



## Full wwPDB EM Validation Report ⓘ

Dec 18, 2022 – 03:33 am GMT

PDB ID : 6ZVT  
EMDB ID : EMD-11470  
Title : C13 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 Full wwPDB EM Validation 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:

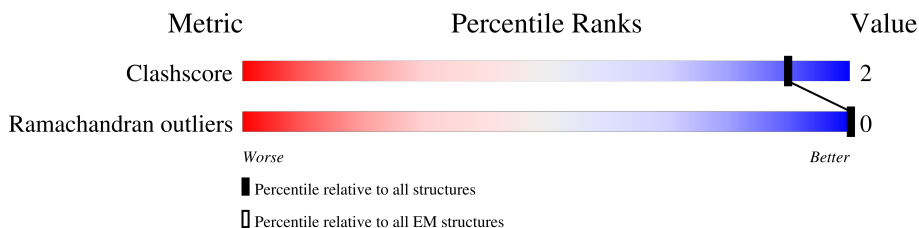
EMDB validation analysis : 0.0.1.dev43  
MolProbity : 4.02b-467  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
MapQ : 1.9.9  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.31.3

# 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 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.



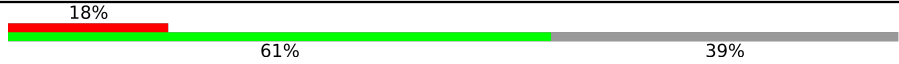
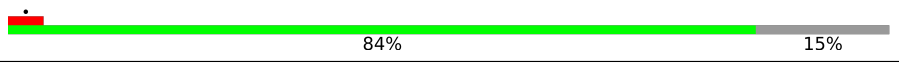
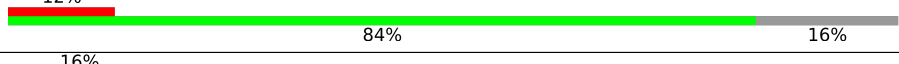
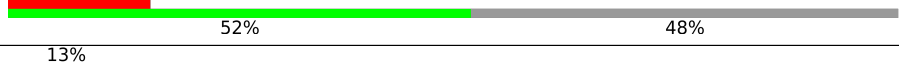

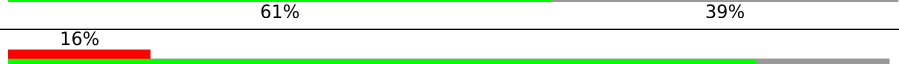
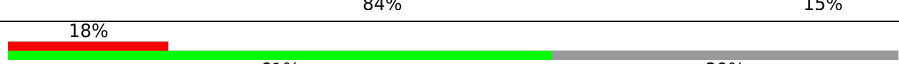
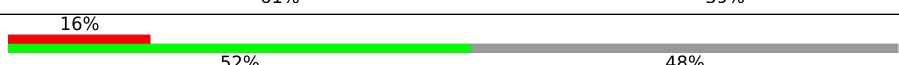
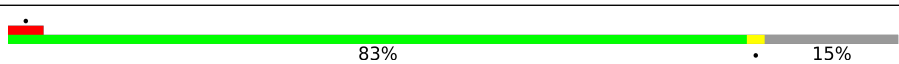


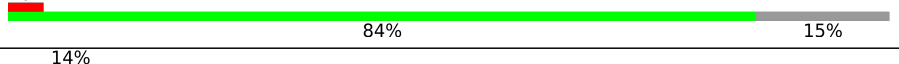
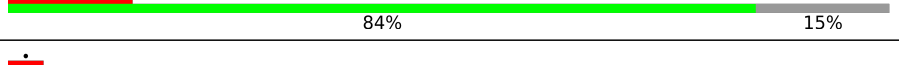

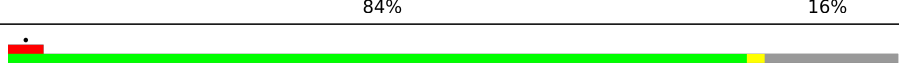




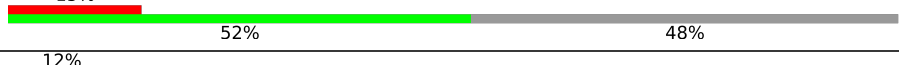





Metric	Whole archive (#Entries)	EM structures (#Entries)
Clashscore	158937	4297
Ramachandran outliers	154571	4023

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	16% 52% 48%
1	AA	258	14% 84% 15%
1	AB	258	84% 15%
1	AC	258	18% 61% 39%
1	B	258	14% 84% 15%
1	BA	258	83% 15%
1	BB	258	13% 84% 16%
1	C	258	83% 15%
1	CA	258	84% 15%

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Mol	Chain	Length	Quality of chain
1	CB	258	
1	D	258	
1	DA	258	
1	DB	258	
1	E	258	
1	EA	258	
1	EB	258	
1	F	258	
1	FA	258	
1	FB	258	
1	G	258	
1	GA	258	
1	GB	258	
1	H	258	
1	HA	258	
1	HB	258	
1	I	258	
1	IA	258	
1	IB	258	
1	J	258	
1	JA	258	
1	JB	258	
1	K	258	
1	KA	258	
1	KB	258	



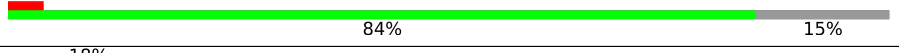
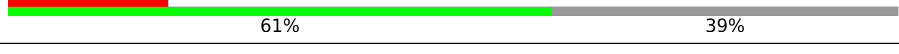



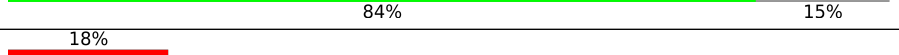
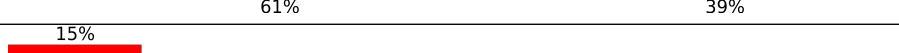
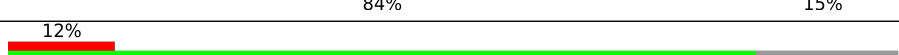


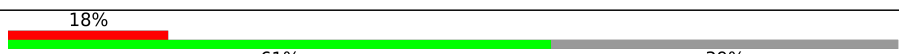
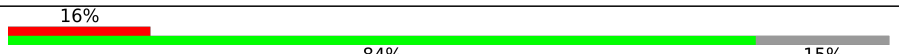
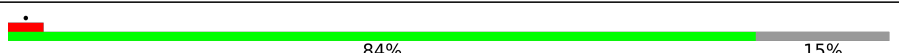
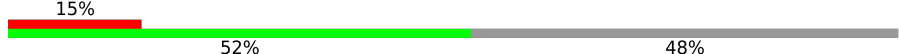



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Mol	Chain	Length	Quality of chain		
1	L	258	18%	61%	39%
1	LA	258	16%	52%	48%
1	LB	258	•	83%	• 15%
1	M	258	15%	52%	48%
1	MA	258	15%	84%	15%
1	MB	258	5%	84%	15%
1	N	258	16%	84%	15%
1	NA	258	•	83%	• 15%
1	NB	258	12%	84%	16%
1	O	258	•	83%	• 15%
1	OA	258	•	84%	15%
1	OB	258	18%	61%	39%
1	P	258	5%	84%	15%
1	PA	258	12%	84%	16%
1	PB	258	15%	52%	48%
1	Q	258	13%	84%	16%
1	QA	258	18%	61%	39%
1	QB	258	16%	84%	15%
1	R	258	18%	61%	39%
1	RA	258	15%	52%	48%
1	RB	258	•	83%	• 15%
1	S	258	16%	52%	48%
1	SA	258	14%	84%	15%
1	SB	258	5%	84%	15%
1	T	258	15%	84%	15%

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Mol	Chain	Length	Quality of chain
1	TA	258	 13% 83% 15%
1	TB	258	 18% 61% 39%
1	UA	258	 12% 84% 16%
1	UB	258	 16% 52% 48%
1	V	258	 5% 84% 15%
1	VA	258	 18% 61% 39%
1	VB	258	 15% 84% 15%
1	W	258	 12% 84% 16%
1	WA	258	 15% 52% 48%
1	WB	258	 18% 61% 39%
1	X	258	 15% 84% 15%
1	XA	258	 12% 84% 16%
1	XB	258	 15% 52% 48%
1	Y	258	 18% 61% 39%
1	YA	258	 16% 84% 15%
1	YB	258	 18% 61% 39%
1	Z	258	 15% 52% 48%
1	ZA	258	 13% 84% 16%
1	ZB	258	 13% 84% 16%

## 2 Entry composition [i](#)

There is only 1 type of molecule in this entry. The entry contains 75205 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
			Total	C	N	O		
1	A	133	661	395	133	133	0	0
1	B	219	1086	648	219	219	0	0
1	C	219	1086	648	219	219	0	0
1	D	219	1086	648	219	219	0	0
1	E	218	1081	645	218	218	0	0
1	F	158	785	469	158	158	0	0
1	G	133	661	395	133	133	0	0
1	H	219	1086	648	219	219	0	0
1	I	219	1086	648	219	219	0	0
1	J	219	1086	648	219	219	0	0
1	K	218	1081	645	218	218	0	0
1	L	158	785	469	158	158	0	0
1	M	133	661	395	133	133	0	0
1	N	219	1086	648	219	219	0	0
1	O	219	1086	648	219	219	0	0
1	P	219	1086	648	219	219	0	0
1	Q	218	1081	645	218	218	0	0

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

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

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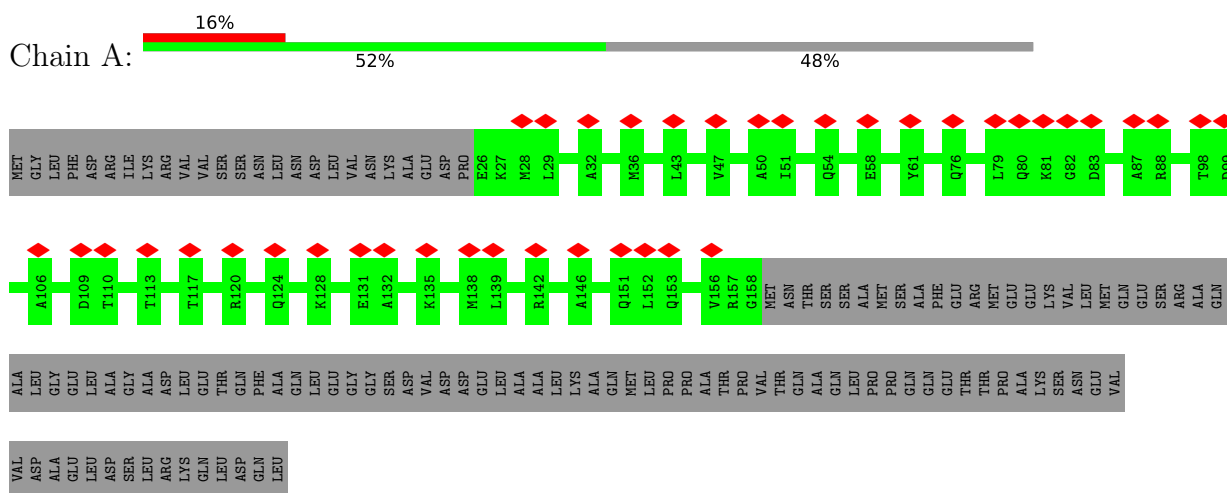
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Mol	Chain	Residues	Atoms				AltConf	Trace
1	IB	158	Total	C	N	O	0	0
			785	469	158	158		
1	JB	133	Total	C	N	O	0	0
			661	395	133	133		
1	KB	219	Total	C	N	O	0	0
			1086	648	219	219		
1	LB	219	Total	C	N	O	0	0
			1086	648	219	219		
1	MB	219	Total	C	N	O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

