



Full wwPDB EM Validation Report ⓘ

Mar 8, 2026 – 02:16 PM UTC

PDB ID : 6E8G / pdb_00006e8g
EMDB ID : EMD-9005
Title : CryoEM reconstruction of IST1-CHMP1B copolymer filament bound to ss-DNA at 2.9 Angstrom resolution
Authors : Talledge, N.; Frost, A.; McCullough, J.
Deposited on : 2018-07-29
Resolution : 2.90 Å(reported)

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

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

A user guide is available at

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

The types of validation reports are described at

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

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

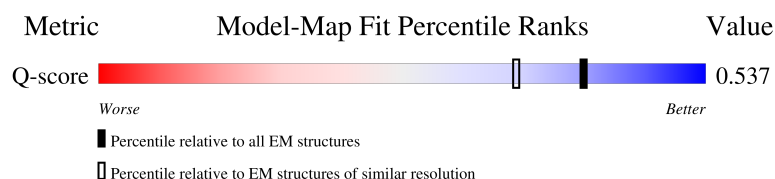
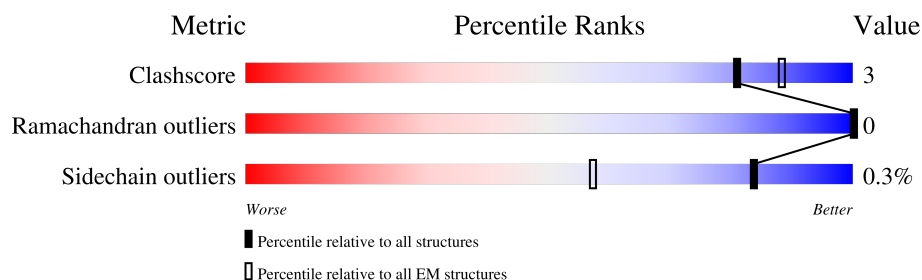
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDb archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh &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

1 Overall quality at a glance

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

The reported resolution of this entry is 2.90 Å.

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



Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	13054 (2.40 - 3.40)

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

Mol	Chain	Length	Quality of chain
1	A	366	7% 43% 54%
1	BA	366	6% 43% 54%
1	BB	366	7% 43% 54%
1	C	366	6% 43% 54%

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Mol	Chain	Length	Quality of chain		
1	DA	366	7%	43%	54%
1	DB	366	7%	43%	54%
1	E	366	7%	43%	54%
1	FA	366	7%	43%	54%
1	FB	366	8%	43%	54%
1	G	366	7%	44%	54%
1	HA	366	7%	43%	54%
1	HB	366	6%	43%	54%
1	I	366	7%	43%	54%
1	JA	366	7%	43%	54%
1	JB	366	7%	44%	54%
1	K	366	7%	43%	54%
1	LA	366	6%	43%	54%
1	LB	366	6%	43%	54%
1	M	366	8%	43%	54%
1	NA	366	7%	43%	54%
1	NB	366	7%	43%	54%
1	O	366	7%	43%	54%
1	PA	366	7%	43%	54%
1	PB	366	8%	43%	54%
1	Q	366	7%	43%	54%
1	RA	366	7%	44%	54%
1	RB	366	7%	43%	54%
1	S	366	7%	43%	54%
1	TA	366	7%	43%	54%

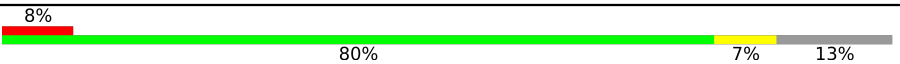
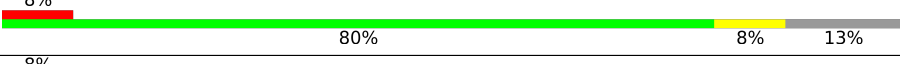
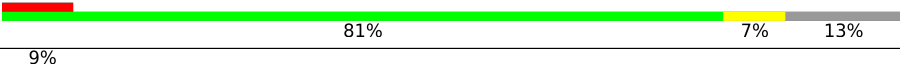
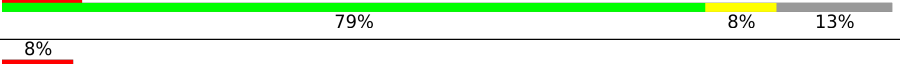
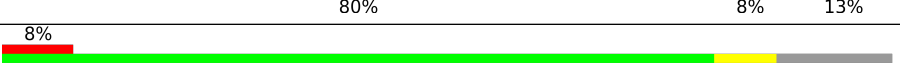
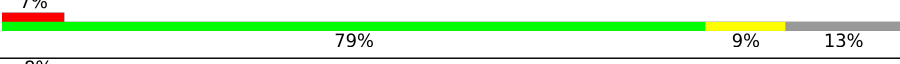
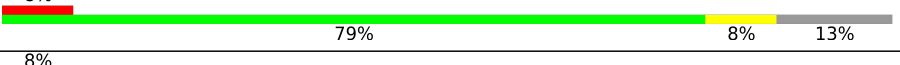
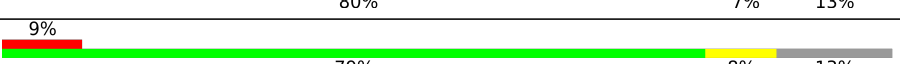
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Mol	Chain	Length	Quality of chain
1	TB	366	
1	V	366	
1	VA	366	
1	X	366	
1	XA	366	
1	Z	366	
1	ZA	366	
2	AA	199	
2	AB	199	
2	B	199	
2	CA	199	
2	CB	199	
2	D	199	
2	EA	199	
2	EB	199	
2	F	199	
2	GA	199	
2	GB	199	
2	H	199	
2	IA	199	
2	IB	199	
2	J	199	
2	KA	199	
2	KB	199	
2	L	199	

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Mol	Chain	Length	Quality of chain
2	MA	199	
2	MB	199	
2	N	199	
2	OA	199	
2	OB	199	
2	P	199	
2	QA	199	
2	QB	199	
2	R	199	
2	SA	199	
2	SB	199	
2	T	199	
2	UA	199	
2	UB	199	
2	W	199	
2	WA	199	
2	Y	199	
2	YA	199	

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 97272 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 IST1 homolog.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	S	167	Total	C	N	O	S	0	0
			1349	862	234	248	5		
1	A	167	Total	C	N	O	S	0	0
			1349	862	234	248	5		
1	C	167	Total	C	N	O	S	0	0
			1349	862	234	248	5		
1	E	167	Total	C	N	O	S	0	0
			1349	862	234	248	5		
1	G	167	Total	C	N	O	S	0	0
			1349	862	234	248	5		
1	I	167	Total	C	N	O	S	0	0
			1349	862	234	248	5		
1	K	167	Total	C	N	O	S	0	0
			1349	862	234	248	5		
1	M	167	Total	C	N	O	S	0	0
			1349	862	234	248	5		
1	O	167	Total	C	N	O	S	0	0
			1349	862	234	248	5		
1	Q	167	Total	C	N	O	S	0	0
			1349	862	234	248	5		
1	V	167	Total	C	N	O	S	0	0
			1349	862	234	248	5		
1	X	167	Total	C	N	O	S	0	0
			1349	862	234	248	5		
1	Z	167	Total	C	N	O	S	0	0
			1349	862	234	248	5		
1	BA	167	Total	C	N	O	S	0	0
			1349	862	234	248	5		
1	DA	167	Total	C	N	O	S	0	0
			1349	862	234	248	5		
1	FA	167	Total	C	N	O	S	0	0
			1349	862	234	248	5		
1	HA	167	Total	C	N	O	S	0	0
			1349	862	234	248	5		

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Mol	Chain	Residues	Atoms					AltConf	Trace
1	JA	167	Total	C	N	O	S	0	0
			1349	862	234	248	5		
1	LA	167	Total	C	N	O	S	0	0
			1349	862	234	248	5		
1	NA	167	Total	C	N	O	S	0	0
			1349	862	234	248	5		
1	PA	167	Total	C	N	O	S	0	0
			1349	862	234	248	5		
1	RA	167	Total	C	N	O	S	0	0
			1349	862	234	248	5		
1	TA	167	Total	C	N	O	S	0	0
			1349	862	234	248	5		
1	VA	167	Total	C	N	O	S	0	0
			1349	862	234	248	5		
1	XA	167	Total	C	N	O	S	0	0
			1349	862	234	248	5		
1	ZA	167	Total	C	N	O	S	0	0
			1349	862	234	248	5		
1	BB	167	Total	C	N	O	S	0	0
			1349	862	234	248	5		
1	DB	167	Total	C	N	O	S	0	0
			1349	862	234	248	5		
1	FB	167	Total	C	N	O	S	0	0
			1349	862	234	248	5		
1	HB	167	Total	C	N	O	S	0	0
			1349	862	234	248	5		
1	JB	167	Total	C	N	O	S	0	0
			1349	862	234	248	5		
1	LB	167	Total	C	N	O	S	0	0
			1349	862	234	248	5		
1	NB	167	Total	C	N	O	S	0	0
			1349	862	234	248	5		
1	PB	167	Total	C	N	O	S	0	0
			1349	862	234	248	5		
1	RB	167	Total	C	N	O	S	0	0
			1349	862	234	248	5		
1	TB	167	Total	C	N	O	S	0	0
			1349	862	234	248	5		

