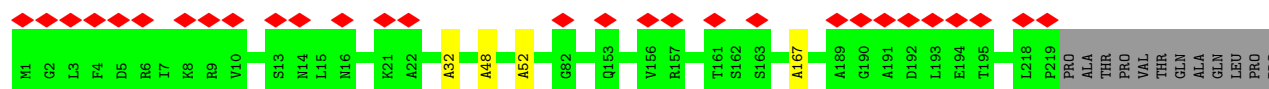
• Molecule 1: vipp1

Chain VB: 9% 52% 48%



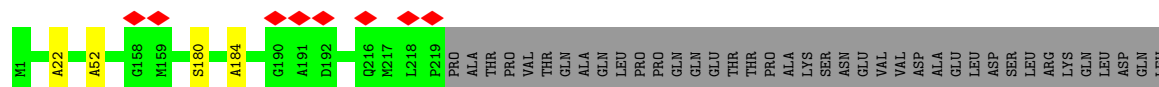
• Molecule 1: vipp1

Chain WB: 11% 83% 15%



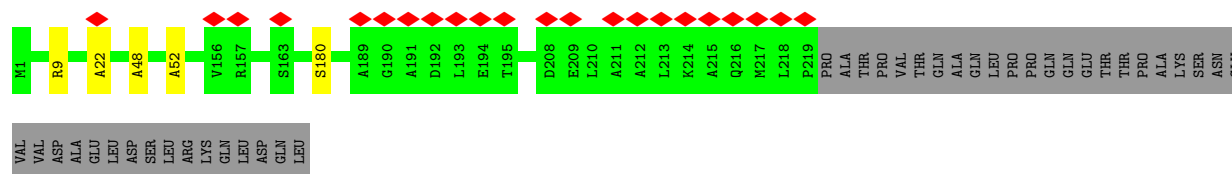
• Molecule 1: vipp1

Chain XB: 83% 15%



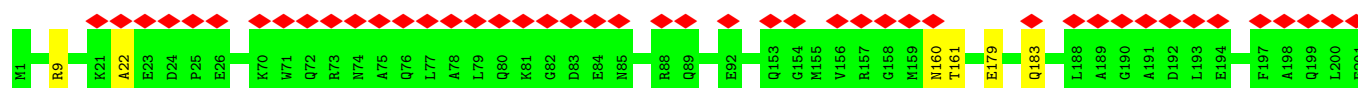
• Molecule 1: vipp1

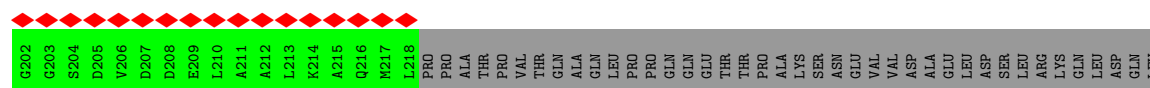
Chain YB: 9% 83% 15%



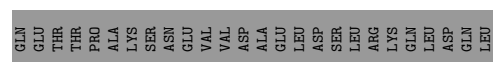
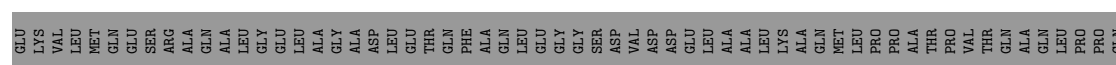
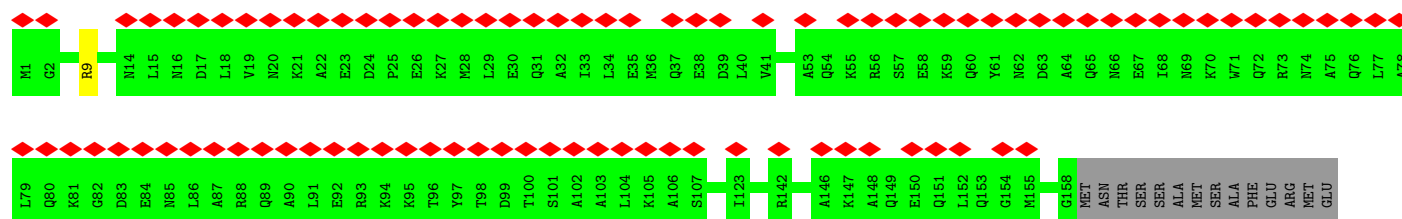
• Molecule 1: vipp1

Chain ZB: 24% 82% 16%

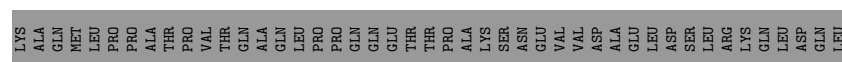
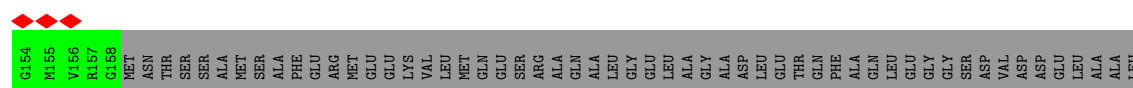
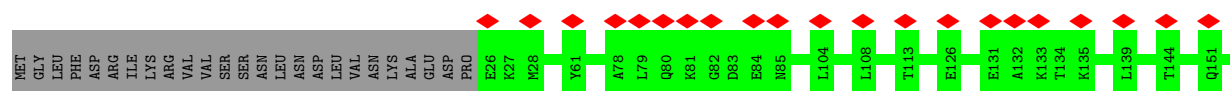




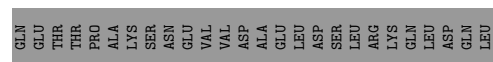
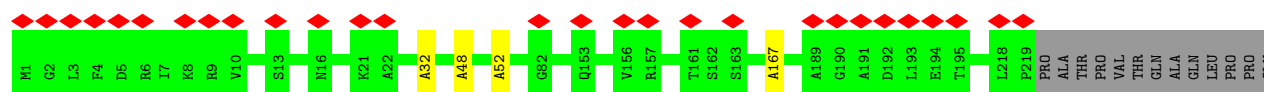
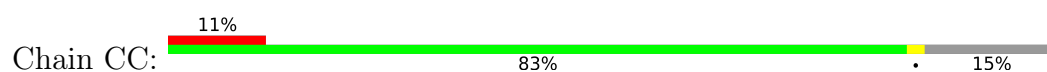
• Molecule 1: vipp1



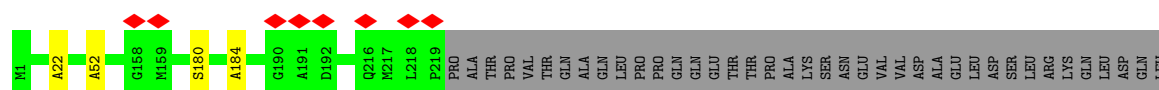
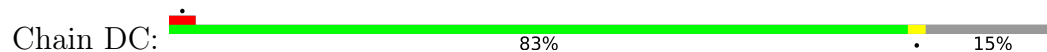
• Molecule 1: vipp1