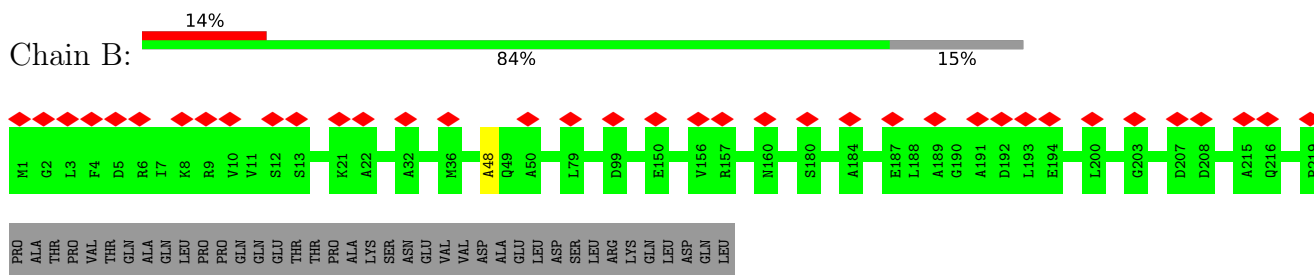
### 3 Residue-property plots

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

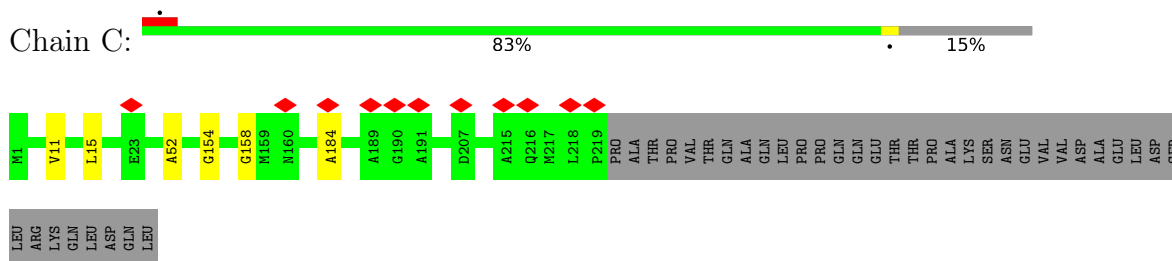
- Molecule 1: Vipp1



- Molecule 1: Vipp1



- Molecule 1: Vipp1

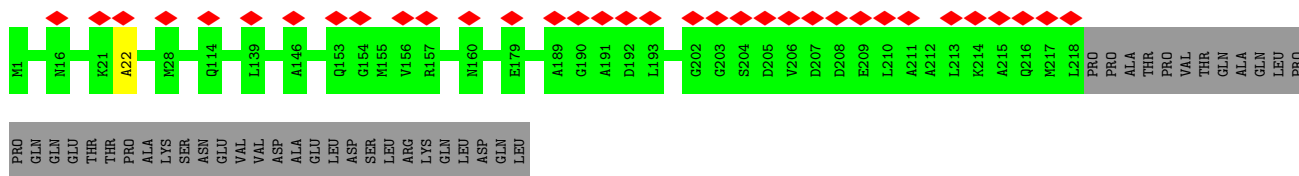


- Molecule 1: Vipp1

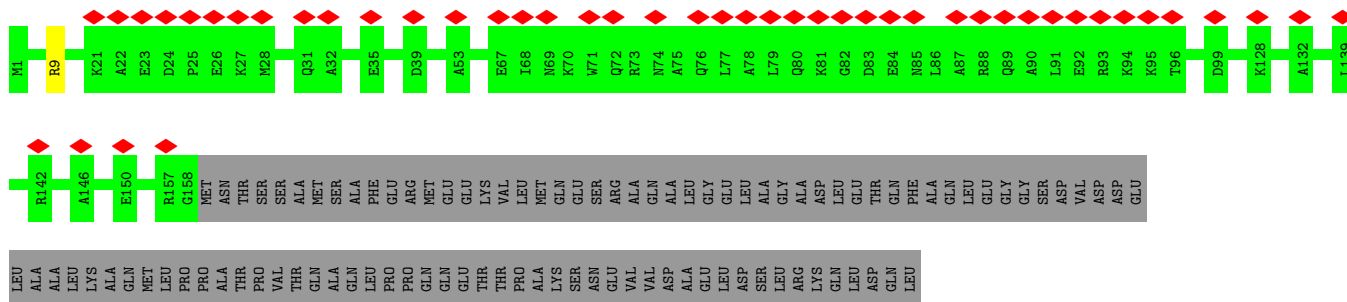




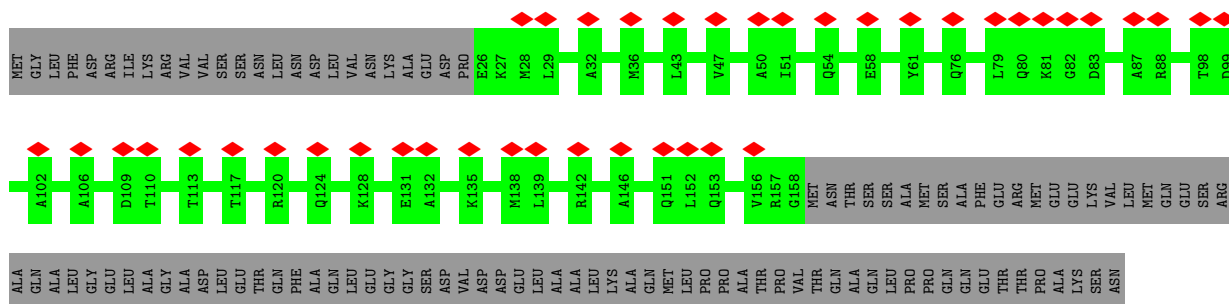




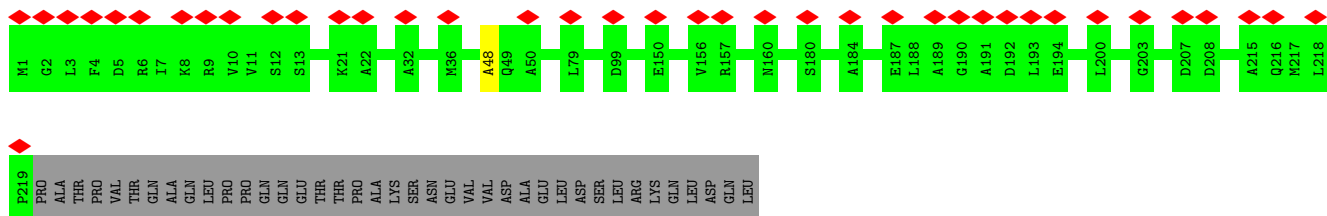
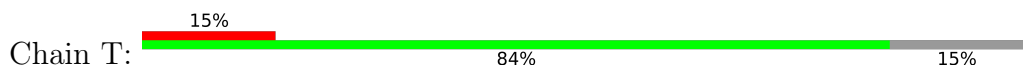
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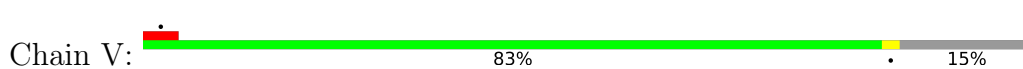
- Molecule 1: Vipp1



- Molecule 1: Vipp1



- Molecule 1: Vipp1



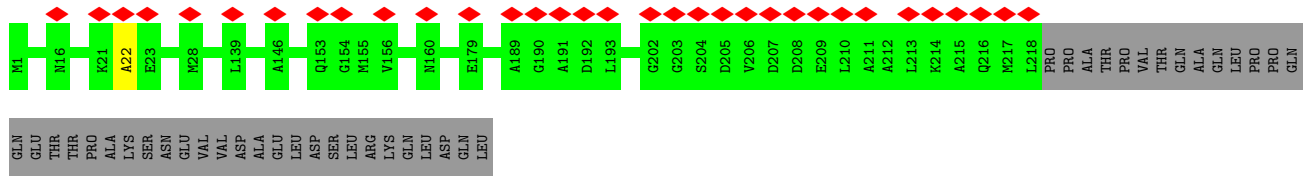
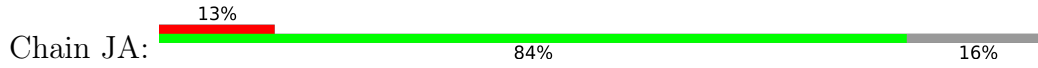




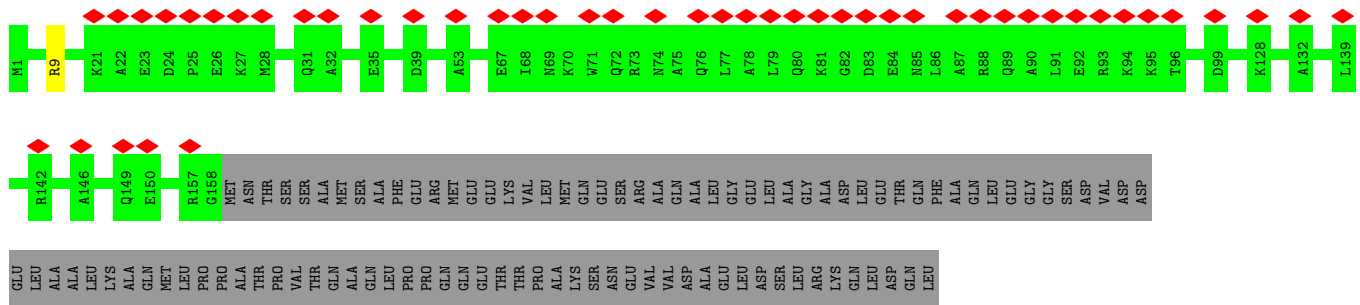




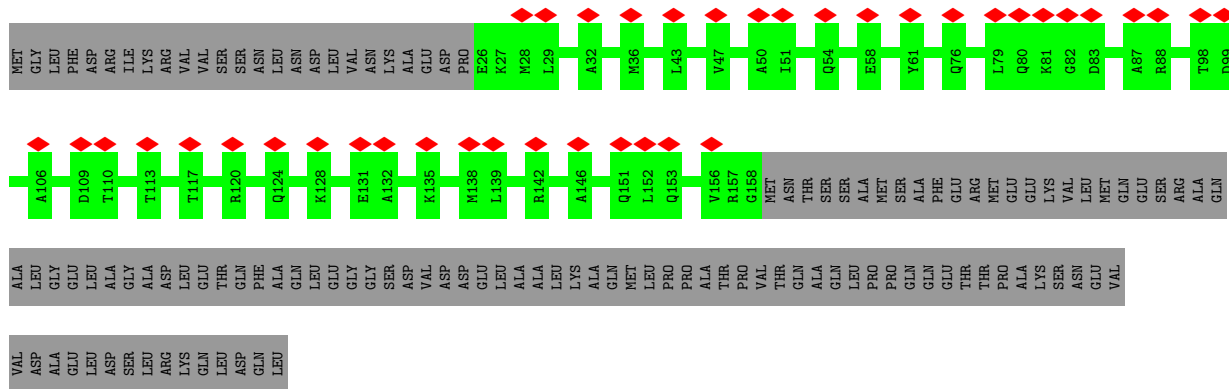
• Molecule 1: Vipp1



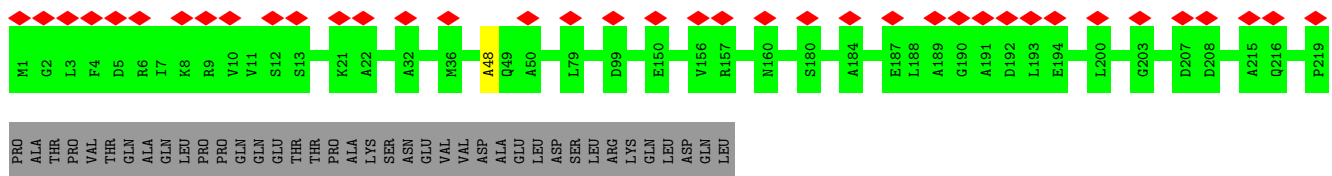
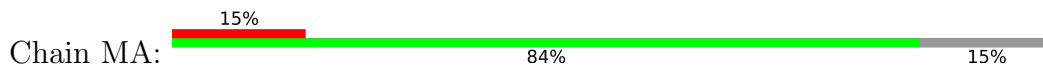
• Molecule 1: Vipp1



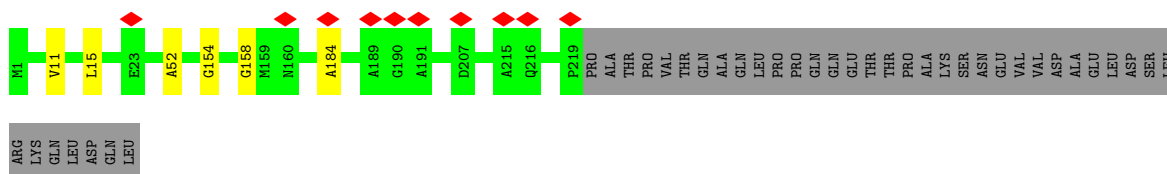
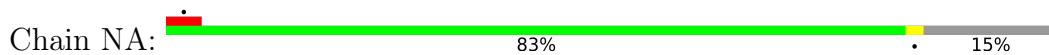
• Molecule 1: Vipp1