- Molecule 2 is a protein called Charged multivesicular body protein 1b.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	T	174	Total	C	N	O	S	0	0
			1353	831	245	264	13		
2	B	174	Total	C	N	O	S	0	0
			1353	831	245	264	13		
2	D	174	Total	C	N	O	S	0	0
			1353	831	245	264	13		
2	F	174	Total	C	N	O	S	0	0
			1353	831	245	264	13		
2	H	174	Total	C	N	O	S	0	0
			1353	831	245	264	13		
2	J	174	Total	C	N	O	S	0	0
			1353	831	245	264	13		
2	L	174	Total	C	N	O	S	0	0
			1353	831	245	264	13		
2	N	174	Total	C	N	O	S	0	0
			1353	831	245	264	13		
2	P	174	Total	C	N	O	S	0	0
			1353	831	245	264	13		
2	R	174	Total	C	N	O	S	0	0
			1353	831	245	264	13		
2	W	174	Total	C	N	O	S	0	0
			1353	831	245	264	13		
2	Y	174	Total	C	N	O	S	0	0
			1353	831	245	264	13		
2	AA	174	Total	C	N	O	S	0	0
			1353	831	245	264	13		
2	CA	174	Total	C	N	O	S	0	0
			1353	831	245	264	13		
2	EA	174	Total	C	N	O	S	0	0
			1353	831	245	264	13		
2	GA	174	Total	C	N	O	S	0	0
			1353	831	245	264	13		
2	IA	174	Total	C	N	O	S	0	0
			1353	831	245	264	13		
2	KA	174	Total	C	N	O	S	0	0
			1353	831	245	264	13		
2	MA	174	Total	C	N	O	S	0	0
			1353	831	245	264	13		
2	OA	174	Total	C	N	O	S	0	0
			1353	831	245	264	13		
2	QA	174	Total	C	N	O	S	0	0
			1353	831	245	264	13		
2	SA	174	Total	C	N	O	S	0	0
			1353	831	245	264	13		

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Mol	Chain	Residues	Atoms					AltConf	Trace
2	UA	174	Total 1353	C 831	N 245	O 264	S 13	0	0
2	WA	174	Total 1353	C 831	N 245	O 264	S 13	0	0
2	YA	174	Total 1353	C 831	N 245	O 264	S 13	0	0
2	AB	174	Total 1353	C 831	N 245	O 264	S 13	0	0
2	CB	174	Total 1353	C 831	N 245	O 264	S 13	0	0
2	EB	174	Total 1353	C 831	N 245	O 264	S 13	0	0
2	GB	174	Total 1353	C 831	N 245	O 264	S 13	0	0
2	IB	174	Total 1353	C 831	N 245	O 264	S 13	0	0
2	KB	174	Total 1353	C 831	N 245	O 264	S 13	0	0
2	MB	174	Total 1353	C 831	N 245	O 264	S 13	0	0
2	OB	174	Total 1353	C 831	N 245	O 264	S 13	0	0
2	QB	174	Total 1353	C 831	N 245	O 264	S 13	0	0
2	SB	174	Total 1353	C 831	N 245	O 264	S 13	0	0
2	UB	174	Total 1353	C 831	N 245	O 264	S 13	0	0

There are 36 discrepancies between the modelled and reference sequences:

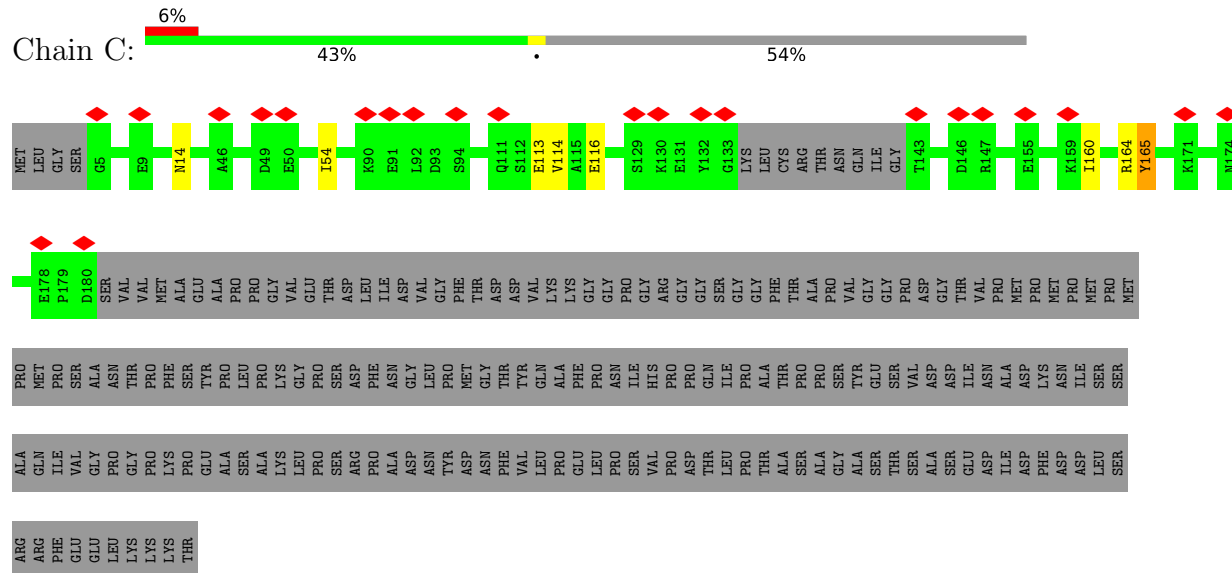
Chain	Residue	Modelled	Actual	Comment	Reference
T	37	GLU	LYS	engineered mutation	UNP Q7LBR1
B	37	GLU	LYS	engineered mutation	UNP Q7LBR1
D	37	GLU	LYS	engineered mutation	UNP Q7LBR1
F	37	GLU	LYS	engineered mutation	UNP Q7LBR1
H	37	GLU	LYS	engineered mutation	UNP Q7LBR1
J	37	GLU	LYS	engineered mutation	UNP Q7LBR1
L	37	GLU	LYS	engineered mutation	UNP Q7LBR1
N	37	GLU	LYS	engineered mutation	UNP Q7LBR1
P	37	GLU	LYS	engineered mutation	UNP Q7LBR1
R	37	GLU	LYS	engineered mutation	UNP Q7LBR1
W	37	GLU	LYS	engineered mutation	UNP Q7LBR1