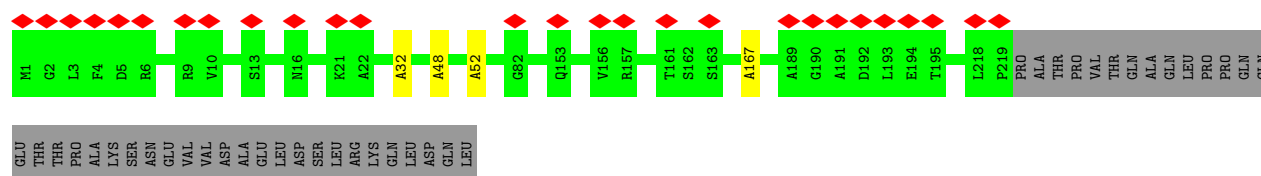
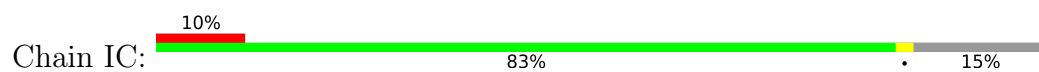
• Molecule 1: vipp1



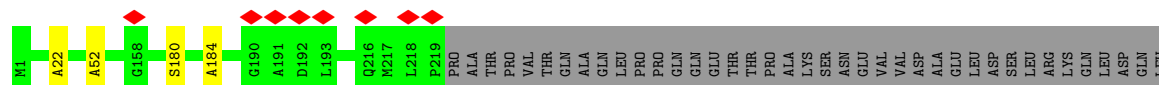
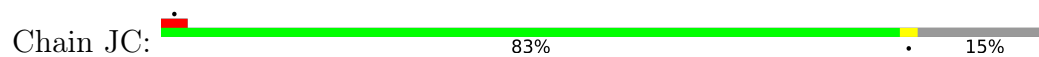
• Molecule 1: vipp1



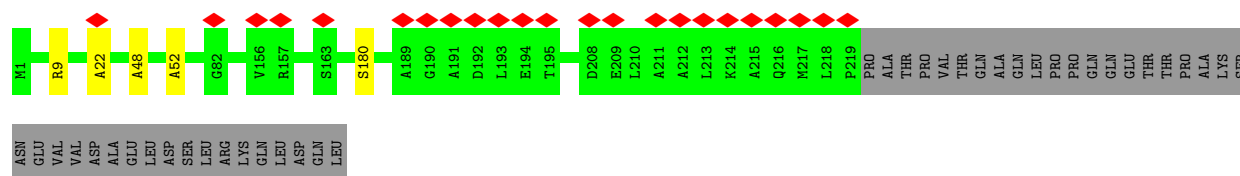
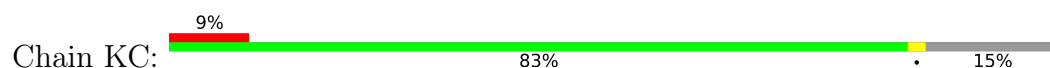




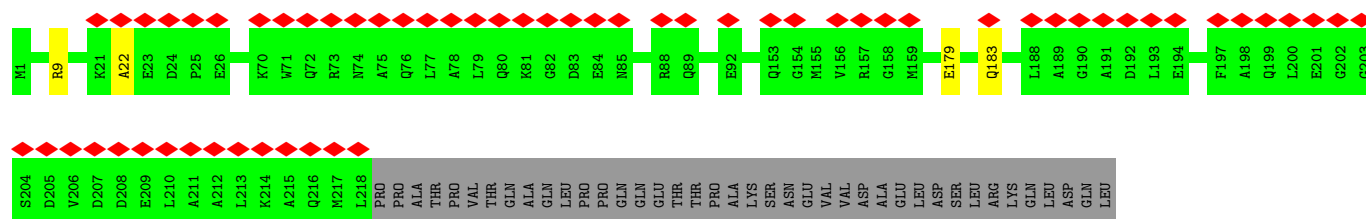
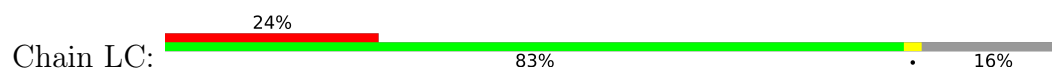
• Molecule 1: vipp1



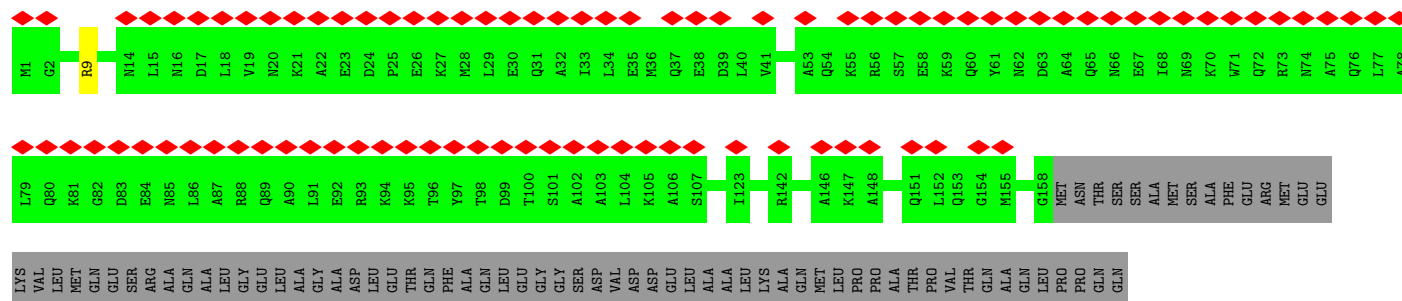
• Molecule 1: vipp1



• Molecule 1: vipp1



• Molecule 1: vipp1





GLU	THR	THR	PRO	ALA	LYS	SER	ASN	GLU	VAL	VAL	ASP	ALA	GLU	LEU	ASP	SER	LEU	ARG	LYS	GLN	LEU	ASP	GLN	LEU
-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----

## 4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	18217	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ( $e^-/\text{\AA}^2$ )	1.5	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.046	Depositor
Minimum map value	-0.006	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.003	Depositor
Recommended contour level	0.0182	Depositor
Map size ( $\text{\AA}$ )	467.04, 467.04, 467.04	wwPDB
Map dimensions	336, 336, 336	wwPDB
Map angles ( $^\circ$ )	90.0, 90.0, 90.0	wwPDB
Pixel spacing ( $\text{\AA}$ )	1.39, 1.39, 1.39	Depositor

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z  > 5$	RMSZ	# $ Z  > 5$
1	A	0.18	0/660	0.37	0/920
1	AA	0.18	0/1085	0.39	0/1512
1	AB	0.19	0/1085	0.39	0/1512
1	AC	0.18	0/784	0.37	0/1093
1	B	0.18	0/1085	0.39	0/1512
1	BA	0.19	0/1085	0.44	0/1512
1	BB	0.21	0/1080	0.42	0/1505
1	BC	0.18	0/660	0.37	0/920
1	C	0.19	0/1085	0.44	0/1512
1	CA	0.19	0/1085	0.39	0/1512
1	CB	0.18	0/784	0.37	0/1093
1	CC	0.18	0/1085	0.39	0/1512
1	D	0.19	0/1085	0.39	0/1512
1	DA	0.21	0/1080	0.42	0/1505
1	DB	0.18	0/660	0.37	0/920
1	DC	0.19	0/1085	0.44	0/1512
1	E	0.21	0/1080	0.42	0/1505
1	EA	0.18	0/784	0.37	0/1093
1	EB	0.18	0/1085	0.39	0/1512
1	EC	0.19	0/1085	0.39	0/1512
1	F	0.18	0/784	0.37	0/1093
1	FA	0.18	0/660	0.37	0/920
1	FB	0.19	0/1085	0.44	0/1512
1	FC	0.21	0/1080	0.42	0/1505
1	G	0.18	0/660	0.37	0/920
1	GA	0.19	0/1085	0.39	0/1512
1	GB	0.19	0/1085	0.39	0/1512
1	GC	0.18	0/784	0.37	0/1093
1	H	0.19	0/1085	0.39	0/1512
1	HA	0.19	0/1085	0.44	0/1512
1	HB	0.21	0/1080	0.42	0/1505
1	HC	0.18	0/660	0.37	0/920
1	I	0.19	0/1085	0.44	0/1512
1	IA	0.19	0/1085	0.39	0/1512