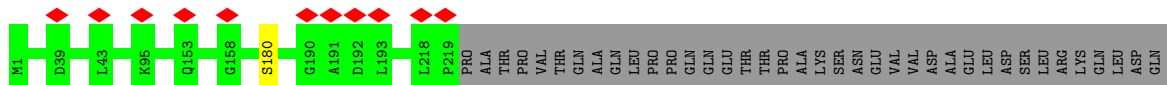
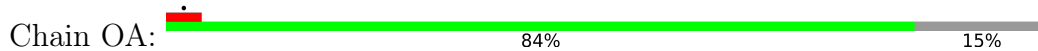
• Molecule 1: Vipp1



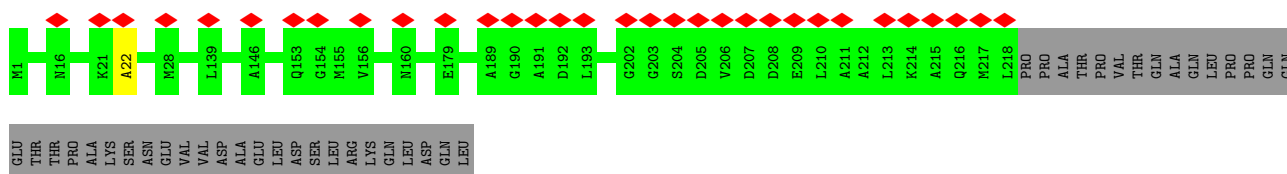
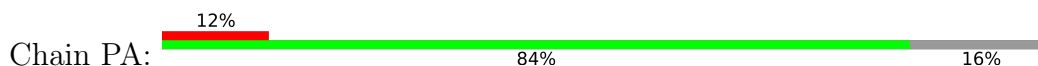
• Molecule 1: Vipp1



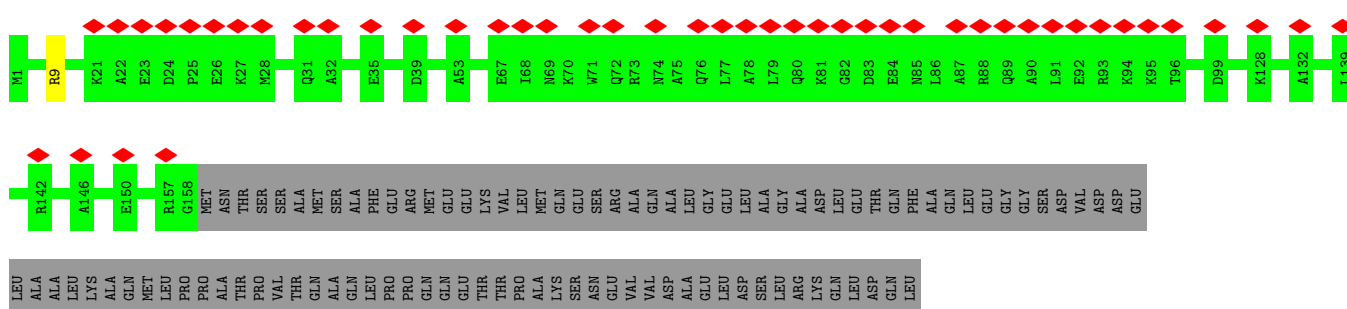
• Molecule 1: Vipp1



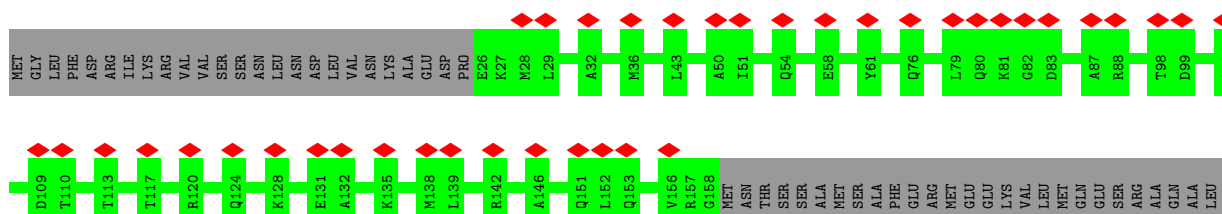
• Molecule 1: Vipp1



• Molecule 1: Vipp1




• Molecule 1: Vipp1

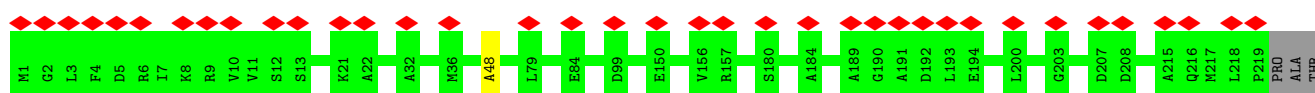


GLY GLU  
GLU LEU  
ALA ASP  
GLY SER  
ALA LEU  
ASP ARG  
LEU LYS  
GLU GLN  
THR LEU  
GLN THR  
PHE ASP  
GLN ASP  
ALA GLN  
LEU LEU  
GLY GLY  
SER SER  
ASP ASP  
VAL VAL  
GLU GLU  
ALA LEU  
ALA LEU  
LYS LEU  
ALA ARG  
GLN MET  
LEU MET  
PRO PRO  
PRO PRO  
ALA THR  
THR THR  
PRO VAL  
THR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

ALA  
GLU  
LEU  
ASP  
SER  
LEU  
ARG  
LYS  
GLN  
LEU  
LEU


- Molecule 1: Vipp1

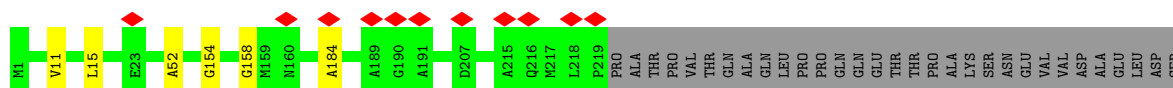
Chain SA: 



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


- Molecule 1: Vipp1

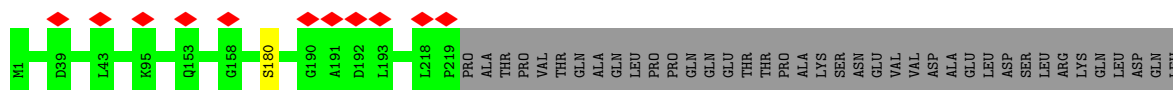
Chain TA: 




LEU  
ARG  
LYS  
GLN  
LEU  
ASP  
GLN  
LEU

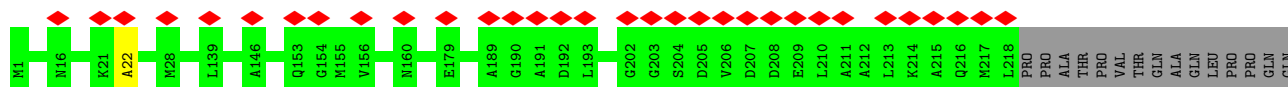
- Molecule 1: Vipp1

Chain UA: 



- Molecule 1: Vipp1

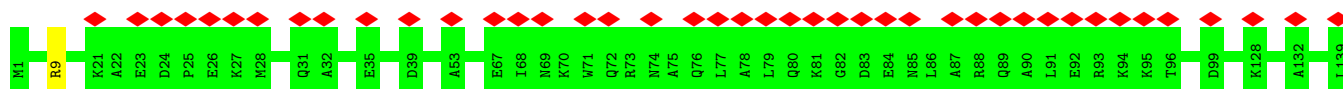
Chain VA: 

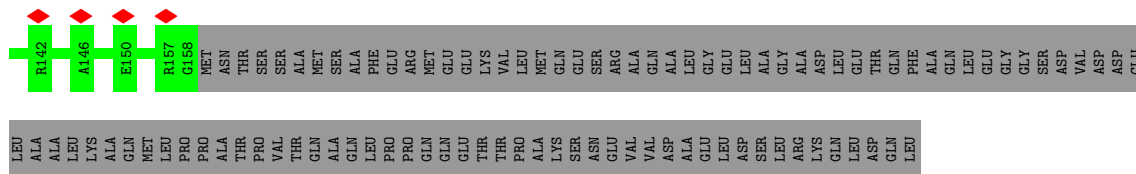


GLU  
THR  
PRO  
ALA  
LYS  
SER  
ASN  
VAL  
VAL  
ASP  
ALA  
GLU  
LEU  
ASP  
SER  
LEU  
ARG  
LYS  
GLN  
LEU  
ASP  
GLN  
LEU

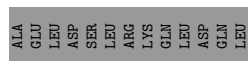
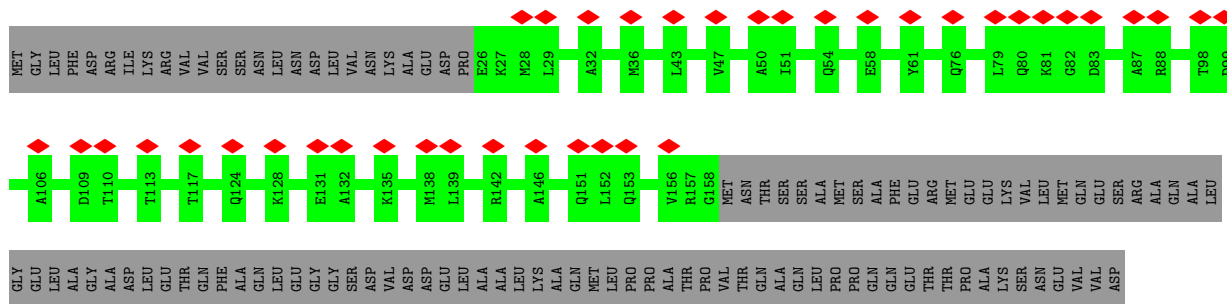
- Molecule 1: Vipp1

Chain WA: 

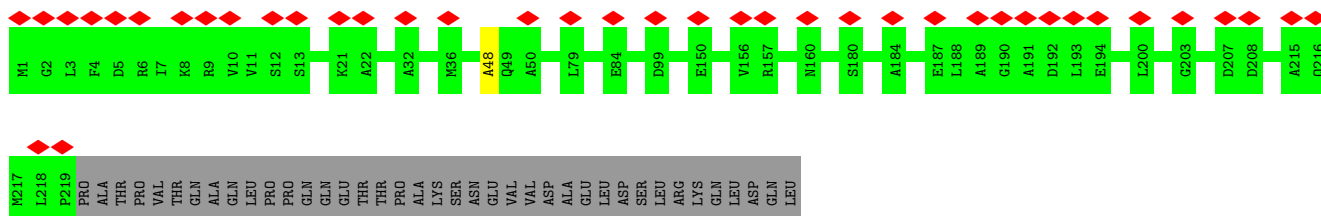
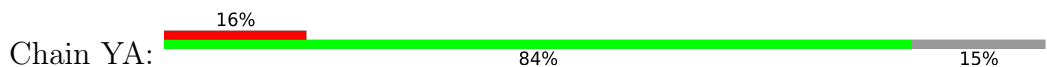




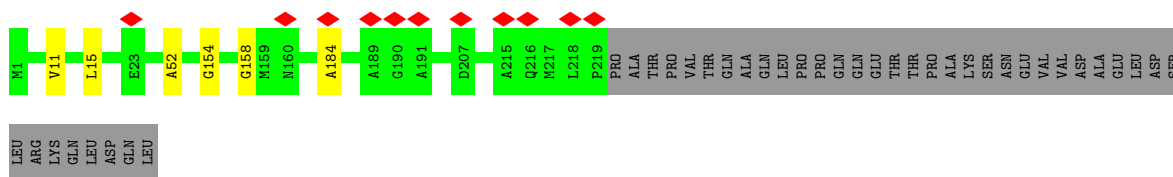
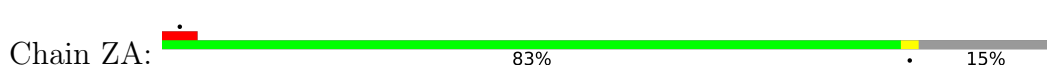
• Molecule 1: Vipp1