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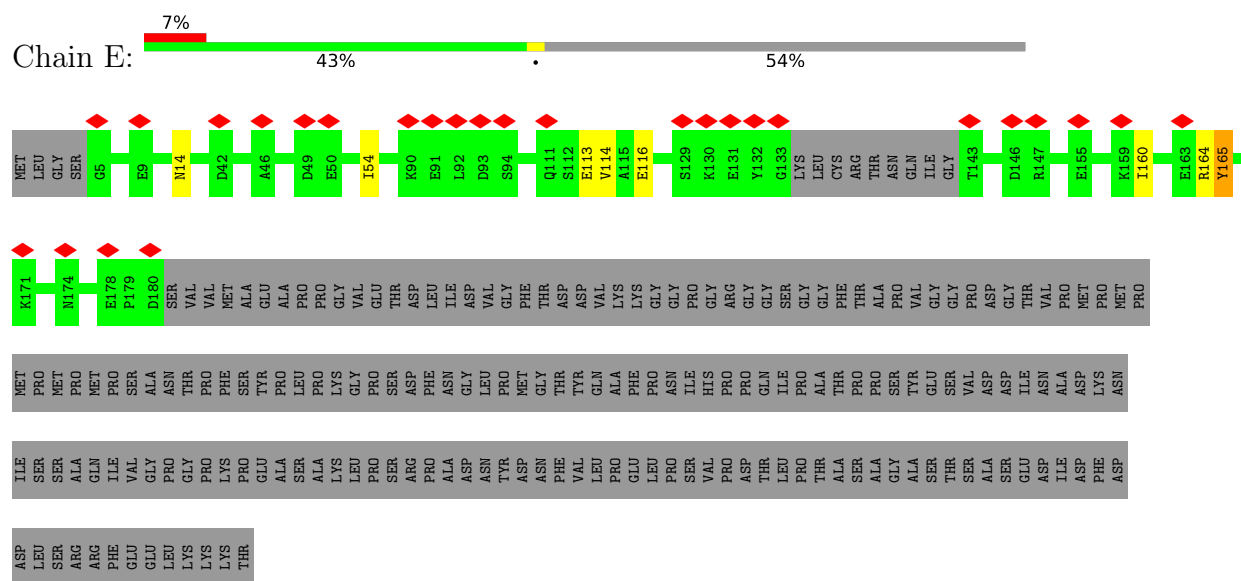
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Chain	Residue	Modelled	Actual	Comment	Reference
Y	37	GLU	LYS	engineered mutation	UNP Q7LBR1
AA	37	GLU	LYS	engineered mutation	UNP Q7LBR1
CA	37	GLU	LYS	engineered mutation	UNP Q7LBR1
EA	37	GLU	LYS	engineered mutation	UNP Q7LBR1
GA	37	GLU	LYS	engineered mutation	UNP Q7LBR1
IA	37	GLU	LYS	engineered mutation	UNP Q7LBR1
KA	37	GLU	LYS	engineered mutation	UNP Q7LBR1
MA	37	GLU	LYS	engineered mutation	UNP Q7LBR1
OA	37	GLU	LYS	engineered mutation	UNP Q7LBR1
QA	37	GLU	LYS	engineered mutation	UNP Q7LBR1
SA	37	GLU	LYS	engineered mutation	UNP Q7LBR1
UA	37	GLU	LYS	engineered mutation	UNP Q7LBR1
WA	37	GLU	LYS	engineered mutation	UNP Q7LBR1
YA	37	GLU	LYS	engineered mutation	UNP Q7LBR1
AB	37	GLU	LYS	engineered mutation	UNP Q7LBR1
CB	37	GLU	LYS	engineered mutation	UNP Q7LBR1
EB	37	GLU	LYS	engineered mutation	UNP Q7LBR1
GB	37	GLU	LYS	engineered mutation	UNP Q7LBR1
IB	37	GLU	LYS	engineered mutation	UNP Q7LBR1
KB	37	GLU	LYS	engineered mutation	UNP Q7LBR1
MB	37	GLU	LYS	engineered mutation	UNP Q7LBR1
OB	37	GLU	LYS	engineered mutation	UNP Q7LBR1
QB	37	GLU	LYS	engineered mutation	UNP Q7LBR1
SB	37	GLU	LYS	engineered mutation	UNP Q7LBR1
UB	37	GLU	LYS	engineered mutation	UNP Q7LBR1

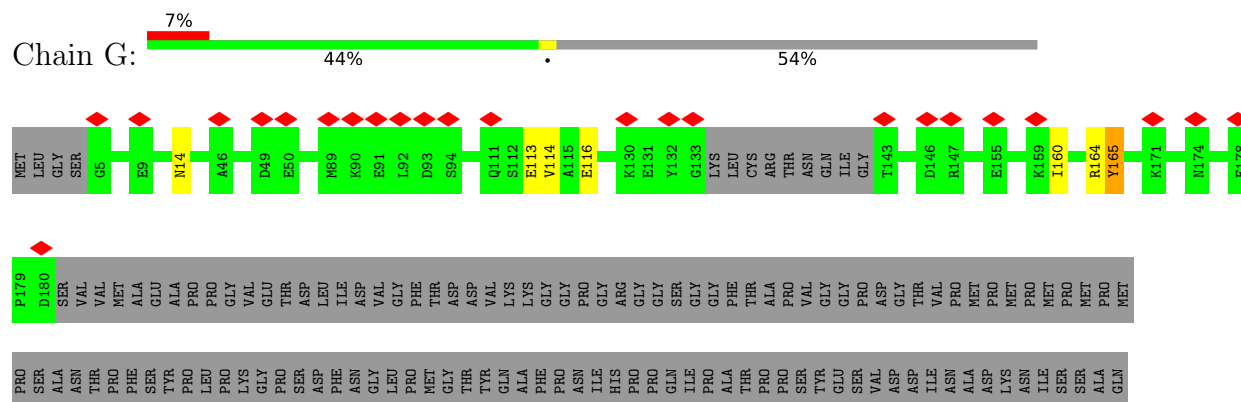
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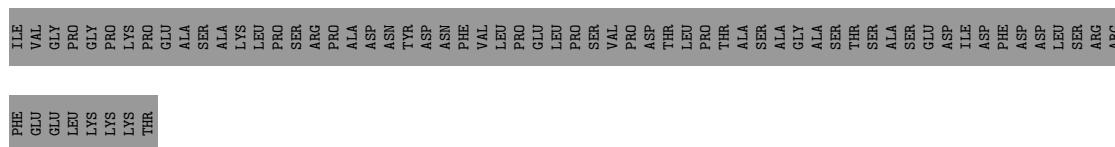


• Molecule 1: IST1 homolog

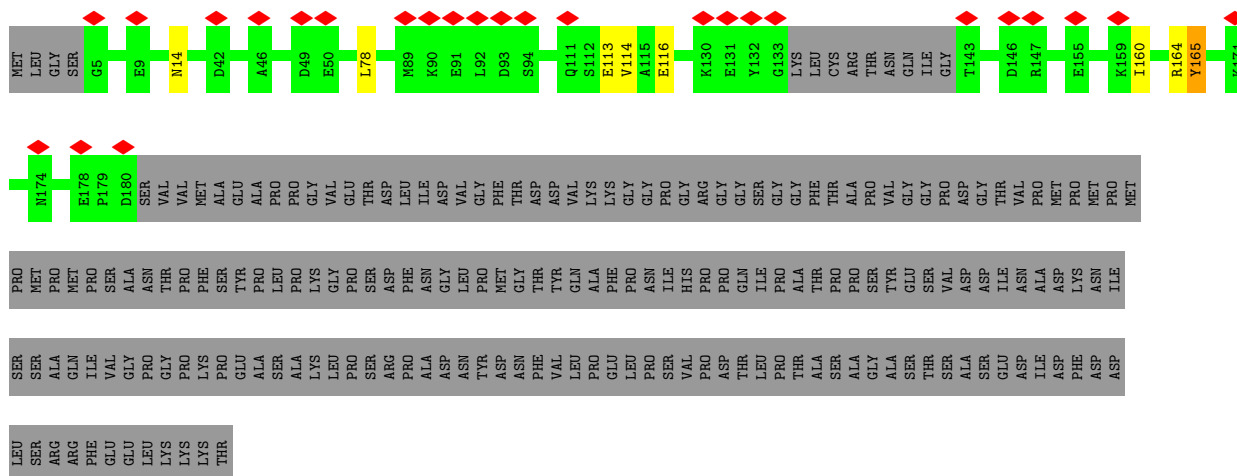


• Molecule 1: IST1 homolog

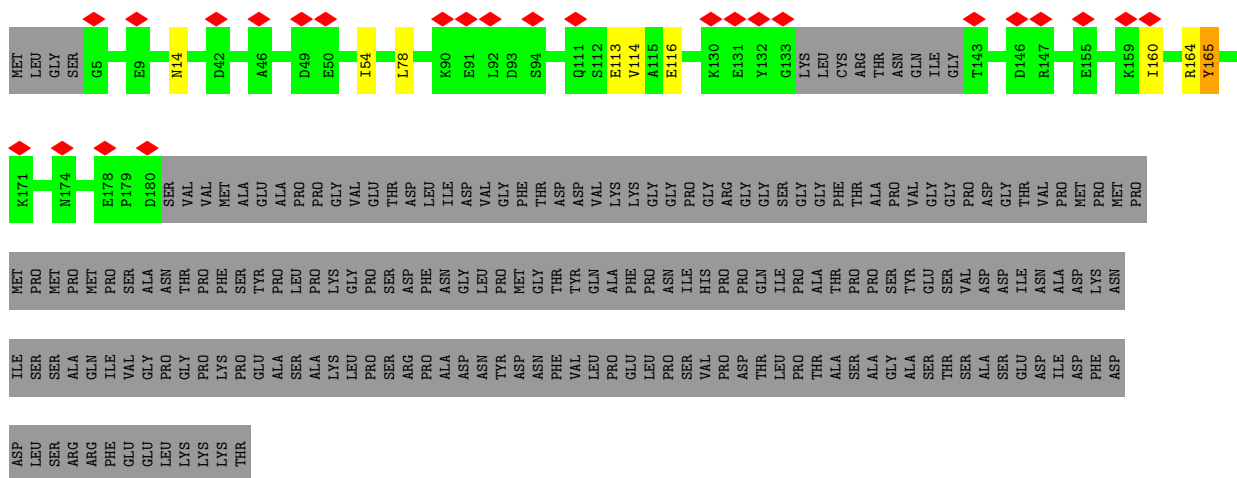
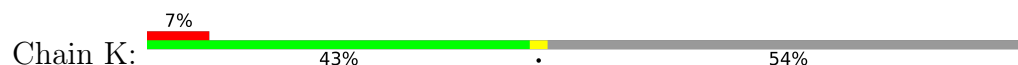