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
1	IB	0.18	0/784	0.37	0/1093
1	IC	0.18	0/1085	0.39	0/1512
1	J	0.19	0/1085	0.39	0/1512
1	JA	0.21	0/1080	0.42	0/1505
1	JB	0.18	0/660	0.37	0/920
1	JC	0.19	0/1085	0.44	0/1512
1	K	0.21	0/1080	0.42	0/1505
1	KA	0.18	0/784	0.37	0/1093
1	KB	0.18	0/1085	0.39	0/1512
1	KC	0.19	0/1085	0.39	0/1512
1	L	0.18	0/784	0.37	0/1093
1	LA	0.18	0/660	0.37	0/920
1	LB	0.19	0/1085	0.44	0/1512
1	LC	0.20	0/1080	0.42	0/1505
1	M	0.18	0/660	0.37	0/920
1	MA	0.18	0/1085	0.39	0/1512
1	MB	0.19	0/1085	0.39	0/1512
1	MC	0.18	0/784	0.37	0/1093
1	N	0.18	0/1085	0.39	0/1512
1	NA	0.19	0/1085	0.44	0/1512
1	NB	0.21	0/1080	0.42	0/1505
1	O	0.19	0/1085	0.44	0/1512
1	OA	0.19	0/1085	0.39	0/1512
1	OB	0.18	0/784	0.37	0/1093
1	P	0.19	0/1085	0.39	0/1512
1	PA	0.21	0/1080	0.42	0/1505
1	PB	0.18	0/660	0.37	0/920
1	Q	0.21	0/1080	0.42	0/1505
1	QA	0.18	0/784	0.37	0/1093
1	QB	0.19	0/1085	0.39	0/1512
1	R	0.18	0/784	0.37	0/1093
1	RA	0.18	0/660	0.37	0/920
1	RB	0.19	0/1085	0.44	0/1512
1	S	0.18	0/660	0.37	0/920
1	SA	0.18	0/1085	0.39	0/1512
1	SB	0.19	0/1085	0.39	0/1512
1	T	0.19	0/1085	0.39	0/1512
1	TA	0.19	0/1085	0.44	0/1512
1	TB	0.21	0/1080	0.42	0/1505
1	UA	0.19	0/1085	0.39	0/1512
1	UB	0.18	0/784	0.37	0/1093
1	V	0.19	0/1085	0.44	0/1512
1	VA	0.20	0/1080	0.42	0/1505

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
1	VB	0.18	0/660	0.37	0/920
1	W	0.19	0/1085	0.39	0/1512
1	WA	0.18	0/784	0.37	0/1093
1	WB	0.18	0/1085	0.39	0/1512
1	X	0.21	0/1080	0.42	0/1505
1	XA	0.18	0/660	0.37	0/920
1	XB	0.19	0/1085	0.44	0/1512
1	Y	0.18	0/784	0.37	0/1093
1	YA	0.18	0/1085	0.39	0/1512
1	YB	0.19	0/1085	0.39	0/1512
1	Z	0.18	0/660	0.37	0/920
1	ZA	0.19	0/1085	0.44	0/1512
1	ZB	0.21	0/1080	0.42	0/1505
All	All	0.19	0/86685	0.40	0/120810

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

## 5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	661	0	320	0	0
1	AA	1086	0	531	13	0
1	AB	1086	0	531	9	0
1	AC	785	0	376	2	0
1	B	1086	0	531	13	0
1	BA	1086	0	531	9	0
1	BB	1081	0	530	9	0
1	BC	661	0	320	0	0
1	C	1086	0	531	9	0
1	CA	1086	0	531	9	0
1	CB	785	0	376	2	0
1	CC	1086	0	531	13	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	1086	0	531	9	0
1	DA	1081	0	530	7	0
1	DB	661	0	320	0	0
1	DC	1086	0	531	9	0
1	E	1081	0	530	7	0
1	EA	785	0	376	2	0
1	EB	1086	0	531	13	0
1	EC	1086	0	531	9	0
1	F	785	0	376	2	0
1	FA	661	0	320	0	0
1	FB	1086	0	531	9	0
1	FC	1081	0	530	8	0
1	G	661	0	320	0	0
1	GA	1086	0	531	13	0
1	GB	1086	0	531	9	0
1	GC	785	0	376	2	0
1	H	1086	0	531	13	0
1	HA	1086	0	531	9	0
1	HB	1081	0	530	9	0
1	HC	661	0	320	0	0
1	I	1086	0	531	9	0
1	IA	1086	0	531	9	0
1	IB	785	0	376	2	0
1	IC	1086	0	531	13	0
1	J	1086	0	531	9	0
1	JA	1081	0	530	7	0
1	JB	661	0	320	0	0
1	JC	1086	0	531	9	0
1	K	1081	0	530	7	0
1	KA	785	0	376	2	0
1	KB	1086	0	531	13	0
1	KC	1086	0	531	9	0
1	L	785	0	376	2	0
1	LA	661	0	320	0	0
1	LB	1086	0	531	9	0
1	LC	1081	0	530	7	0
1	M	661	0	320	0	0
1	MA	1086	0	531	13	0
1	MB	1086	0	531	9	0
1	MC	785	0	376	2	0
1	N	1086	0	531	12	0
1	NA	1086	0	531	9	0

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	NB	1081	0	530	9	0
1	O	1086	0	531	9	0
1	OA	1086	0	531	9	0
1	OB	785	0	376	2	0
1	P	1086	0	531	9	0
1	PA	1081	0	530	9	0
1	PB	661	0	320	0	0
1	Q	1081	0	530	7	0
1	QA	785	0	376	2	0
1	QB	1086	0	531	13	0
1	R	785	0	376	2	0
1	RA	661	0	320	0	0
1	RB	1086	0	531	9	0
1	S	661	0	320	0	0
1	SA	1086	0	531	13	0
1	SB	1086	0	531	9	0
1	T	1086	0	531	12	0
1	TA	1086	0	531	9	0
1	TB	1081	0	530	9	0
1	UA	1086	0	531	9	0
1	UB	785	0	376	2	0
1	V	1086	0	531	9	0
1	VA	1081	0	530	9	0
1	VB	661	0	320	0	0
1	W	1086	0	531	9	0
1	WA	785	0	376	2	0
1	WB	1086	0	531	13	0
1	X	1081	0	530	7	0
1	XA	661	0	320	0	0
1	XB	1086	0	531	9	0
1	Y	785	0	376	2	0
1	YA	1086	0	531	13	0
1	YB	1086	0	531	9	0
1	Z	661	0	320	0	0
1	ZA	1086	0	531	9	0
1	ZB	1081	0	530	8	0
All	All	86775	0	42285	313	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 2.