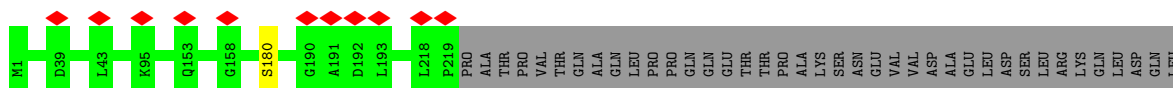
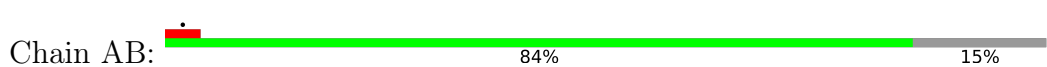
• Molecule 1: Vipp1



• Molecule 1: Vipp1

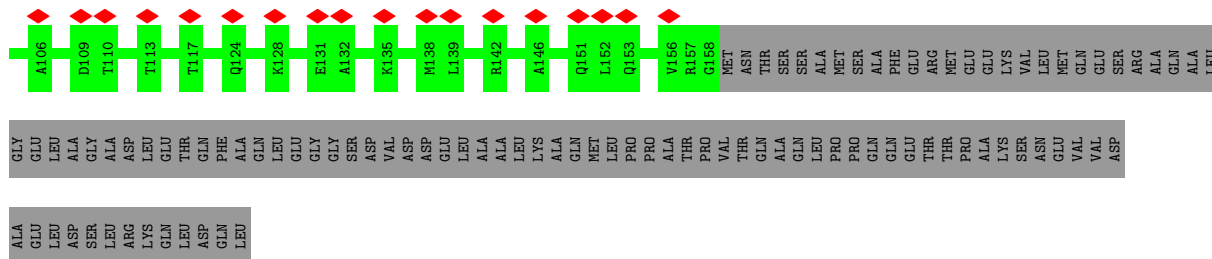


• Molecule 1: Vipp1

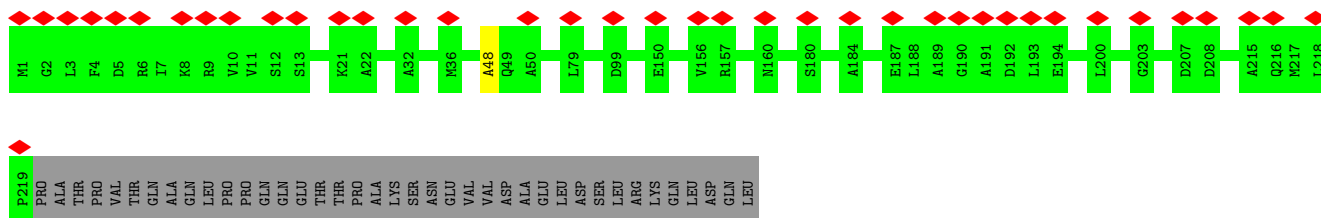
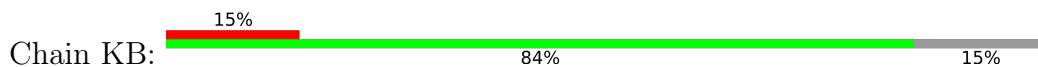




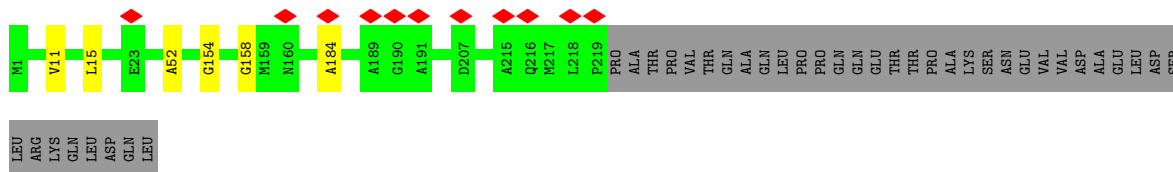
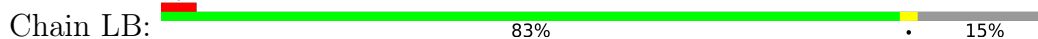




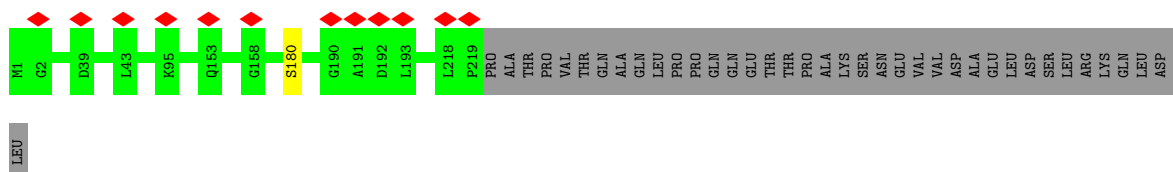
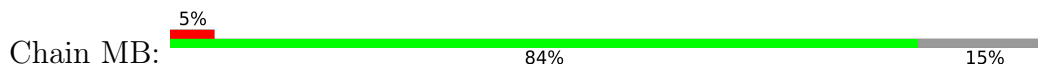
● Molecule 1: Vipp1



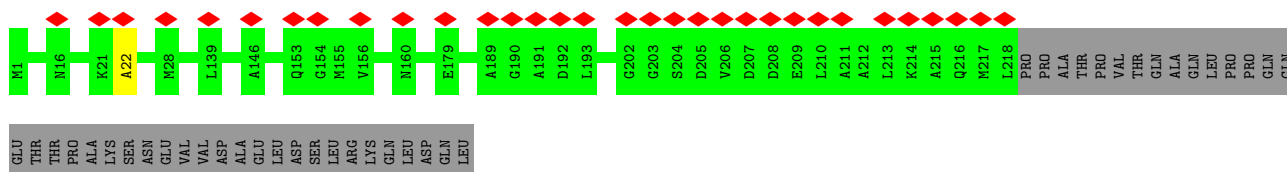
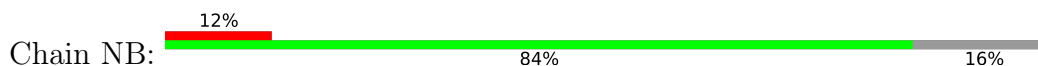
● Molecule 1: Vipp1



● Molecule 1: Vipp1



● Molecule 1: Vipp1











## 4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	6920	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	0.055	Depositor
Minimum map value	-0.009	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.003	Depositor
Recommended contour level	0.0187	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	0.19	0/660	0.37	0/920
1	AA	0.18	0/1085	0.35	0/1512
1	AB	0.19	0/1085	0.39	0/1512
1	AC	0.19	0/784	0.39	0/1093
1	B	0.18	0/1085	0.35	0/1512
1	BA	0.20	0/1085	0.39	0/1512
1	BB	0.20	0/1080	0.41	0/1505
1	C	0.20	0/1085	0.39	0/1512
1	CA	0.19	0/1085	0.39	0/1512
1	CB	0.20	0/784	0.39	0/1093
1	D	0.19	0/1085	0.39	0/1512
1	DA	0.20	0/1080	0.41	0/1505
1	DB	0.20	0/660	0.37	0/920
1	E	0.20	0/1080	0.41	0/1505
1	EA	0.20	0/784	0.39	0/1093
1	EB	0.18	0/1085	0.35	0/1512
1	F	0.20	0/784	0.39	0/1093
1	FA	0.19	0/660	0.37	0/920
1	FB	0.20	0/1085	0.39	0/1512
1	G	0.19	0/660	0.37	0/920
1	GA	0.18	0/1085	0.35	0/1512
1	GB	0.19	0/1085	0.39	0/1512
1	H	0.18	0/1085	0.35	0/1512
1	HA	0.20	0/1085	0.39	0/1512
1	HB	0.20	0/1080	0.41	0/1505
1	I	0.20	0/1085	0.39	0/1512
1	IA	0.19	0/1085	0.39	0/1512
1	IB	0.20	0/784	0.39	0/1093
1	J	0.19	0/1085	0.39	0/1512
1	JA	0.20	0/1080	0.41	0/1505
1	JB	0.20	0/660	0.37	0/920
1	K	0.20	0/1080	0.41	0/1505
1	KA	0.20	0/784	0.39	0/1093
1	KB	0.18	0/1085	0.35	0/1512

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
1	L	0.20	0/784	0.39	0/1093
1	LA	0.20	0/660	0.37	0/920
1	LB	0.20	0/1085	0.39	0/1512
1	M	0.19	0/660	0.37	0/920
1	MA	0.18	0/1085	0.35	0/1512
1	MB	0.19	0/1085	0.39	0/1512
1	N	0.18	0/1085	0.35	0/1512
1	NA	0.20	0/1085	0.39	0/1512
1	NB	0.20	0/1080	0.41	0/1505
1	O	0.20	0/1085	0.39	0/1512
1	OA	0.19	0/1085	0.39	0/1512
1	OB	0.20	0/784	0.39	0/1093
1	P	0.19	0/1085	0.39	0/1512
1	PA	0.20	0/1080	0.41	0/1505
1	PB	0.20	0/660	0.37	0/920
1	Q	0.20	0/1080	0.41	0/1505
1	QA	0.20	0/784	0.39	0/1093
1	QB	0.18	0/1085	0.35	0/1512
1	R	0.20	0/784	0.39	0/1093
1	RA	0.19	0/660	0.37	0/920
1	RB	0.20	0/1085	0.39	0/1512
1	S	0.20	0/660	0.37	0/920
1	SA	0.18	0/1085	0.35	0/1512
1	SB	0.19	0/1085	0.39	0/1512
1	T	0.18	0/1085	0.35	0/1512
1	TA	0.20	0/1085	0.39	0/1512
1	TB	0.20	0/1080	0.41	0/1505
1	UA	0.19	0/1085	0.39	0/1512
1	UB	0.20	0/784	0.39	0/1093
1	V	0.20	0/1085	0.39	0/1512
1	VA	0.20	0/1080	0.41	0/1505
1	VB	0.20	0/660	0.37	0/920
1	W	0.19	0/1085	0.39	0/1512
1	WA	0.19	0/784	0.39	0/1093
1	WB	0.18	0/1085	0.35	0/1512
1	X	0.20	0/1080	0.41	0/1505
1	XA	0.19	0/660	0.37	0/920
1	XB	0.20	0/1085	0.39	0/1512
1	Y	0.20	0/784	0.39	0/1093
1	YA	0.18	0/1085	0.35	0/1512
1	YB	0.19	0/1085	0.39	0/1512
1	Z	0.19	0/660	0.37	0/920
1	ZA	0.20	0/1085	0.39	0/1512

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
1	ZB	0.20	0/1080	0.41	0/1505
All	All	0.19	0/75127	0.38	0/104702

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

## 5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	661	0	320	0	0
1	AA	1086	0	531	1	0
1	AB	1086	0	531	3	0
1	AC	785	0	376	2	0
1	B	1086	0	531	1	0
1	BA	1086	0	531	12	0
1	BB	1081	0	530	2	0
1	C	1086	0	531	12	0
1	CA	1086	0	531	3	0
1	CB	785	0	376	2	0
1	D	1086	0	531	3	0
1	DA	1081	0	530	2	0
1	DB	661	0	320	0	0
1	E	1081	0	530	2	0
1	EA	785	0	376	2	0
1	EB	1086	0	531	1	0
1	F	785	0	376	2	0
1	FA	661	0	320	0	0
1	FB	1086	0	531	13	0
1	G	661	0	320	0	0
1	GA	1086	0	531	1	0
1	GB	1086	0	531	3	0
1	H	1086	0	531	1	0
1	HA	1086	0	531	11	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	HB	1081	0	530	2	0
1	I	1086	0	531	13	0
1	IA	1086	0	531	3	0
1	IB	785	0	376	2	0
1	J	1086	0	531	3	0
1	JA	1081	0	530	2	0
1	JB	661	0	320	0	0
1	K	1081	0	530	2	0
1	KA	785	0	376	2	0
1	KB	1086	0	531	1	0
1	L	785	0	376	2	0
1	LA	661	0	320	0	0
1	LB	1086	0	531	13	0
1	M	661	0	320	0	0
1	MA	1086	0	531	1	0
1	MB	1086	0	531	3	0
1	N	1086	0	531	1	0
1	NA	1086	0	531	12	0
1	NB	1081	0	530	2	0
1	O	1086	0	531	13	0
1	OA	1086	0	531	3	0
1	OB	785	0	376	2	0
1	P	1086	0	531	3	0
1	PA	1081	0	530	2	0
1	PB	661	0	320	0	0
1	Q	1081	0	530	2	0
1	QA	785	0	376	2	0
1	QB	1086	0	531	1	0
1	R	785	0	376	2	0
1	RA	661	0	320	0	0
1	RB	1086	0	531	13	0
1	S	661	0	320	0	0
1	SA	1086	0	531	1	0
1	SB	1086	0	531	3	0
1	T	1086	0	531	1	0
1	TA	1086	0	531	12	0
1	TB	1081	0	530	2	0
1	UA	1086	0	531	3	0
1	UB	785	0	376	2	0
1	V	1086	0	531	12	0
1	VA	1081	0	530	2	0
1	VB	661	0	320	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	W	1086	0	531	3	0
1	WA	785	0	376	2	0
1	WB	1086	0	531	1	0
1	X	1081	0	530	2	0
1	XA	661	0	320	0	0
1	XB	1086	0	531	12	0
1	Y	785	0	376	2	0
1	YA	1086	0	531	1	0
1	YB	1086	0	531	3	0
1	Z	661	0	320	0	0
1	ZA	1086	0	531	13	0
1	ZB	1081	0	530	2	0
All	All	75205	0	36647	187	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.