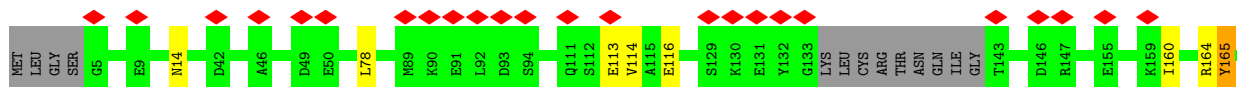
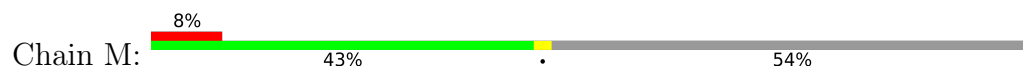
● Molecule 1: IST1 homolog

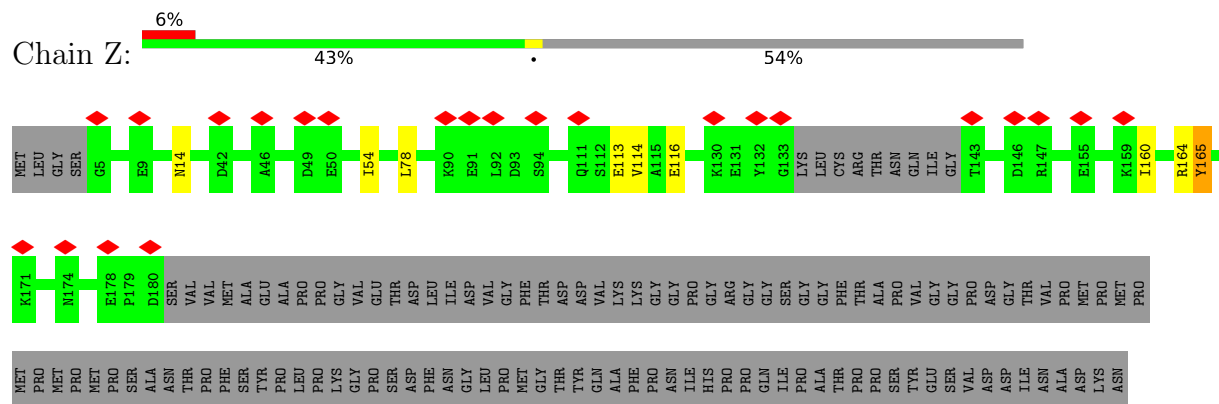


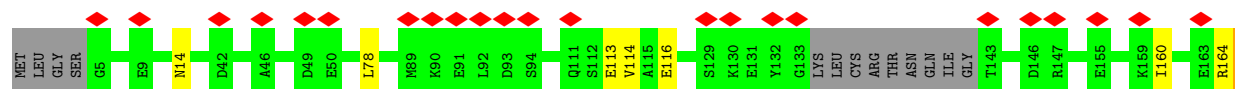
● Molecule 1: IST1 homolog

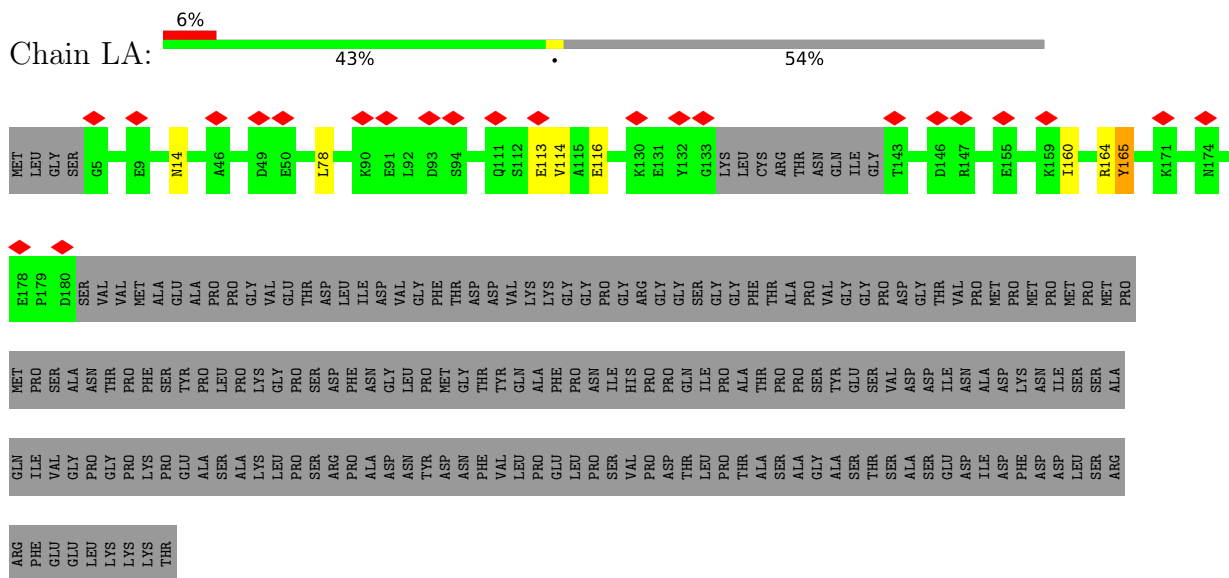


● Molecule 1: IST1 homolog

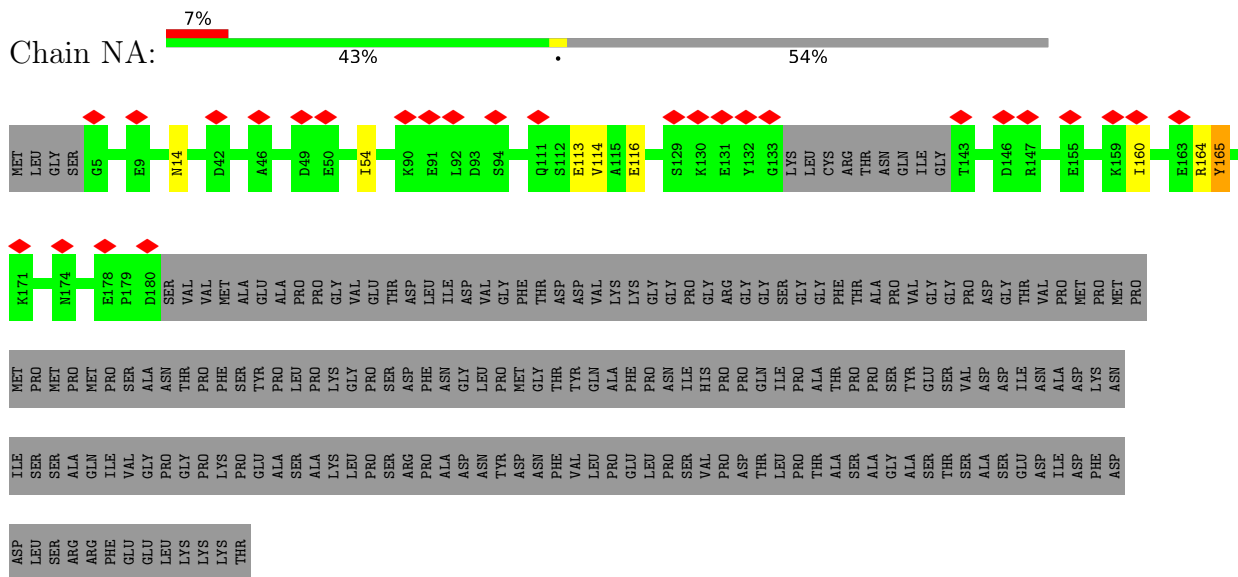