The worst 5 of 313 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:ZA:184:ALA:HB1	1:KB:48:ALA:HB1	1.21	1.18
1:TA:184:ALA:HB1	1:EB:48:ALA:HB1	1.21	1.18
1:NA:184:ALA:HB1	1:YA:48:ALA:HB1	1.23	1.17
1:HA:184:ALA:HB1	1:SA:48:ALA:HB1	1.22	1.16
1:LB:184:ALA:HB1	1:WB:48:ALA:HB1	1.22	1.15

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	131/258 (51%)	131 (100%)	0	0	100	100
1	AA	217/258 (84%)	216 (100%)	1 (0%)	0	100	100
1	AB	217/258 (84%)	216 (100%)	1 (0%)	0	100	100
1	AC	156/258 (60%)	156 (100%)	0	0	100	100
1	B	217/258 (84%)	216 (100%)	1 (0%)	0	100	100
1	BA	217/258 (84%)	217 (100%)	0	0	100	100
1	BB	216/258 (84%)	215 (100%)	1 (0%)	0	100	100
1	BC	131/258 (51%)	131 (100%)	0	0	100	100
1	C	217/258 (84%)	217 (100%)	0	0	100	100
1	CA	217/258 (84%)	216 (100%)	1 (0%)	0	100	100
1	CB	156/258 (60%)	156 (100%)	0	0	100	100
1	CC	217/258 (84%)	216 (100%)	1 (0%)	0	100	100
1	D	217/258 (84%)	216 (100%)	1 (0%)	0	100	100
1	DA	216/258 (84%)	215 (100%)	1 (0%)	0	100	100
1	DB	131/258 (51%)	131 (100%)	0	0	100	100
1	DC	217/258 (84%)	217 (100%)	0	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	E	216/258 (84%)	215 (100%)	1 (0%)	0	100	100
1	EA	156/258 (60%)	156 (100%)	0	0	100	100
1	EB	217/258 (84%)	216 (100%)	1 (0%)	0	100	100
1	EC	217/258 (84%)	216 (100%)	1 (0%)	0	100	100
1	F	156/258 (60%)	156 (100%)	0	0	100	100
1	FA	131/258 (51%)	131 (100%)	0	0	100	100
1	FB	217/258 (84%)	217 (100%)	0	0	100	100
1	FC	216/258 (84%)	215 (100%)	1 (0%)	0	100	100
1	G	131/258 (51%)	131 (100%)	0	0	100	100
1	GA	217/258 (84%)	216 (100%)	1 (0%)	0	100	100
1	GB	217/258 (84%)	216 (100%)	1 (0%)	0	100	100
1	GC	156/258 (60%)	156 (100%)	0	0	100	100
1	H	217/258 (84%)	216 (100%)	1 (0%)	0	100	100
1	HA	217/258 (84%)	217 (100%)	0	0	100	100
1	HB	216/258 (84%)	215 (100%)	1 (0%)	0	100	100
1	HC	131/258 (51%)	131 (100%)	0	0	100	100
1	I	217/258 (84%)	217 (100%)	0	0	100	100
1	IA	217/258 (84%)	216 (100%)	1 (0%)	0	100	100
1	IB	156/258 (60%)	156 (100%)	0	0	100	100
1	IC	217/258 (84%)	216 (100%)	1 (0%)	0	100	100
1	J	217/258 (84%)	216 (100%)	1 (0%)	0	100	100
1	JA	216/258 (84%)	215 (100%)	1 (0%)	0	100	100
1	JB	131/258 (51%)	131 (100%)	0	0	100	100
1	JC	217/258 (84%)	217 (100%)	0	0	100	100
1	K	216/258 (84%)	215 (100%)	1 (0%)	0	100	100
1	KA	156/258 (60%)	156 (100%)	0	0	100	100
1	KB	217/258 (84%)	216 (100%)	1 (0%)	0	100	100
1	KC	217/258 (84%)	216 (100%)	1 (0%)	0	100	100
1	L	156/258 (60%)	156 (100%)	0	0	100	100
1	LA	131/258 (51%)	131 (100%)	0	0	100	100
1	LB	217/258 (84%)	217 (100%)	0	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	LC	216/258 (84%)	215 (100%)	1 (0%)	0	100	100
1	M	131/258 (51%)	131 (100%)	0	0	100	100
1	MA	217/258 (84%)	216 (100%)	1 (0%)	0	100	100
1	MB	217/258 (84%)	216 (100%)	1 (0%)	0	100	100
1	MC	156/258 (60%)	156 (100%)	0	0	100	100
1	N	217/258 (84%)	216 (100%)	1 (0%)	0	100	100
1	NA	217/258 (84%)	217 (100%)	0	0	100	100
1	NB	216/258 (84%)	215 (100%)	1 (0%)	0	100	100
1	O	217/258 (84%)	217 (100%)	0	0	100	100
1	OA	217/258 (84%)	216 (100%)	1 (0%)	0	100	100
1	OB	156/258 (60%)	156 (100%)	0	0	100	100
1	P	217/258 (84%)	216 (100%)	1 (0%)	0	100	100
1	PA	216/258 (84%)	215 (100%)	1 (0%)	0	100	100
1	PB	131/258 (51%)	131 (100%)	0	0	100	100
1	Q	216/258 (84%)	215 (100%)	1 (0%)	0	100	100
1	QA	156/258 (60%)	156 (100%)	0	0	100	100
1	QB	217/258 (84%)	216 (100%)	1 (0%)	0	100	100
1	R	156/258 (60%)	156 (100%)	0	0	100	100
1	RA	131/258 (51%)	131 (100%)	0	0	100	100
1	RB	217/258 (84%)	217 (100%)	0	0	100	100
1	S	131/258 (51%)	131 (100%)	0	0	100	100
1	SA	217/258 (84%)	216 (100%)	1 (0%)	0	100	100
1	SB	217/258 (84%)	216 (100%)	1 (0%)	0	100	100
1	T	217/258 (84%)	216 (100%)	1 (0%)	0	100	100
1	TA	217/258 (84%)	217 (100%)	0	0	100	100
1	TB	216/258 (84%)	215 (100%)	1 (0%)	0	100	100
1	UA	217/258 (84%)	216 (100%)	1 (0%)	0	100	100
1	UB	156/258 (60%)	156 (100%)	0	0	100	100
1	V	217/258 (84%)	217 (100%)	0	0	100	100
1	VA	216/258 (84%)	215 (100%)	1 (0%)	0	100	100
1	VB	131/258 (51%)	131 (100%)	0	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	W	217/258 (84%)	216 (100%)	1 (0%)	0	100	100
1	WA	156/258 (60%)	156 (100%)	0	0	100	100
1	WB	217/258 (84%)	216 (100%)	1 (0%)	0	100	100
1	X	216/258 (84%)	215 (100%)	1 (0%)	0	100	100
1	XA	131/258 (51%)	131 (100%)	0	0	100	100
1	XB	217/258 (84%)	217 (100%)	0	0	100	100
1	Y	156/258 (60%)	156 (100%)	0	0	100	100
1	YA	217/258 (84%)	216 (100%)	1 (0%)	0	100	100
1	YB	217/258 (84%)	216 (100%)	1 (0%)	0	100	100
1	Z	131/258 (51%)	131 (100%)	0	0	100	100
1	ZA	217/258 (84%)	217 (100%)	0	0	100	100
1	ZB	216/258 (84%)	215 (100%)	1 (0%)	0	100	100
All	All	17310/23220 (74%)	17265 (100%)	45 (0%)	0	100	100

There are no Ramachandran outliers to report.

### 5.3.2 Protein sidechains [i](#)

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

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

There are no ligands in this entry.