All (187) 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:BA:154:GLY:O	1:BA:158:GLY:CA	2.11	0.98
1:XB:154:GLY:O	1:XB:158:GLY:CA	2.11	0.98
1:HA:154:GLY:O	1:HA:158:GLY:CA	2.11	0.98
1:C:154:GLY:O	1:C:158:GLY:CA	2.11	0.98
1:RB:154:GLY:O	1:RB:158:GLY:CA	2.12	0.98
1:ZA:154:GLY:O	1:ZA:158:GLY:CA	2.11	0.98
1:V:154:GLY:O	1:V:158:GLY:CA	2.11	0.98
1:FB:154:GLY:O	1:FB:158:GLY:CA	2.11	0.98
1:I:154:GLY:O	1:I:158:GLY:CA	2.11	0.97
1:NA:154:GLY:O	1:NA:158:GLY:CA	2.11	0.97
1:TA:154:GLY:O	1:TA:158:GLY:CA	2.11	0.97
1:O:154:GLY:O	1:O:158:GLY:CA	2.11	0.97
1:LB:154:GLY:O	1:LB:158:GLY:CA	2.11	0.97
1:V:154:GLY:O	1:V:158:GLY:HA3	1.66	0.96
1:RB:154:GLY:O	1:RB:158:GLY:HA3	1.66	0.96
1:I:154:GLY:O	1:I:158:GLY:HA3	1.66	0.96
1:C:154:GLY:O	1:C:158:GLY:HA3	1.66	0.96
1:O:154:GLY:O	1:O:158:GLY:HA3	1.66	0.96
1:XB:154:GLY:O	1:XB:158:GLY:HA3	1.66	0.95
1:HA:154:GLY:O	1:HA:158:GLY:HA3	1.66	0.95
1:FB:154:GLY:O	1:FB:158:GLY:HA3	1.66	0.95

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:FB:154:GLY:O	1:FB:158:GLY:N	2.01	0.94
1:XB:154:GLY:O	1:XB:158:GLY:N	2.00	0.94
1:NA:154:GLY:O	1:NA:158:GLY:HA3	1.66	0.94
1:ZA:154:GLY:O	1:ZA:158:GLY:HA3	1.66	0.94
1:BA:154:GLY:O	1:BA:158:GLY:HA3	1.66	0.94
1:I:154:GLY:O	1:I:158:GLY:N	2.01	0.94
1:TA:154:GLY:O	1:TA:158:GLY:N	2.00	0.94
1:LB:154:GLY:O	1:LB:158:GLY:HA3	1.66	0.94
1:RB:154:GLY:O	1:RB:158:GLY:N	2.00	0.94
1:LB:154:GLY:O	1:LB:158:GLY:N	2.01	0.94
1:V:154:GLY:O	1:V:158:GLY:N	2.00	0.93
1:HA:154:GLY:O	1:HA:158:GLY:N	2.00	0.93
1:NA:154:GLY:O	1:NA:158:GLY:N	2.01	0.93
1:ZA:154:GLY:O	1:ZA:158:GLY:N	2.00	0.93
1:O:154:GLY:O	1:O:158:GLY:N	2.00	0.93
1:C:154:GLY:O	1:C:158:GLY:N	2.00	0.93
1:BA:154:GLY:O	1:BA:158:GLY:N	2.01	0.92
1:TA:154:GLY:O	1:TA:158:GLY:HA3	1.66	0.91
1:NA:52:ALA:CB	1:AB:180:SER:CB	2.51	0.89
1:D:180:SER:CB	1:RB:52:ALA:CB	2.51	0.89
1:HA:52:ALA:CB	1:UA:180:SER:CB	2.51	0.89
1:I:52:ALA:CB	1:W:180:SER:CB	2.51	0.89
1:C:52:ALA:CB	1:P:180:SER:CB	2.51	0.88
1:TA:52:ALA:CB	1:GB:180:SER:CB	2.51	0.88
1:BA:52:ALA:CB	1:OA:180:SER:CB	2.51	0.88
1:LB:52:ALA:CB	1:YB:180:SER:CB	2.51	0.88
1:ZA:52:ALA:CB	1:MB:180:SER:CB	2.51	0.88
1:O:52:ALA:CB	1:CA:180:SER:CB	2.51	0.88
1:J:180:SER:CB	1:XB:52:ALA:CB	2.51	0.88
1:V:52:ALA:CB	1:IA:180:SER:CB	2.51	0.88
1:FB:52:ALA:CB	1:SB:180:SER:CB	2.51	0.87
1:O:52:ALA:HB2	1:CA:180:SER:CB	2.13	0.77
1:J:180:SER:CB	1:XB:52:ALA:HB2	2.13	0.77
1:ZA:52:ALA:HB2	1:MB:180:SER:CB	2.14	0.77
1:FB:52:ALA:HB2	1:SB:180:SER:CB	2.13	0.77
1:D:180:SER:CB	1:RB:52:ALA:HB2	2.14	0.77
1:HA:52:ALA:HB2	1:UA:180:SER:CB	2.14	0.77
1:BA:52:ALA:HB2	1:OA:180:SER:CB	2.14	0.76
1:TA:52:ALA:HB2	1:GB:180:SER:CB	2.14	0.76
1:I:52:ALA:HB2	1:W:180:SER:CB	2.14	0.76
1:V:52:ALA:HB2	1:IA:180:SER:CB	2.14	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:52:ALA:HB2	1:P:180:SER:CB	2.14	0.76
1:NA:52:ALA:HB2	1:AB:180:SER:CB	2.14	0.75
1:LB:52:ALA:HB2	1:YB:180:SER:CB	2.15	0.75
1:LB:52:ALA:HB1	1:YB:180:SER:CB	2.23	0.68
1:TA:52:ALA:HB1	1:GB:180:SER:CB	2.25	0.67
1:D:180:SER:CB	1:RB:52:ALA:HB1	2.25	0.67
1:ZA:52:ALA:HB1	1:MB:180:SER:CB	2.25	0.66
1:NA:52:ALA:HB1	1:AB:180:SER:CB	2.25	0.66
1:FB:52:ALA:HB1	1:SB:180:SER:CB	2.25	0.66
1:C:52:ALA:HB1	1:P:180:SER:CB	2.24	0.65
1:V:52:ALA:HB1	1:IA:180:SER:CB	2.24	0.65
1:HA:52:ALA:HB1	1:UA:180:SER:CB	2.24	0.65
1:BA:52:ALA:HB1	1:OA:180:SER:CB	2.24	0.65
1:J:180:SER:CB	1:XB:52:ALA:HB1	2.26	0.65
1:I:52:ALA:HB1	1:W:180:SER:CB	2.25	0.64
1:O:52:ALA:HB1	1:CA:180:SER:CB	2.25	0.64
1:TB:22:ALA:HB2	1:UB:9:ARG:CB	2.29	0.63
1:BB:22:ALA:HB2	1:CB:9:ARG:CB	2.29	0.62
1:HB:22:ALA:HB2	1:IB:9:ARG:CB	2.29	0.62
1:NB:22:ALA:HB2	1:OB:9:ARG:CB	2.29	0.62
1:ZB:22:ALA:HB2	1:AC:9:ARG:CB	2.29	0.62
1:PA:22:ALA:HB2	1:QA:9:ARG:CB	2.29	0.62
1:VA:22:ALA:HB2	1:WA:9:ARG:CB	2.29	0.62
1:E:22:ALA:HB2	1:F:9:ARG:CB	2.29	0.62
1:JA:22:ALA:HB2	1:KA:9:ARG:CB	2.29	0.62
1:DA:22:ALA:HB2	1:EA:9:ARG:CB	2.29	0.61
1:X:22:ALA:HB2	1:Y:9:ARG:CB	2.29	0.61
1:K:22:ALA:HB2	1:L:9:ARG:CB	2.29	0.61
1:Q:22:ALA:HB2	1:R:9:ARG:CB	2.29	0.60
1:BA:154:GLY:HA2	1:BA:158:GLY:HA2	1.85	0.59
1:HA:154:GLY:HA2	1:HA:158:GLY:HA2	1.85	0.59
1:V:154:GLY:HA2	1:V:158:GLY:HA2	1.85	0.59
1:NA:154:GLY:HA2	1:NA:158:GLY:HA2	1.85	0.59
1:LB:154:GLY:HA2	1:LB:158:GLY:HA2	1.85	0.58
1:TA:154:GLY:HA2	1:TA:158:GLY:HA2	1.85	0.58
1:RB:154:GLY:HA2	1:RB:158:GLY:HA2	1.85	0.58
1:O:154:GLY:HA2	1:O:158:GLY:HA2	1.85	0.58
1:FB:154:GLY:HA2	1:FB:158:GLY:HA2	1.85	0.58
1:XB:154:GLY:HA2	1:XB:158:GLY:HA2	1.85	0.58
1:ZA:154:GLY:HA2	1:ZA:158:GLY:HA2	1.85	0.57
1:C:154:GLY:HA2	1:C:158:GLY:HA2	1.85	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:154:GLY:HA2	1:I:158:GLY:HA2	1.85	0.57
1:PA:22:ALA:CB	1:QA:9:ARG:CB	2.85	0.55
1:HB:22:ALA:CB	1:IB:9:ARG:CB	2.85	0.55
1:NB:22:ALA:CB	1:OB:9:ARG:CB	2.85	0.55
1:TB:22:ALA:CB	1:UB:9:ARG:CB	2.85	0.55
1:JA:22:ALA:CB	1:KA:9:ARG:CB	2.85	0.54
1:VA:22:ALA:CB	1:WA:9:ARG:CB	2.85	0.54
1:ZB:22:ALA:CB	1:AC:9:ARG:CB	2.85	0.54
1:E:22:ALA:CB	1:F:9:ARG:CB	2.85	0.54
1:DA:22:ALA:CB	1:EA:9:ARG:CB	2.85	0.54
1:HA:11:VAL:O	1:HA:15:LEU:N	2.34	0.54
1:I:11:VAL:O	1:I:15:LEU:N	2.34	0.54
1:X:22:ALA:CB	1:Y:9:ARG:CB	2.85	0.54
1:BB:22:ALA:CB	1:CB:9:ARG:CB	2.85	0.54
1:Q:22:ALA:CB	1:R:9:ARG:CB	2.85	0.54
1:K:22:ALA:CB	1:L:9:ARG:CB	2.85	0.53
1:C:11:VAL:O	1:C:15:LEU:N	2.34	0.53
1:NA:11:VAL:O	1:NA:15:LEU:N	2.34	0.53
1:XB:11:VAL:O	1:XB:15:LEU:N	2.34	0.51
1:TA:11:VAL:O	1:TA:15:LEU:N	2.34	0.51
1:V:11:VAL:O	1:V:15:LEU:N	2.34	0.50
1:BA:11:VAL:O	1:BA:15:LEU:N	2.34	0.50
1:RB:11:VAL:O	1:RB:15:LEU:N	2.34	0.50
1:O:11:VAL:O	1:O:15:LEU:N	2.34	0.49
1:LB:11:VAL:O	1:LB:15:LEU:N	2.34	0.49
1:ZA:11:VAL:O	1:ZA:15:LEU:N	2.34	0.48
1:FB:11:VAL:O	1:FB:15:LEU:N	2.34	0.48
1:NA:154:GLY:HA2	1:NA:158:GLY:CA	2.45	0.47
1:TA:154:GLY:HA2	1:TA:158:GLY:CA	2.45	0.47
1:HA:154:GLY:HA2	1:HA:158:GLY:CA	2.45	0.47
1:ZA:154:GLY:HA2	1:ZA:158:GLY:CA	2.45	0.47
1:I:154:GLY:HA2	1:I:158:GLY:CA	2.45	0.46
1:C:154:GLY:HA2	1:C:158:GLY:CA	2.45	0.46
1:BA:154:GLY:HA2	1:BA:158:GLY:CA	2.45	0.46
1:FB:154:GLY:HA2	1:FB:158:GLY:CA	2.45	0.46
1:XB:154:GLY:HA2	1:XB:158:GLY:CA	2.45	0.46
1:I:154:GLY:CA	1:I:158:GLY:HA2	2.46	0.46
1:O:154:GLY:CA	1:O:158:GLY:HA2	2.46	0.46
1:LB:154:GLY:CA	1:LB:158:GLY:HA2	2.46	0.46
1:RB:154:GLY:CA	1:RB:158:GLY:HA2	2.46	0.46
1:V:154:GLY:HA2	1:V:158:GLY:CA	2.45	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:RB:154:GLY:HA2	1:RB:158:GLY:CA	2.45	0.46
1:BA:154:GLY:CA	1:BA:158:GLY:HA2	2.46	0.46
1:ZA:154:GLY:CA	1:ZA:158:GLY:HA2	2.46	0.46
1:HA:154:GLY:CA	1:HA:158:GLY:HA2	2.46	0.45
1:TA:154:GLY:CA	1:TA:158:GLY:HA2	2.46	0.45
1:XB:154:GLY:CA	1:XB:158:GLY:HA2	2.46	0.45
1:C:154:GLY:CA	1:C:158:GLY:HA2	2.46	0.45
1:I:154:GLY:C	1:I:158:GLY:CA	2.84	0.45
1:V:154:GLY:CA	1:V:158:GLY:HA2	2.46	0.45
1:LB:154:GLY:HA2	1:LB:158:GLY:CA	2.45	0.45
1:O:154:GLY:HA2	1:O:158:GLY:CA	2.45	0.45
1:FB:154:GLY:CA	1:FB:158:GLY:HA2	2.46	0.45
1:O:154:GLY:C	1:O:158:GLY:CA	2.84	0.44
1:NA:154:GLY:CA	1:NA:158:GLY:HA2	2.46	0.44
1:FB:154:GLY:C	1:FB:158:GLY:CA	2.84	0.44
1:EB:48:ALA:HA	1:RB:184:ALA:HB1	1.99	0.44
1:GA:48:ALA:HA	1:TA:184:ALA:HB1	2.00	0.43
1:SA:48:ALA:HA	1:FB:184:ALA:HB1	2.00	0.43
1:AA:48:ALA:HA	1:NA:184:ALA:HB1	2.00	0.43
1:T:48:ALA:HA	1:HA:184:ALA:HB1	2.00	0.43
1:KB:48:ALA:HA	1:XB:184:ALA:HB1	1.99	0.43
1:V:154:GLY:C	1:V:158:GLY:CA	2.84	0.43
1:MA:48:ALA:HA	1:ZA:184:ALA:HB1	2.01	0.43
1:NA:154:GLY:C	1:NA:158:GLY:CA	2.84	0.42
1:YA:48:ALA:HA	1:LB:184:ALA:HB1	2.01	0.42
1:LB:154:GLY:C	1:LB:158:GLY:CA	2.84	0.42
1:B:48:ALA:HA	1:O:184:ALA:HB1	2.00	0.42
1:I:184:ALA:HB1	1:WB:48:ALA:HA	2.00	0.42
1:N:48:ALA:HA	1:BA:184:ALA:HB1	2.00	0.42
1:RB:154:GLY:C	1:RB:158:GLY:CA	2.84	0.41
1:C:184:ALA:HB1	1:QB:48:ALA:HA	2.01	0.41
1:FB:154:GLY:CA	1:FB:158:GLY:CA	2.99	0.41
1:RB:154:GLY:CA	1:RB:158:GLY:CA	2.99	0.41
1:C:154:GLY:CA	1:C:158:GLY:CA	2.99	0.41
1:I:154:GLY:CA	1:I:158:GLY:CA	2.99	0.41
1:ZA:154:GLY:C	1:ZA:158:GLY:CA	2.84	0.41
1:H:48:ALA:HA	1:V:184:ALA:HB1	2.02	0.40
1:O:154:GLY:CA	1:O:158:GLY:CA	2.99	0.40
1:XB:154:GLY:CA	1:XB:158:GLY:CA	2.99	0.40
1:TA:154:GLY:C	1:TA:158:GLY:CA	2.84	0.40
1:LB:154:GLY:CA	1:LB:158:GLY:CA	2.99	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BA:154:GLY:C	1:BA:158:GLY:CA	2.84	0.40
1:ZA:154:GLY:CA	1:ZA:158:GLY:CA	2.99	0.40