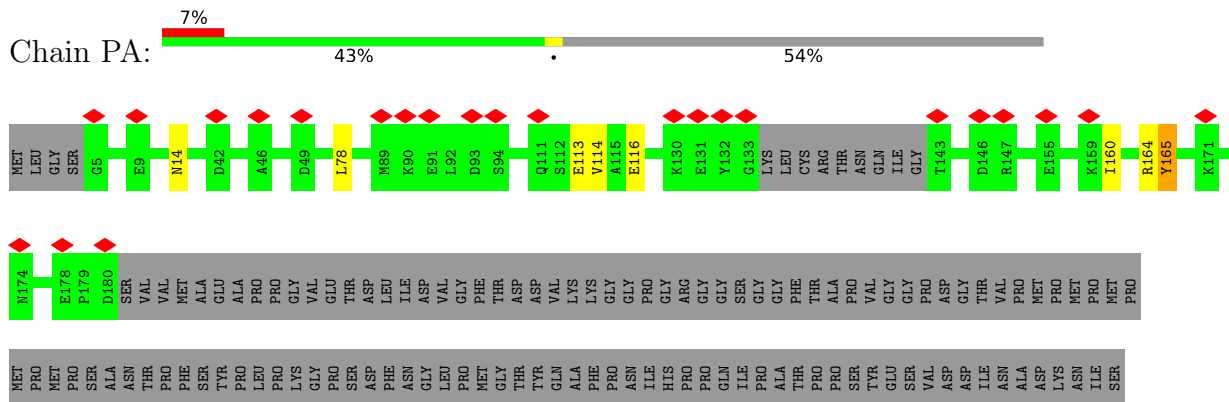




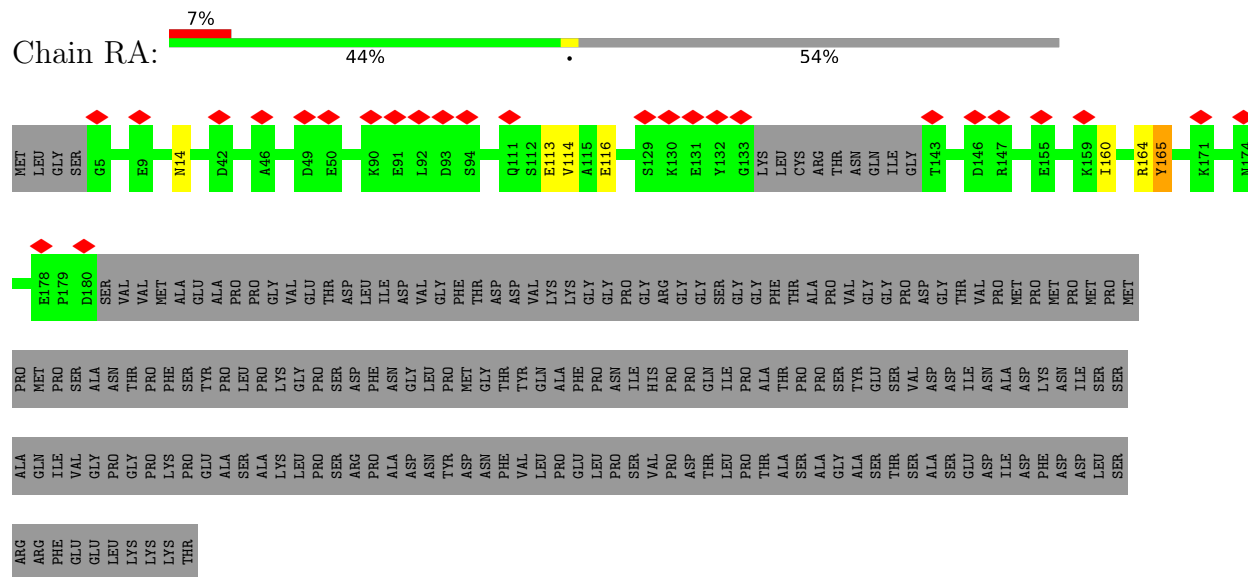
- Molecule 1: IST1 homolog



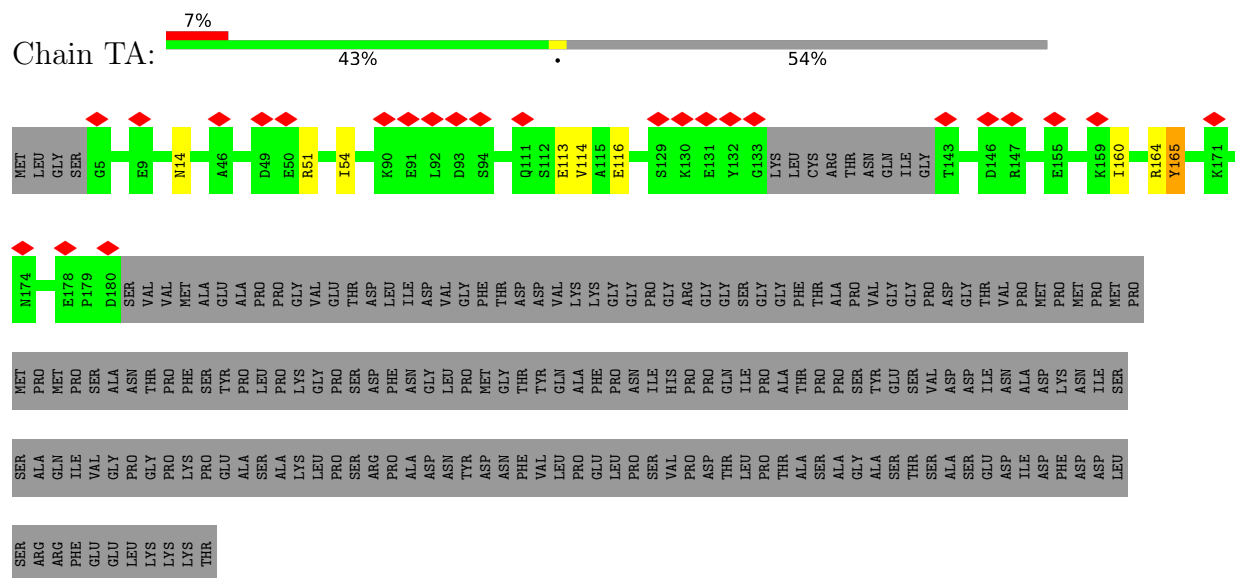
- Molecule 1: IST1 homolog



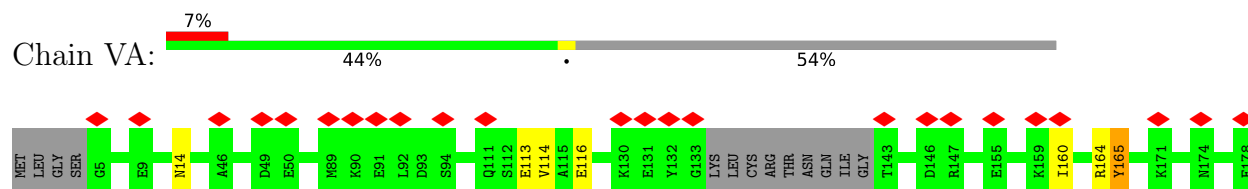
- Molecule 1: IST1 homolog

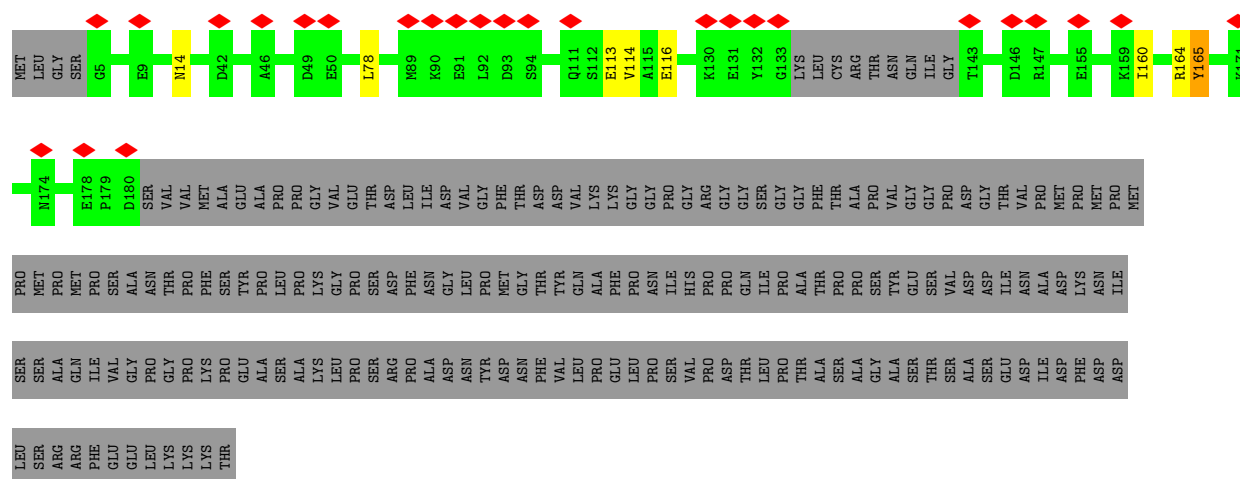
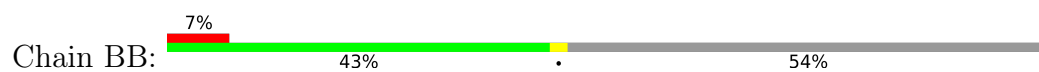