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

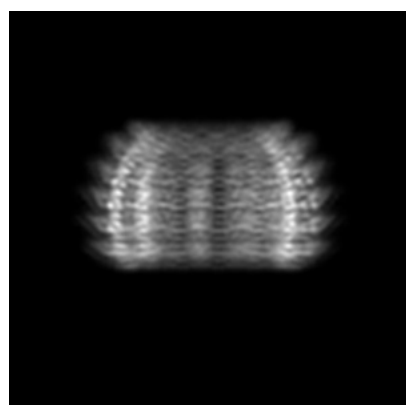
## 6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-11481. These allow visual inspection of the internal detail of the map and identification of artifacts.

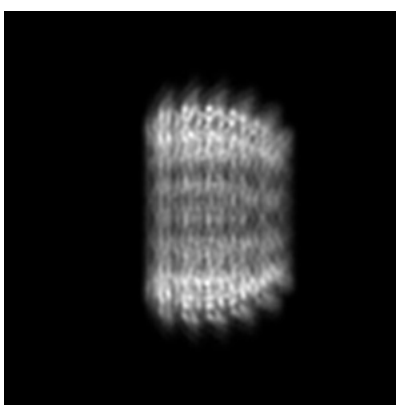
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

### 6.1 Orthogonal projections [i](#)

#### 6.1.1 Primary map



X



Y

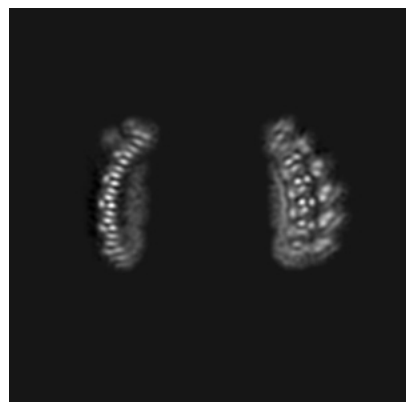


Z

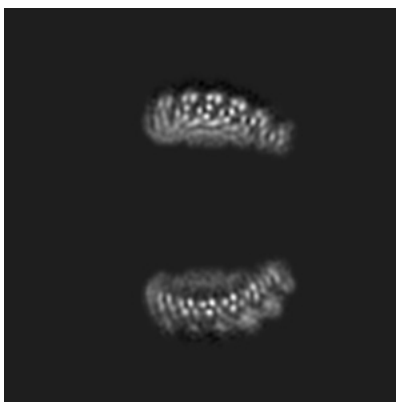
The images above show the map projected in three orthogonal directions.

### 6.2 Central slices [i](#)

#### 6.2.1 Primary map



X Index: 168



Y Index: 168



Z Index: 168

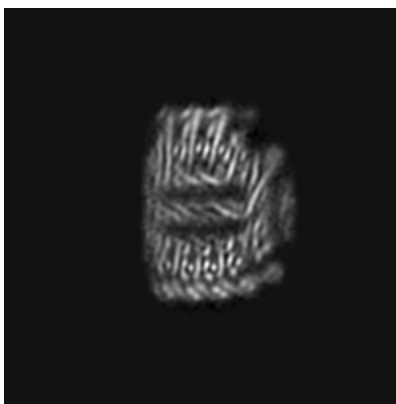
The images above show central slices of the map in three orthogonal directions.

## 6.3 Largest variance slices [i](#)

### 6.3.1 Primary map



X Index: 108



Y Index: 235

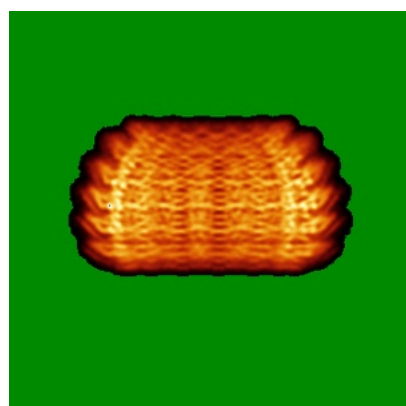


Z Index: 172

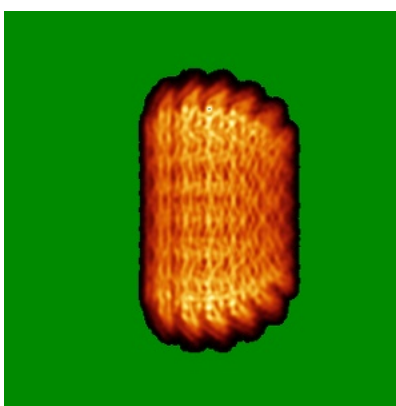
The images above show the largest variance slices of the map in three orthogonal directions.

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

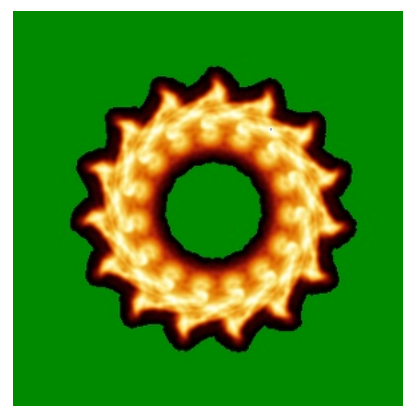
### 6.4.1 Primary map



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



The images above show the 3D surface view of the map at the recommended contour level 0.0182. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

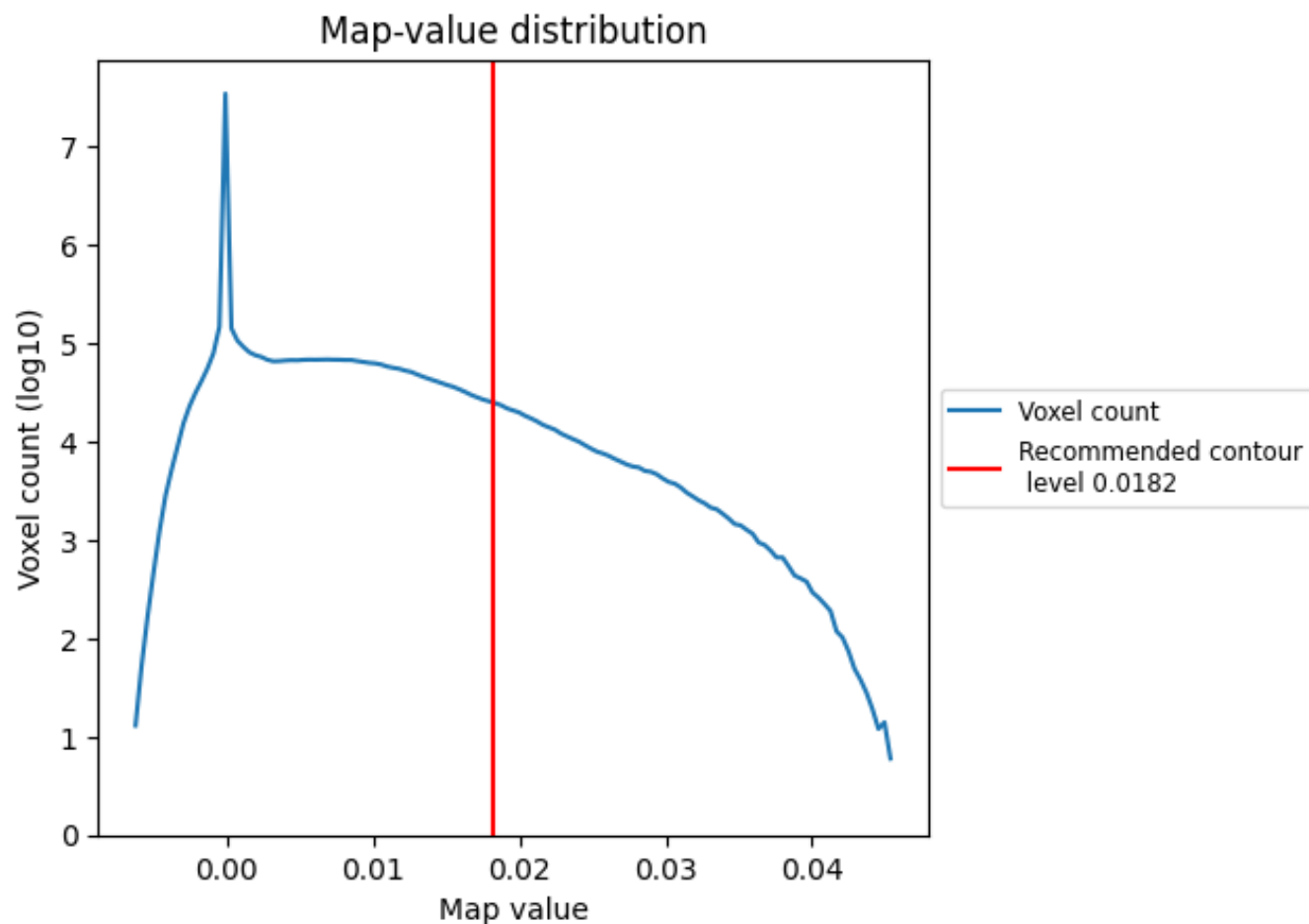
## 6.6 Mask visualisation [i](#)