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%)	129 (98%)	2 (2%)	0	100	100
1	AA	217/258 (84%)	210 (97%)	7 (3%)	0	100	100
1	AB	217/258 (84%)	217 (100%)	0	0	100	100
1	AC	156/258 (60%)	156 (100%)	0	0	100	100
1	B	217/258 (84%)	210 (97%)	7 (3%)	0	100	100
1	BA	217/258 (84%)	216 (100%)	1 (0%)	0	100	100
1	BB	216/258 (84%)	216 (100%)	0	0	100	100
1	C	217/258 (84%)	216 (100%)	1 (0%)	0	100	100
1	CA	217/258 (84%)	217 (100%)	0	0	100	100
1	CB	156/258 (60%)	156 (100%)	0	0	100	100
1	D	217/258 (84%)	217 (100%)	0	0	100	100
1	DA	216/258 (84%)	216 (100%)	0	0	100	100
1	DB	131/258 (51%)	129 (98%)	2 (2%)	0	100	100
1	E	216/258 (84%)	216 (100%)	0	0	100	100
1	EA	156/258 (60%)	156 (100%)	0	0	100	100
1	EB	217/258 (84%)	210 (97%)	7 (3%)	0	100	100
1	F	156/258 (60%)	156 (100%)	0	0	100	100
1	FA	131/258 (51%)	129 (98%)	2 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	FB	217/258 (84%)	216 (100%)	1 (0%)	0	100	100
1	G	131/258 (51%)	129 (98%)	2 (2%)	0	100	100
1	GA	217/258 (84%)	210 (97%)	7 (3%)	0	100	100
1	GB	217/258 (84%)	217 (100%)	0	0	100	100
1	H	217/258 (84%)	210 (97%)	7 (3%)	0	100	100
1	HA	217/258 (84%)	216 (100%)	1 (0%)	0	100	100
1	HB	216/258 (84%)	216 (100%)	0	0	100	100
1	I	217/258 (84%)	216 (100%)	1 (0%)	0	100	100
1	IA	217/258 (84%)	217 (100%)	0	0	100	100
1	IB	156/258 (60%)	156 (100%)	0	0	100	100
1	J	217/258 (84%)	217 (100%)	0	0	100	100
1	JA	216/258 (84%)	216 (100%)	0	0	100	100
1	JB	131/258 (51%)	129 (98%)	2 (2%)	0	100	100
1	K	216/258 (84%)	216 (100%)	0	0	100	100
1	KA	156/258 (60%)	156 (100%)	0	0	100	100
1	KB	217/258 (84%)	210 (97%)	7 (3%)	0	100	100
1	L	156/258 (60%)	156 (100%)	0	0	100	100
1	LA	131/258 (51%)	129 (98%)	2 (2%)	0	100	100
1	LB	217/258 (84%)	216 (100%)	1 (0%)	0	100	100
1	M	131/258 (51%)	129 (98%)	2 (2%)	0	100	100
1	MA	217/258 (84%)	210 (97%)	7 (3%)	0	100	100
1	MB	217/258 (84%)	217 (100%)	0	0	100	100
1	N	217/258 (84%)	210 (97%)	7 (3%)	0	100	100
1	NA	217/258 (84%)	216 (100%)	1 (0%)	0	100	100
1	NB	216/258 (84%)	216 (100%)	0	0	100	100
1	O	217/258 (84%)	216 (100%)	1 (0%)	0	100	100
1	OA	217/258 (84%)	217 (100%)	0	0	100	100
1	OB	156/258 (60%)	156 (100%)	0	0	100	100
1	P	217/258 (84%)	217 (100%)	0	0	100	100
1	PA	216/258 (84%)	216 (100%)	0	0	100	100
1	PB	131/258 (51%)	129 (98%)	2 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	Q	216/258 (84%)	216 (100%)	0	0	100	100
1	QA	156/258 (60%)	156 (100%)	0	0	100	100
1	QB	217/258 (84%)	210 (97%)	7 (3%)	0	100	100
1	R	156/258 (60%)	156 (100%)	0	0	100	100
1	RA	131/258 (51%)	129 (98%)	2 (2%)	0	100	100
1	RB	217/258 (84%)	216 (100%)	1 (0%)	0	100	100
1	S	131/258 (51%)	129 (98%)	2 (2%)	0	100	100
1	SA	217/258 (84%)	210 (97%)	7 (3%)	0	100	100
1	SB	217/258 (84%)	217 (100%)	0	0	100	100
1	T	217/258 (84%)	210 (97%)	7 (3%)	0	100	100
1	TA	217/258 (84%)	216 (100%)	1 (0%)	0	100	100
1	TB	216/258 (84%)	216 (100%)	0	0	100	100
1	UA	217/258 (84%)	217 (100%)	0	0	100	100
1	UB	156/258 (60%)	156 (100%)	0	0	100	100
1	V	217/258 (84%)	216 (100%)	1 (0%)	0	100	100
1	VA	216/258 (84%)	216 (100%)	0	0	100	100
1	VB	131/258 (51%)	129 (98%)	2 (2%)	0	100	100
1	W	217/258 (84%)	217 (100%)	0	0	100	100
1	WA	156/258 (60%)	156 (100%)	0	0	100	100
1	WB	217/258 (84%)	210 (97%)	7 (3%)	0	100	100
1	X	216/258 (84%)	216 (100%)	0	0	100	100
1	XA	131/258 (51%)	129 (98%)	2 (2%)	0	100	100
1	XB	217/258 (84%)	216 (100%)	1 (0%)	0	100	100
1	Y	156/258 (60%)	156 (100%)	0	0	100	100
1	YA	217/258 (84%)	210 (97%)	7 (3%)	0	100	100
1	YB	217/258 (84%)	217 (100%)	0	0	100	100
1	Z	131/258 (51%)	129 (98%)	2 (2%)	0	100	100
1	ZA	217/258 (84%)	216 (100%)	1 (0%)	0	100	100
1	ZB	216/258 (84%)	216 (100%)	0	0	100	100
All	All	15002/20124 (74%)	14872 (99%)	130 (1%)	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.

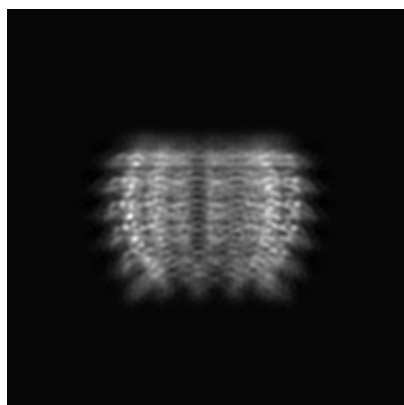
## 6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-11470. These allow visual inspection of the internal detail of the map and identification of artifacts.

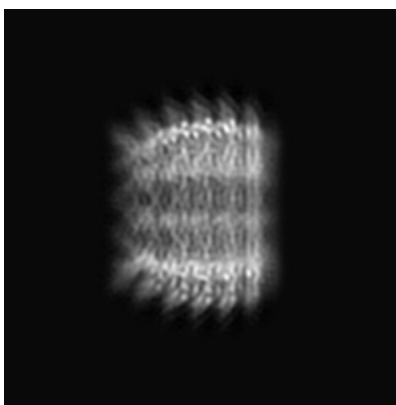
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

### 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

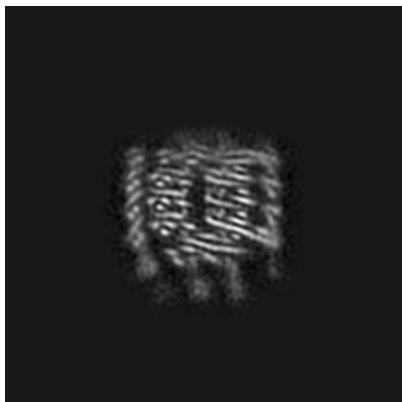


Z Index: 168

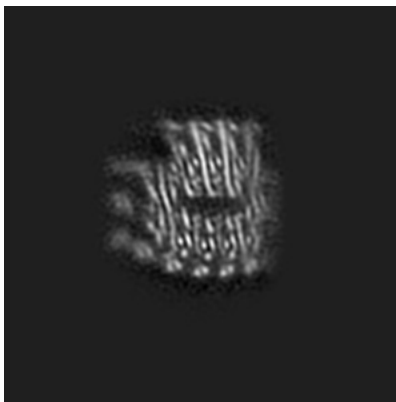
The images above show central slices of the map in three orthogonal directions.