- Molecule 1: IST1 homolog

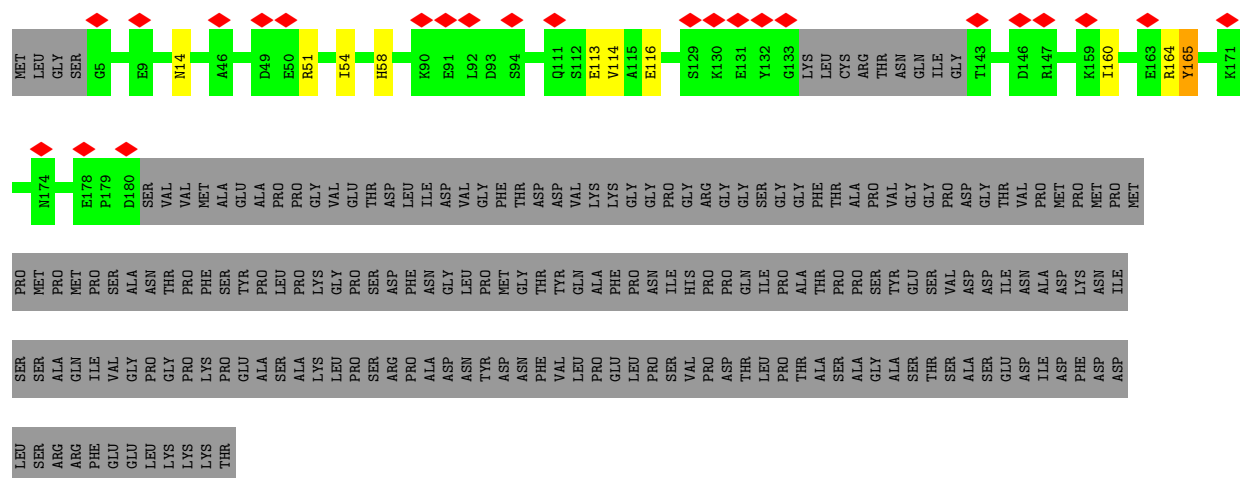


- Molecule 1: IST1 homolog

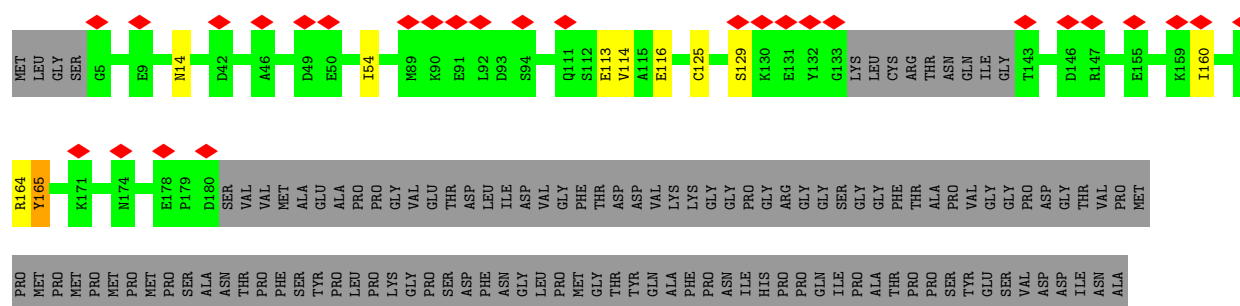
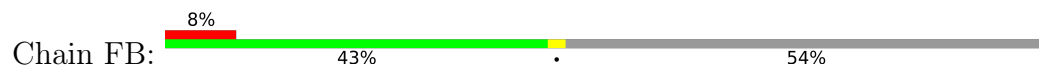