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

## 7 Map analysis [i](#)

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

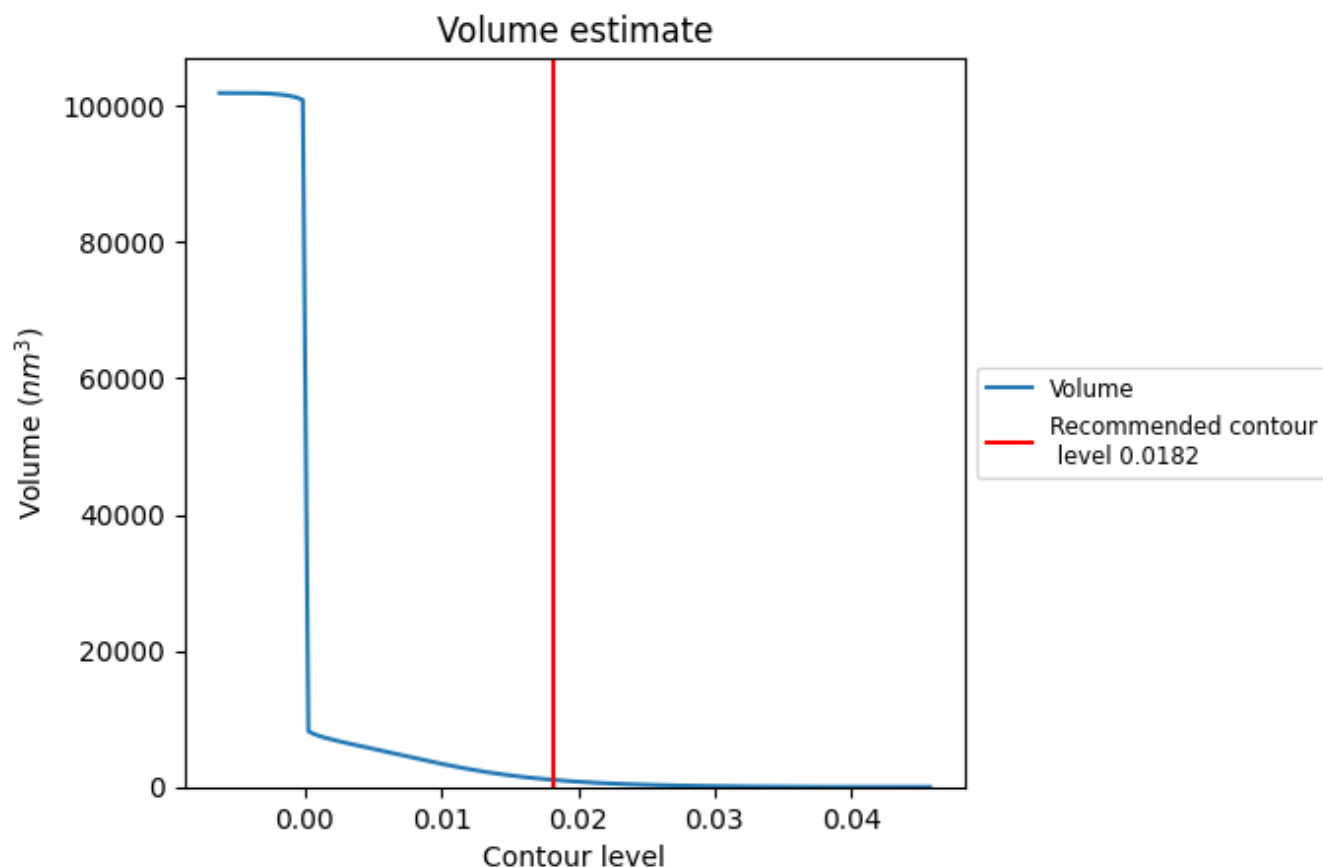
### 7.1 Map-value distribution [i](#)



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.



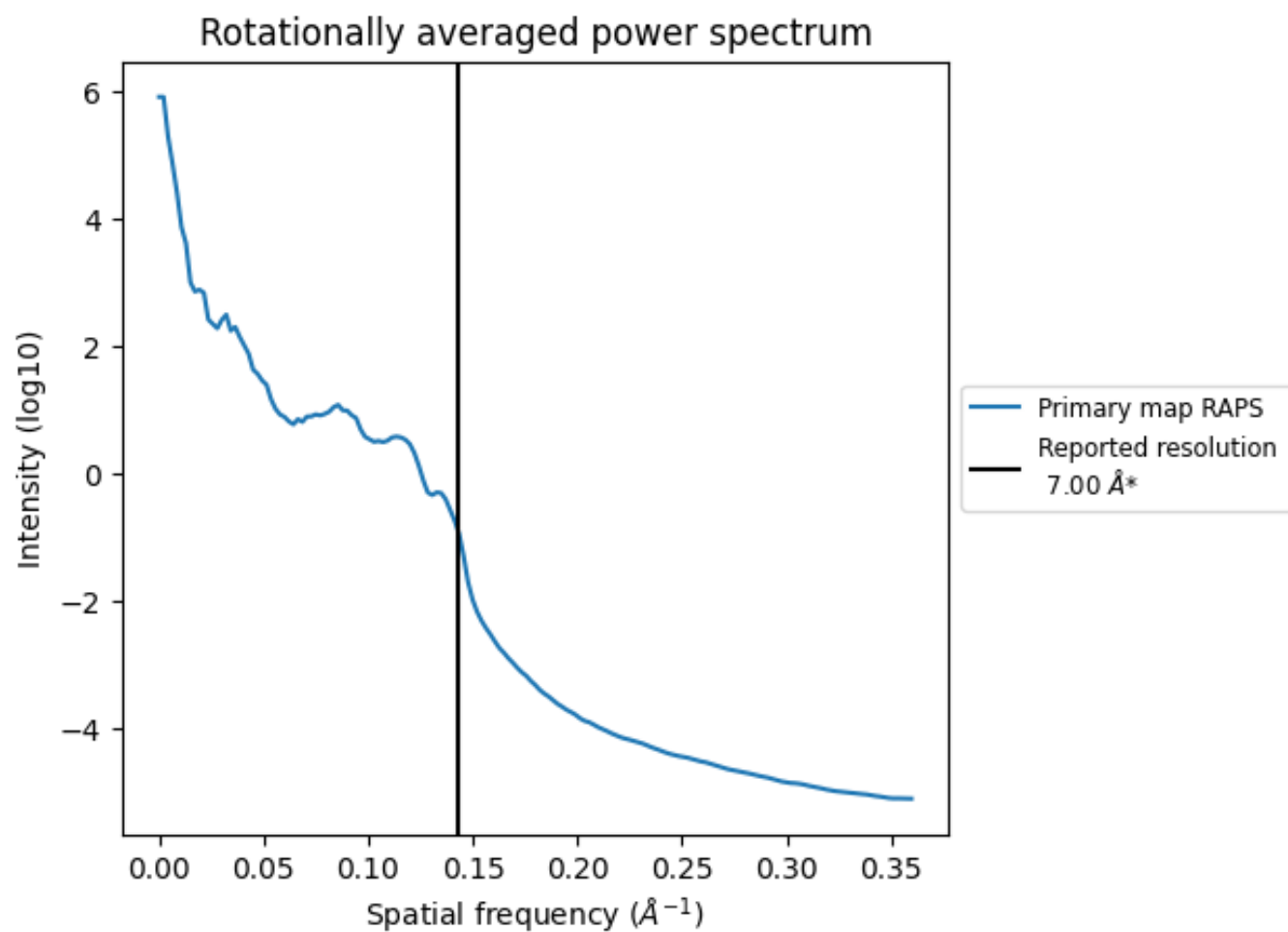
## 7.2 Volume estimate [i](#)