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

### 6.3.1 Primary map



X Index: 232



Y Index: 105

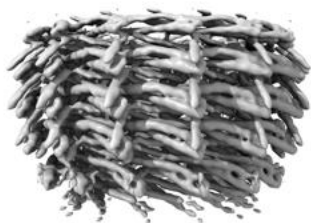


Z Index: 171

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

## 6.4 Orthogonal surface views [i](#)

### 6.4.1 Primary map



X



Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.0187. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

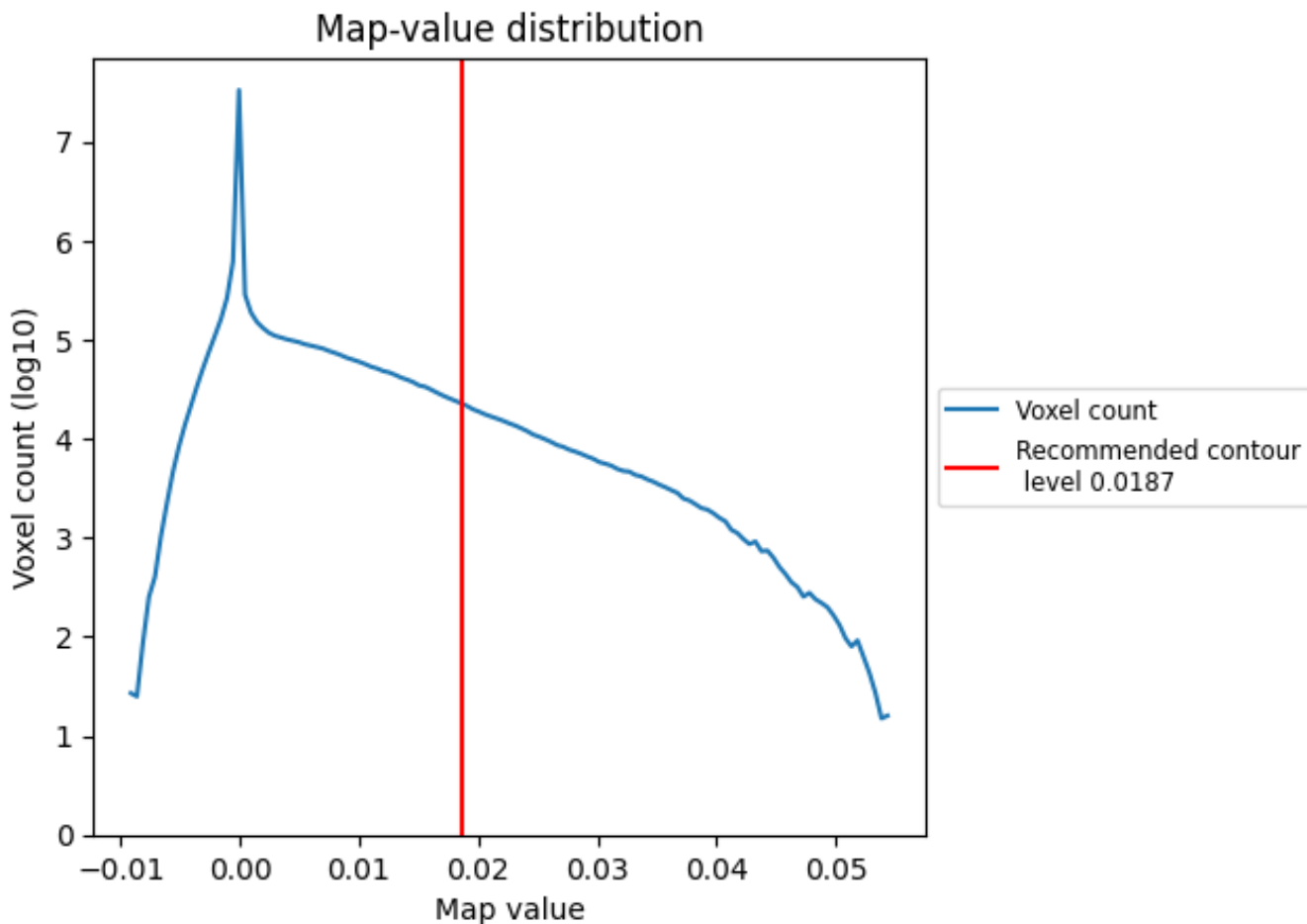
## 6.5 Mask visualisation

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

## 7 Map analysis [i](#)

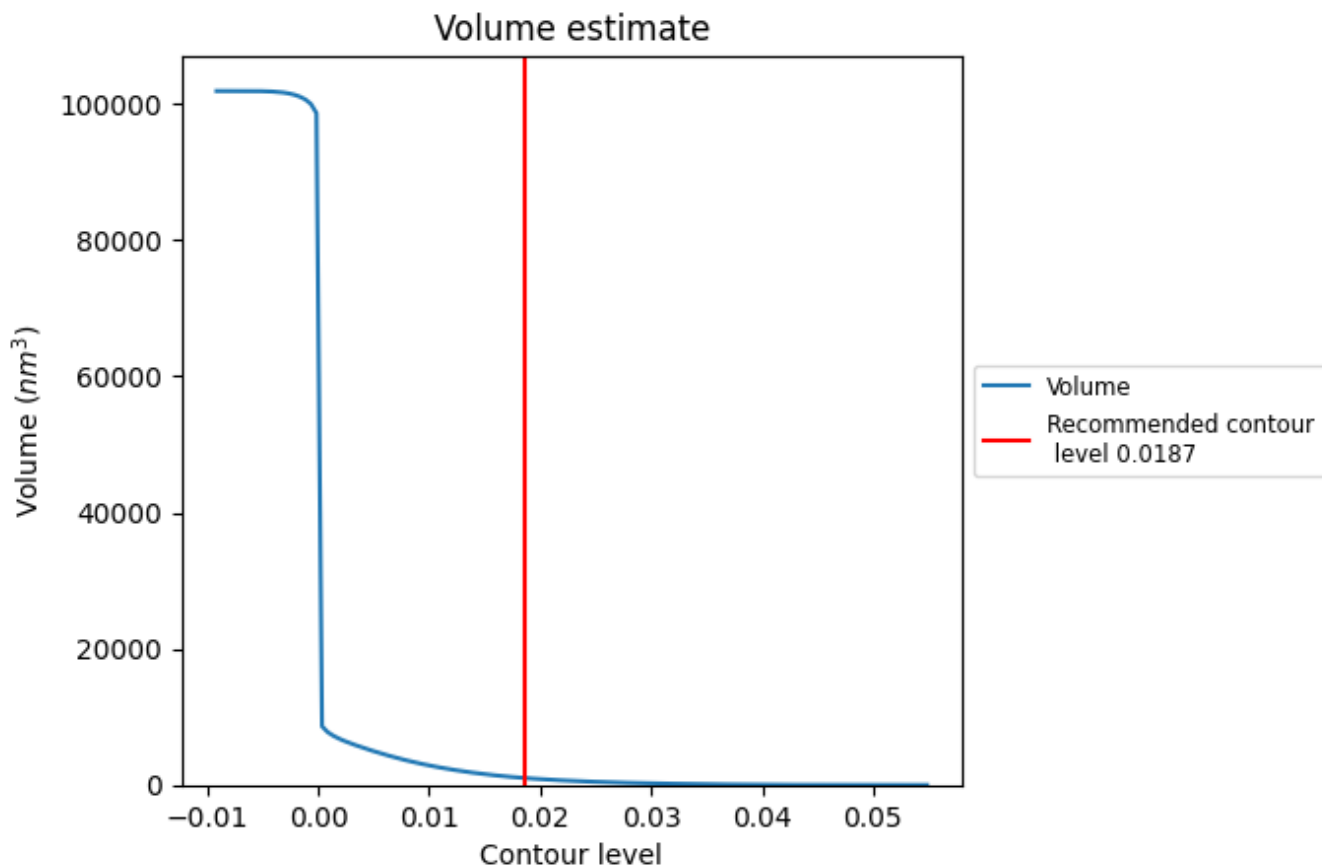
This section contains the results of statistical analysis of the map.

### 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.

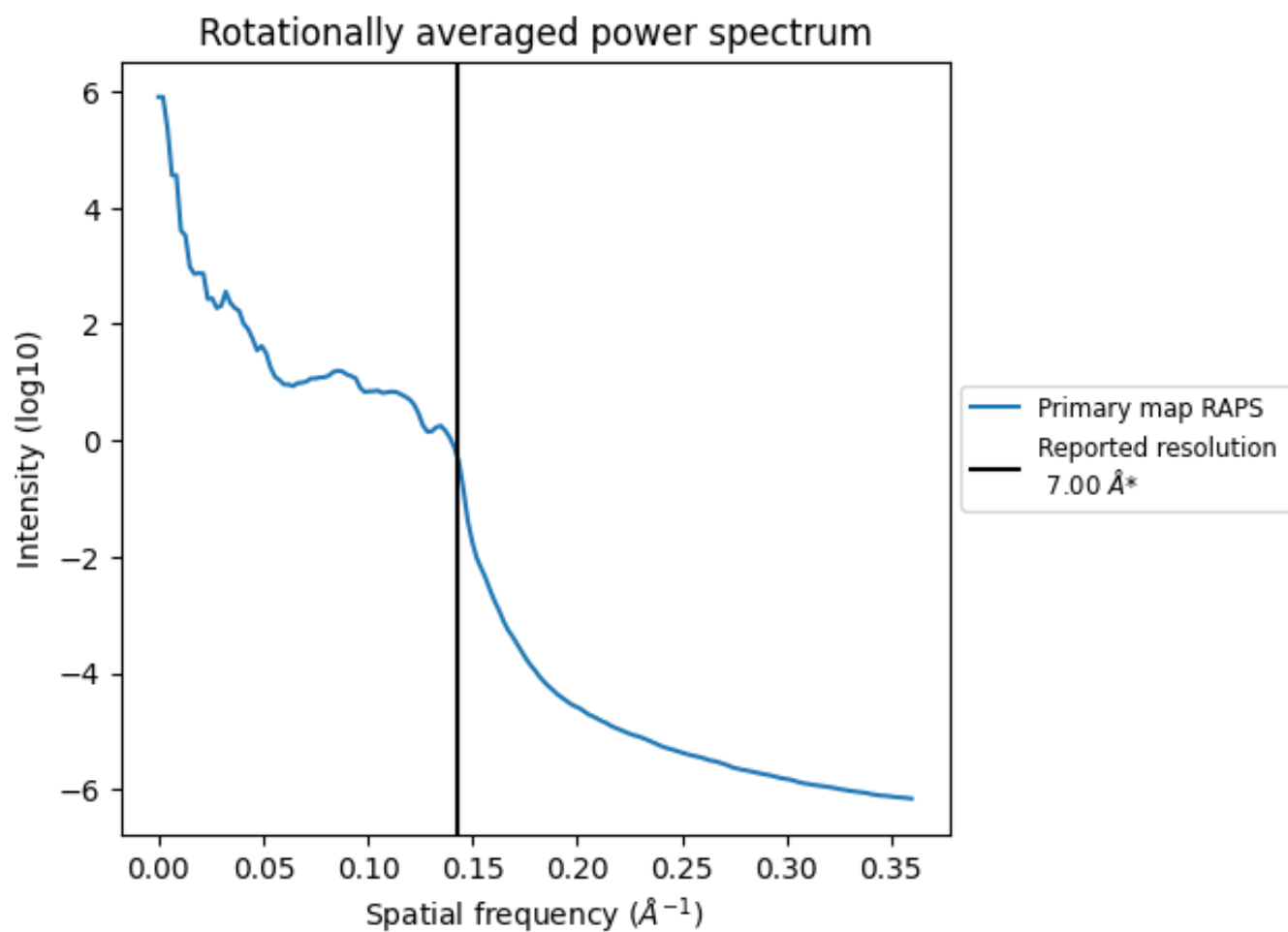
## 7.2 Volume estimate [\(i\)](#)