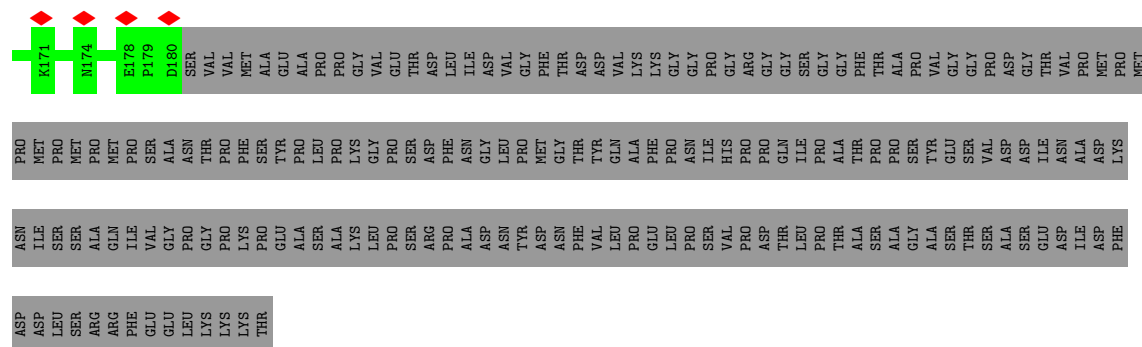


• Molecule 1: IST1 homolog

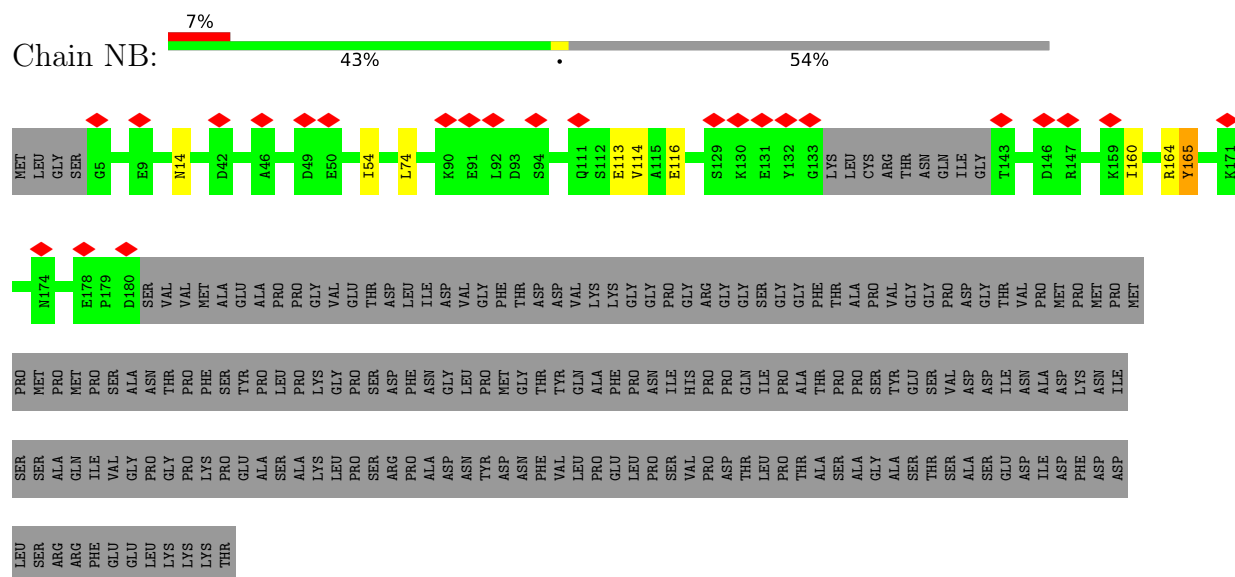


• Molecule 1: IST1 homolog

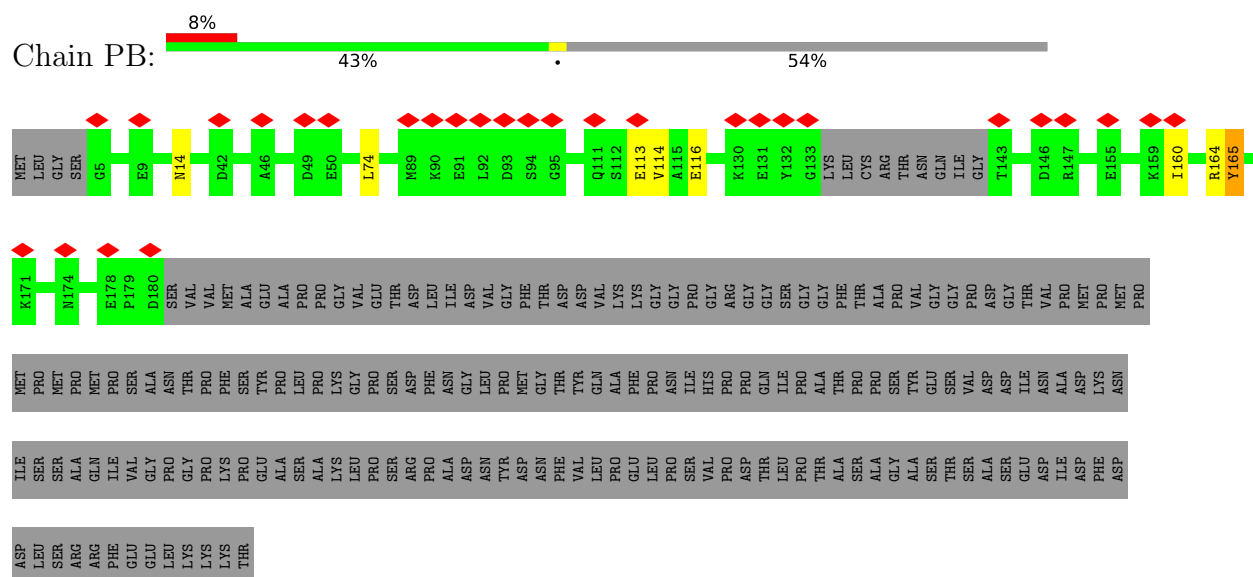




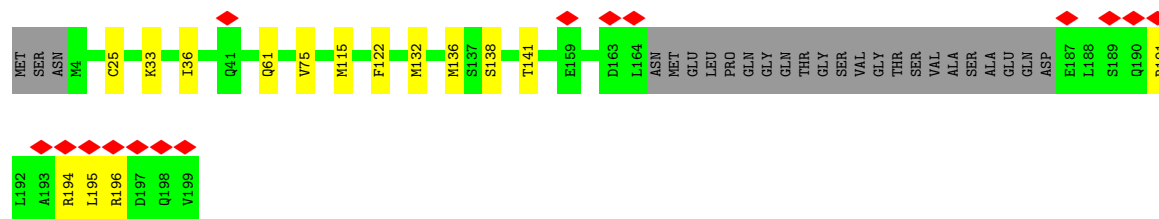
- Molecule 1: IST1 homolog



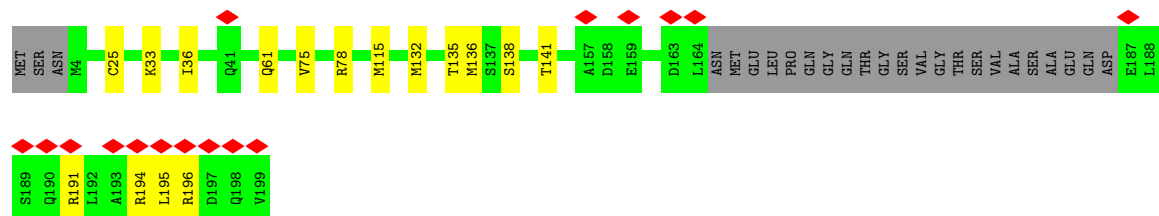
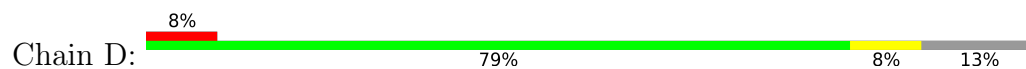
- Molecule 1: IST1 homolog



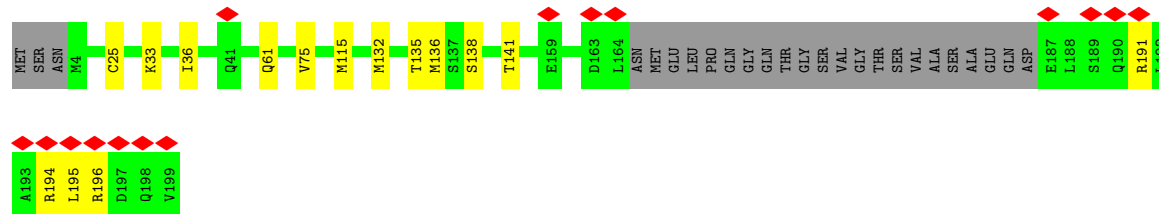
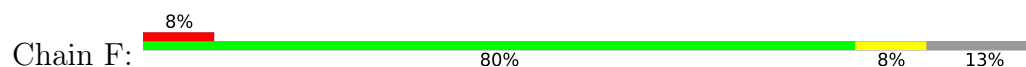
- Molecule 1: IST1 homolog



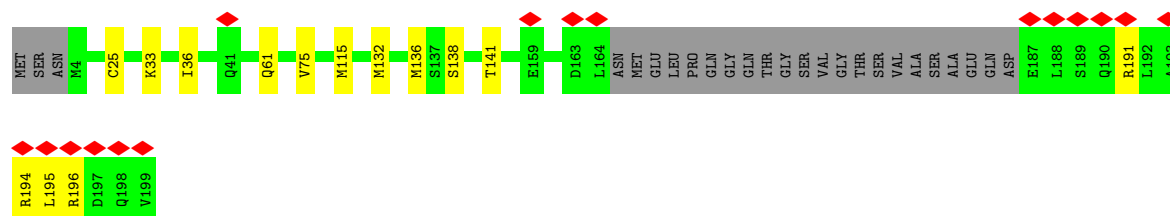
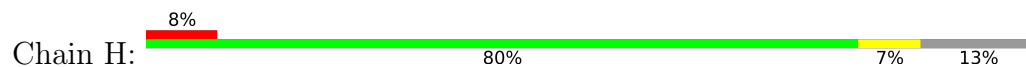
- Molecule 2: Charged multivesicular body protein 1b



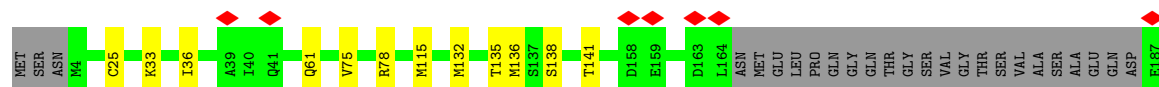
- Molecule 2: Charged multivesicular body protein 1b



- Molecule 2: Charged multivesicular body protein 1b

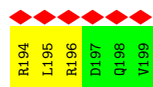
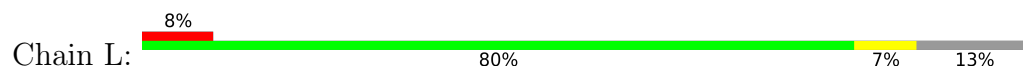


- Molecule 2: Charged multivesicular body protein 1b

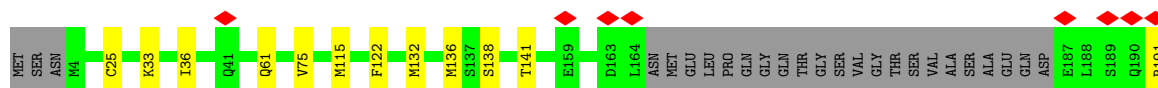
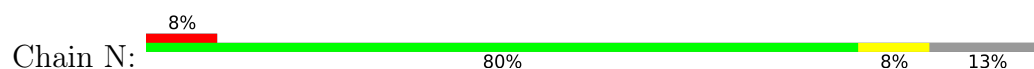




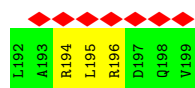
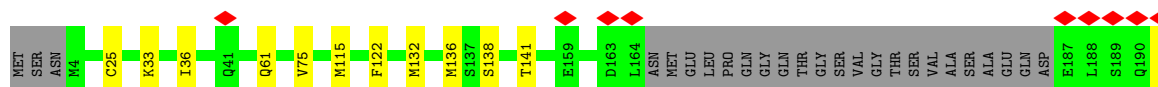
- Molecule 2: Charged multivesicular body protein 1b