The volume at the recommended contour level is 1040  $\text{nm}^3$ ; this corresponds to an approximate mass of 939 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 0.143 Å<sup>-1</sup>

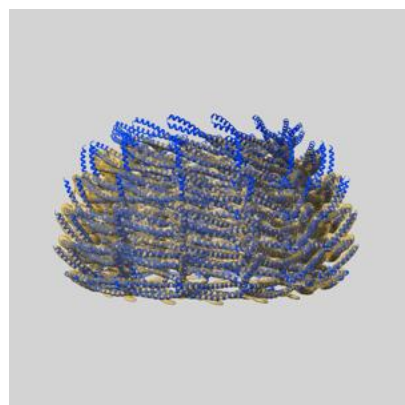
## 8 Fourier-Shell correlation

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

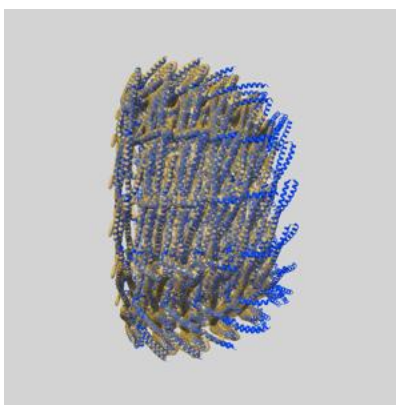
## 9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-11481 and PDB model 6ZW5. Per-residue inclusion information can be found in section 3 on page 12.

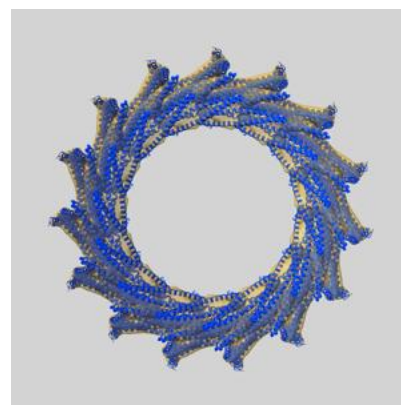
### 9.1 Map-model overlay [i](#)



X



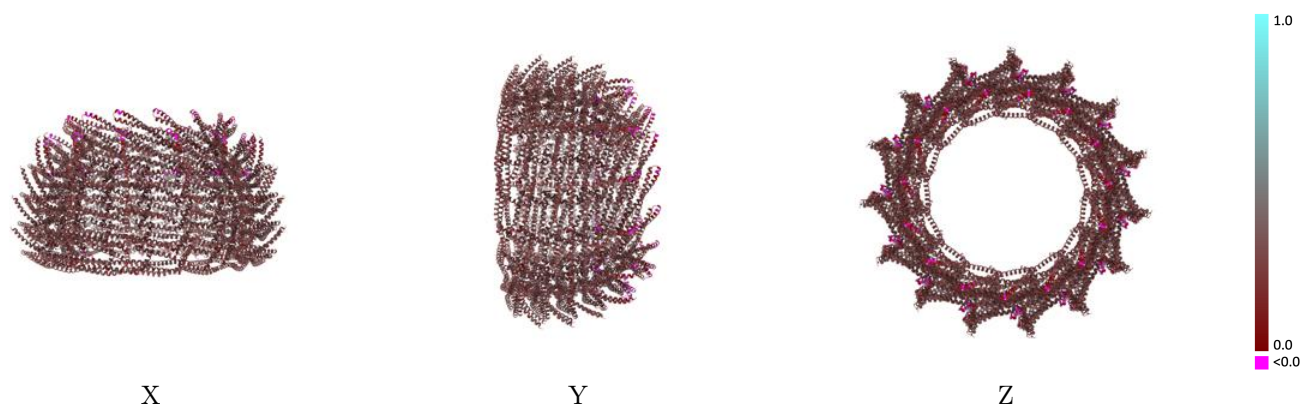
Y



Z

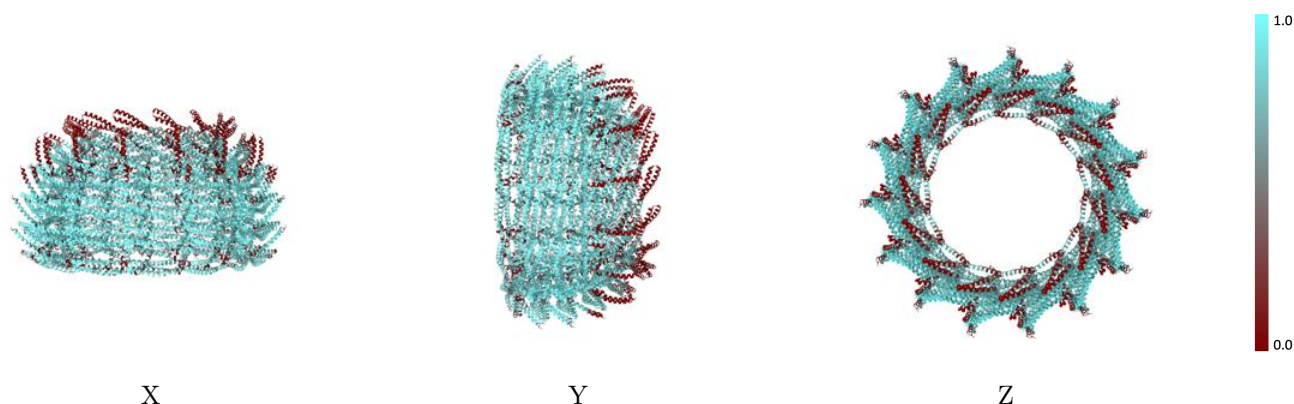
The images above show the 3D surface view of the map at the recommended contour level 0.0182 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