The volume at the recommended contour level is 1006  $\text{nm}^3$ ; this corresponds to an approximate mass of 909 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 0.143 Å<sup>-1</sup>

## 8 Fourier-Shell correlation

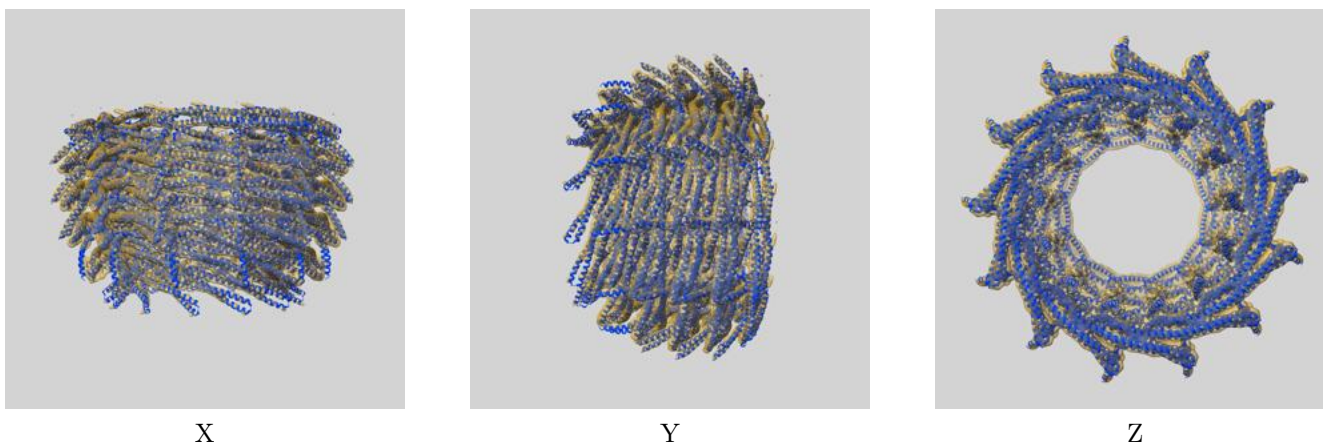
This section was not generated. No FSC curve or half-maps provided.



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

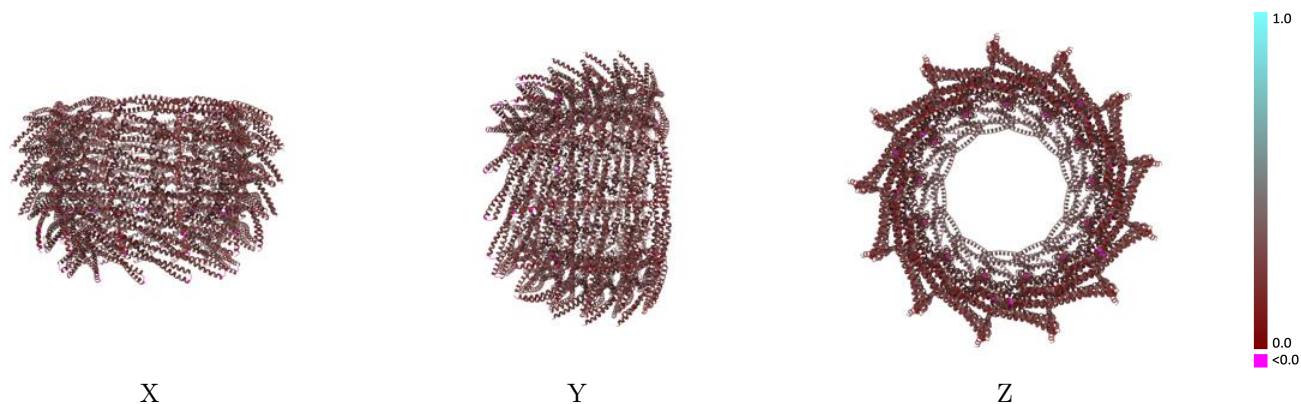
This section contains information regarding the fit between EMDB map EMD-11470 and PDB model 6ZVT. Per-residue inclusion information can be found in section 3 on page 10.

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



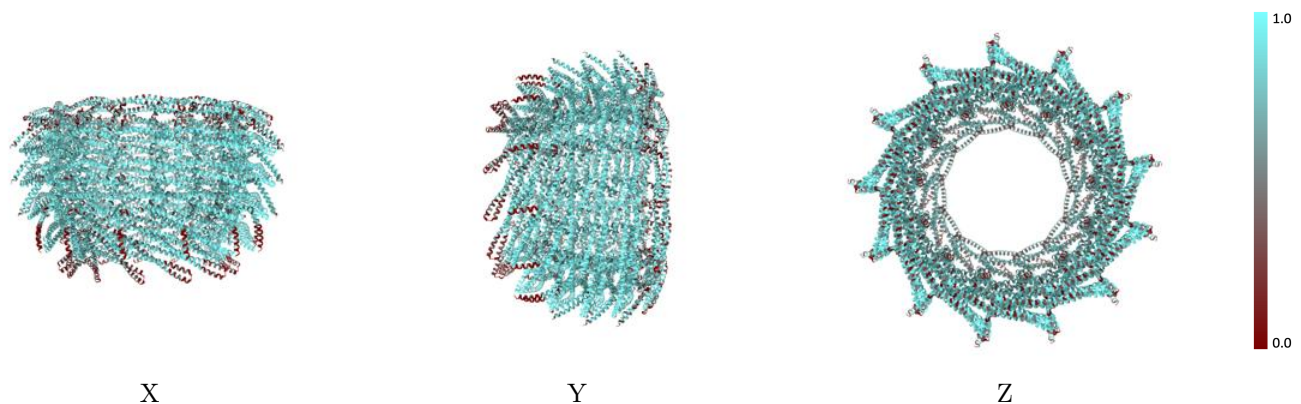
The images above show the 3D surface view of the map at the recommended contour level 0.0187 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



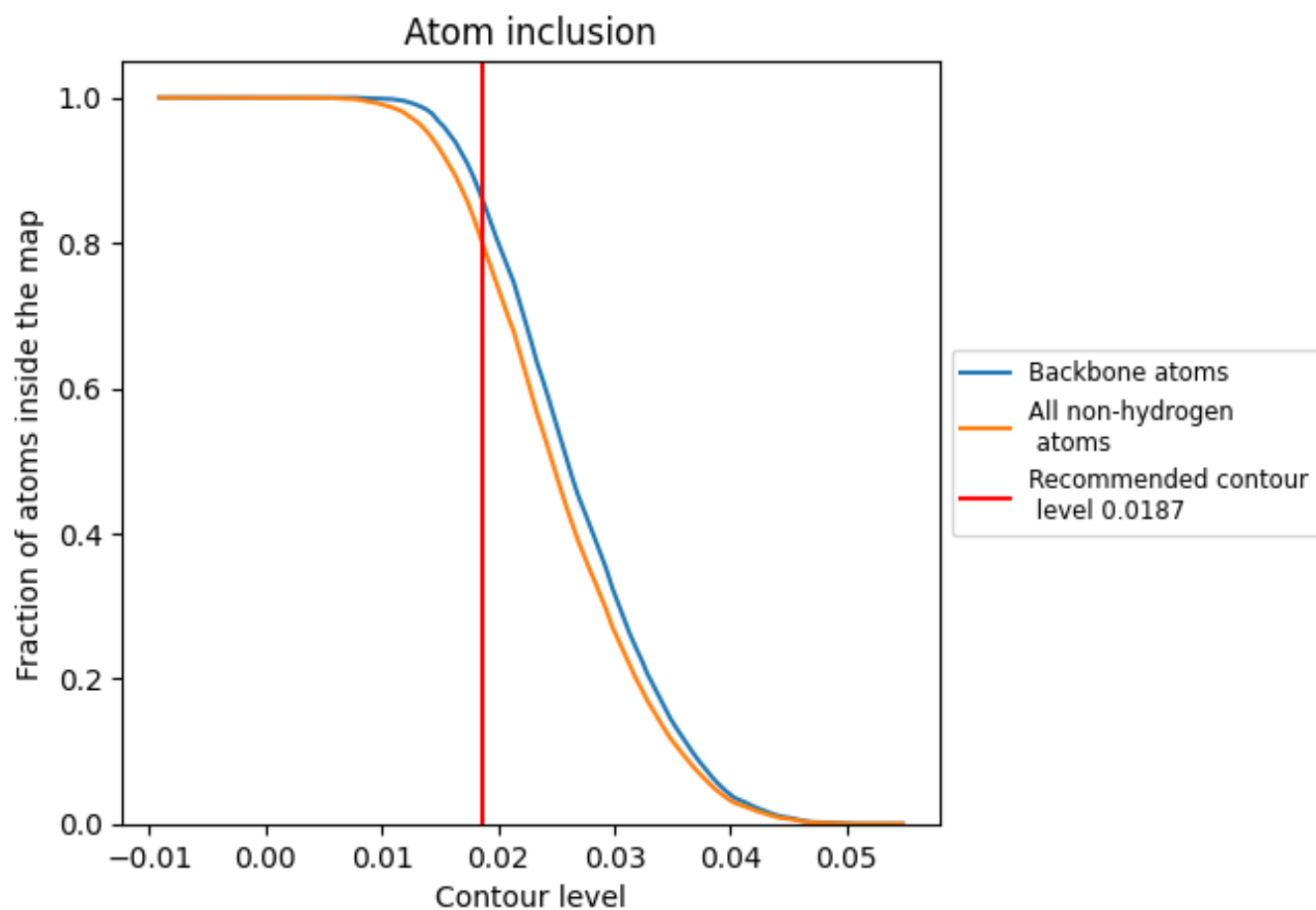
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.0187).







































































## 9.4 Atom inclusion [i](#)



At the recommended contour level, 86% of all backbone atoms, 80% 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 (0.0187) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.7981	 0.2580
A	 0.6959	 0.2520
AA	 0.7818	 0.2520
AB	 0.8840	 0.2690
AC	 0.6662	 0.2560
B	 0.7845	 0.2520
BA	 0.8840	 0.2620
BB	 0.7956	 0.2550
C	 0.8877	 0.2640
CA	 0.8831	 0.2660
CB	 0.6713	 0.2560
D	 0.8867	 0.2690
DA	 0.7974	 0.2560
DB	 0.6899	 0.2530
E	 0.8011	 0.2560
EA	 0.6701	 0.2560
EB	 0.7799	 0.2550
F	 0.6688	 0.2560
FA	 0.6929	 0.2550
FB	 0.8831	 0.2630
G	 0.6959	 0.2540
GA	 0.7790	 0.2530
GB	 0.8831	 0.2680
H	 0.7845	 0.2520
HA	 0.8849	 0.2640
HB	 0.7965	 0.2550
I	 0.8840	 0.2630
IA	 0.8849	 0.2670
IB	 0.6701	 0.2560
J	 0.8840	 0.2650
JA	 0.7956	 0.2580
JB	 0.6944	 0.2520
K	 0.7956	 0.2550
KA	 0.6688	 0.2540
KB	 0.7808	 0.2500







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Chain	Atom inclusion	Q-score
L	0.6726	0.2540
LA	0.6929	0.2550
LB	0.8849	0.2630
M	0.6944	0.2530
MA	0.7836	0.2540
MB	0.8821	0.2660
N	0.7818	0.2500
NA	0.8840	0.2650
NB	0.7956	0.2540
O	0.8858	0.2620
OA	0.8840	0.2700
OB	0.6726	0.2560
P	0.8803	0.2680
PA	0.8011	0.2570
PB	0.6959	0.2520
Q	0.7974	0.2570
QA	0.6713	0.2560
QB	0.7799	0.2520
R	0.6726	0.2550
RA	0.6959	0.2510
RB	0.8840	0.2640
S	0.6929	0.2540
SA	0.7864	0.2540
SB	0.8831	0.2670
T	0.7808	0.2520
TA	0.8849	0.2650
TB	0.7965	0.2550
UA	0.8849	0.2690
UB	0.6688	0.2550
V	0.8821	0.2640
VA	0.8011	0.2570
VB	0.6914	0.2530
W	0.8821	0.2660
WA	0.6726	0.2550
WB	0.7799	0.2520
X	0.7993	0.2570
XA	0.6944	0.2540
XB	0.8849	0.2620
Y	0.6752	0.2550
YA	0.7790	0.2540
YB	0.8831	0.2660
Z	0.6974	0.2530

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Chain	Atom inclusion	Q-score
ZA	 0.8831	 0.2660
ZB	 0.8002	 0.2550