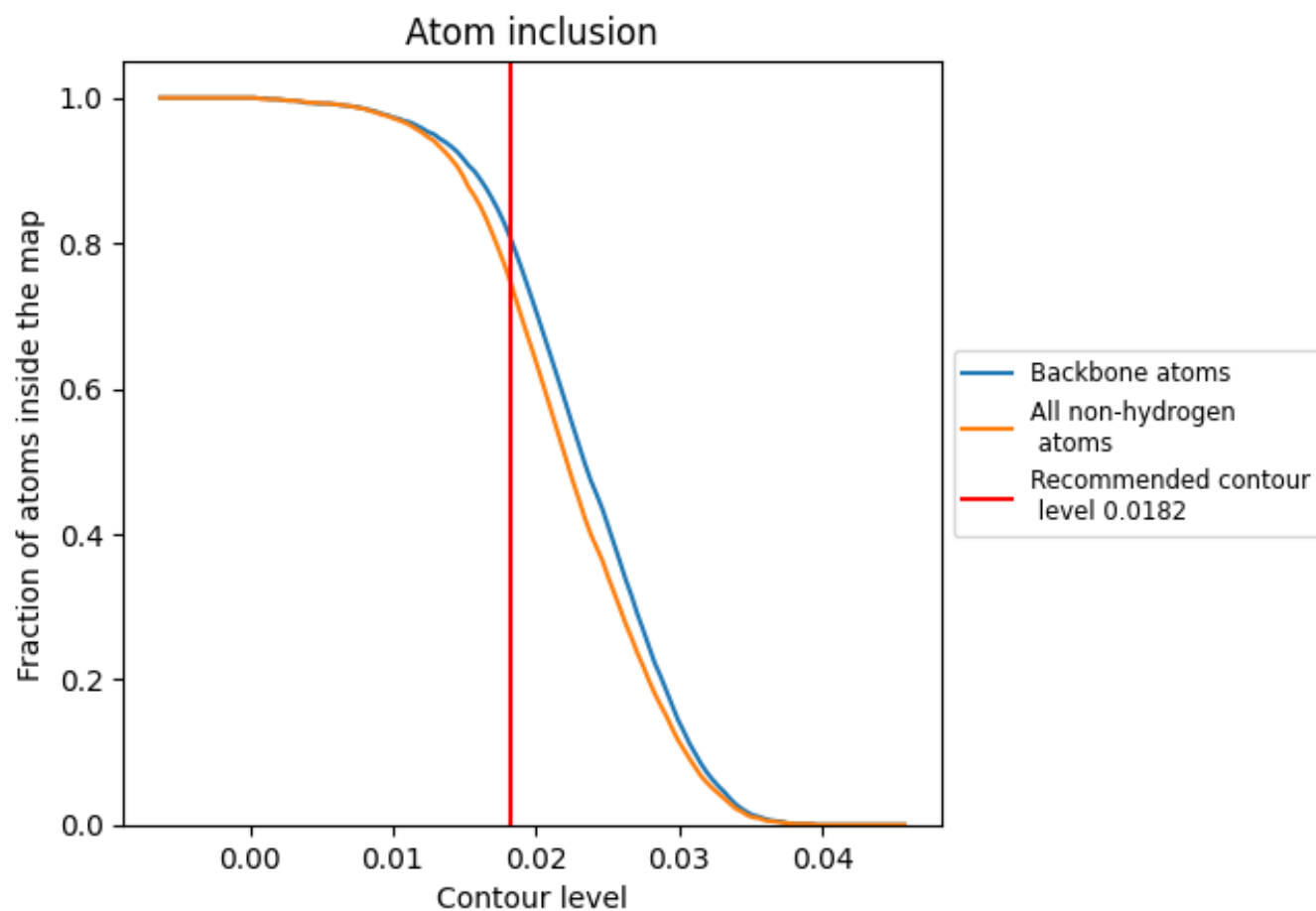
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.0182).







































































## 9.4 Atom inclusion [i](#)



At the recommended contour level, 81% of all backbone atoms, 75% 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 (0.0182) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.7480	 0.2580
A	 0.7730	 0.2660
AA	 0.8060	 0.2610
AB	 0.8590	 0.2680
AC	 0.3730	 0.2370
B	 0.8240	 0.2610
BA	 0.9160	 0.2720
BB	 0.6770	 0.2410
BC	 0.7440	 0.2690
C	 0.9210	 0.2730
CA	 0.8540	 0.2690
CB	 0.3950	 0.2390
CC	 0.8200	 0.2590
D	 0.8560	 0.2700
DA	 0.6650	 0.2410
DB	 0.7550	 0.2680
DC	 0.9160	 0.2730
E	 0.6660	 0.2420
EA	 0.3950	 0.2350
EB	 0.8200	 0.2620
EC	 0.8510	 0.2690
F	 0.3910	 0.2350
FA	 0.7170	 0.2690
FB	 0.9200	 0.2740
FC	 0.6540	 0.2410
G	 0.7620	 0.2660
GA	 0.8080	 0.2590
GB	 0.8530	 0.2700
GC	 0.3860	 0.2350
H	 0.8190	 0.2590
HA	 0.9180	 0.2710
HB	 0.6690	 0.2440
HC	 0.7650	 0.2670
I	 0.9210	 0.2730
IA	 0.8580	 0.2690



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





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Chain	Atom inclusion	Q-score
IB	 0.3780	 0.2380
IC	 0.8250	 0.2590
J	 0.8590	 0.2700
JA	 0.6700	 0.2400
JB	 0.7500	 0.2700
JC	 0.9180	 0.2720
K	 0.6680	 0.2420
KA	 0.3910	 0.2360
KB	 0.8160	 0.2600
KC	 0.8470	 0.2690
L	 0.3960	 0.2360
LA	 0.7440	 0.2700
LB	 0.9170	 0.2740
LC	 0.6580	 0.2400
M	 0.7370	 0.2660
MA	 0.8110	 0.2620
MB	 0.8470	 0.2690
MC	 0.3860	 0.2360
N	 0.8080	 0.2580
NA	 0.9170	 0.2740
NB	 0.6530	 0.2430
O	 0.9180	 0.2710
OA	 0.8540	 0.2690
OB	 0.3720	 0.2380
P	 0.8540	 0.2670
PA	 0.6680	 0.2410
PB	 0.7440	 0.2680
Q	 0.6670	 0.2390
QA	 0.3960	 0.2360
QB	 0.8080	 0.2600
R	 0.3920	 0.2330
RA	 0.7580	 0.2700
RB	 0.9150	 0.2720
S	 0.7310	 0.2690
SA	 0.8220	 0.2610
SB	 0.8460	 0.2680
T	 0.8040	 0.2600
TA	 0.9210	 0.2720
TB	 0.6540	 0.2410
UA	 0.8520	 0.2690
UB	 0.3640	 0.2380
V	 0.9150	 0.2730

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Chain	Atom inclusion	Q-score
VA	 0.6720	 0.2430
VB	 0.7440	 0.2650
W	 0.8470	 0.2690
WA	 0.3990	 0.2390
WB	 0.8140	 0.2570
X	 0.6640	 0.2410
XA	 0.7640	 0.2680
XB	 0.9180	 0.2710
Y	 0.3980	 0.2350
YA	 0.8230	 0.2610
YB	 0.8510	 0.2670
Z	 0.7290	 0.2660
ZA	 0.9240	 0.2720
ZB	 0.6540	 0.2410