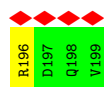
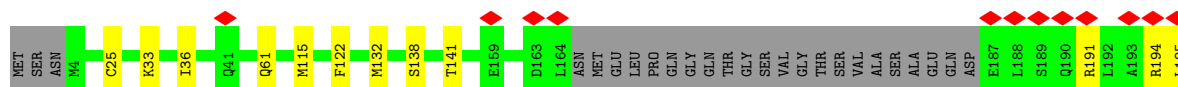
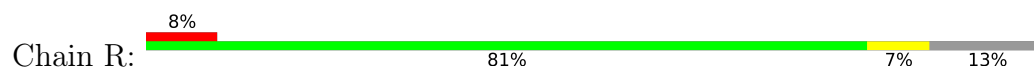
- Molecule 2: Charged multivesicular body protein 1b



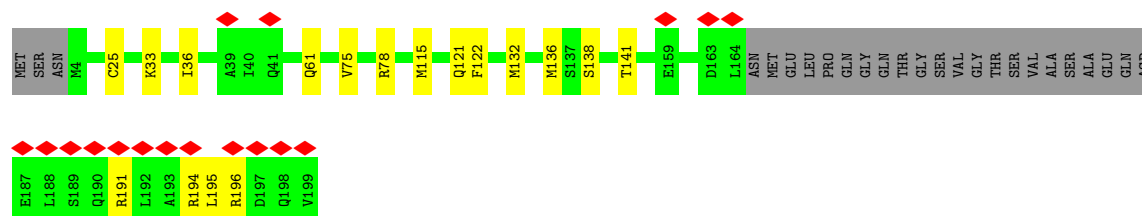
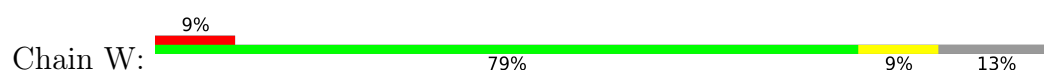
- Molecule 2: Charged multivesicular body protein 1b



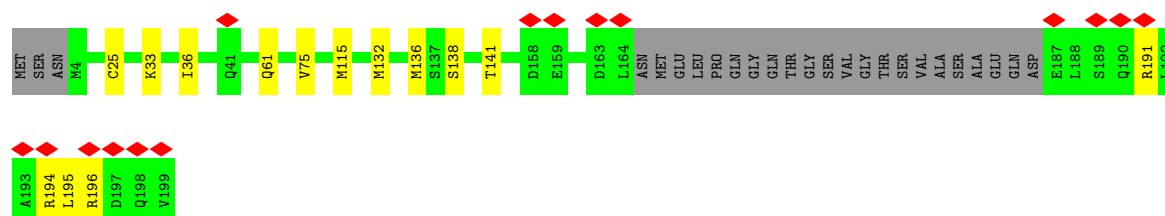
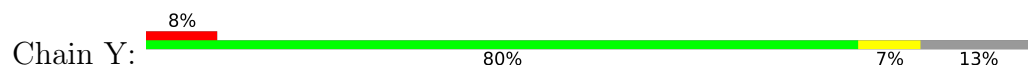
- Molecule 2: Charged multivesicular body protein 1b



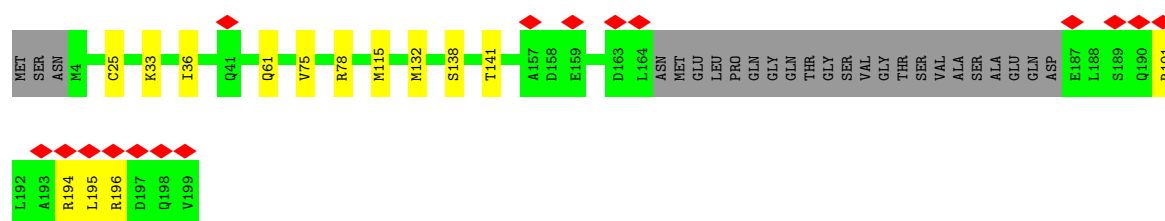
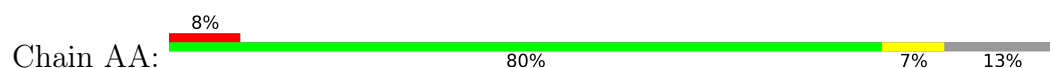
- Molecule 2: Charged multivesicular body protein 1b



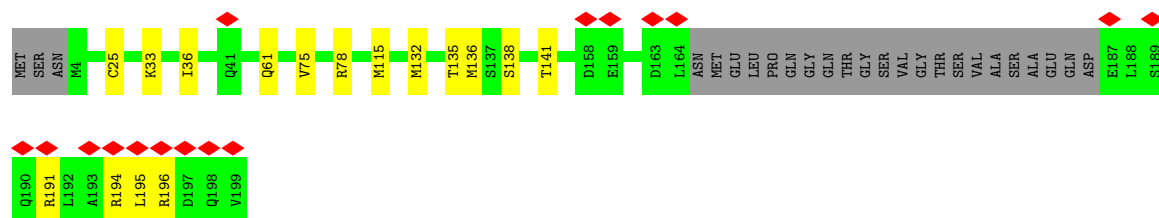
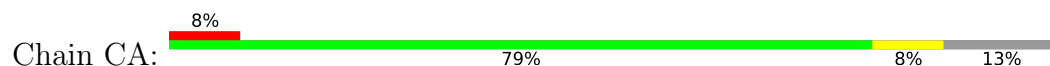
- Molecule 2: Charged multivesicular body protein 1b



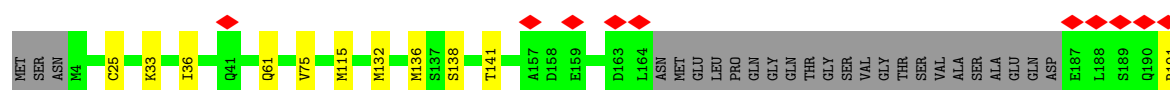
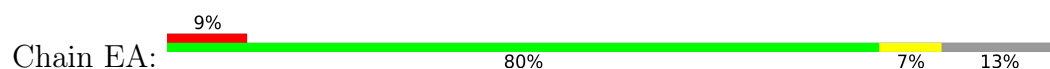
- Molecule 2: Charged multivesicular body protein 1b

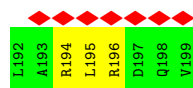


- Molecule 2: Charged multivesicular body protein 1b

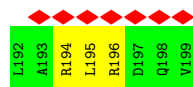
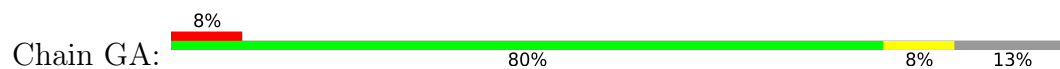


- Molecule 2: Charged multivesicular body protein 1b

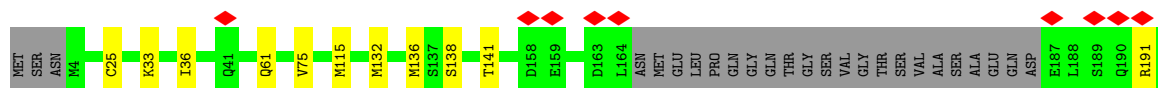
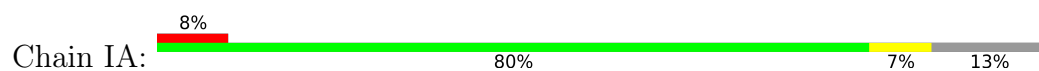




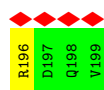
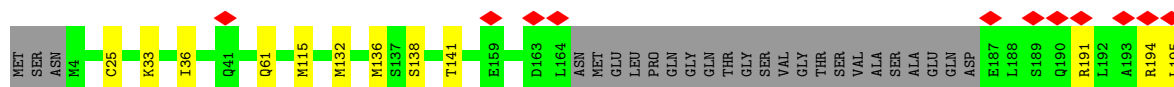
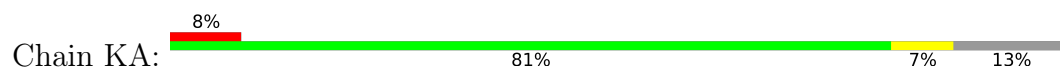
- Molecule 2: Charged multivesicular body protein 1b



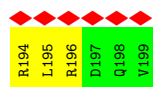
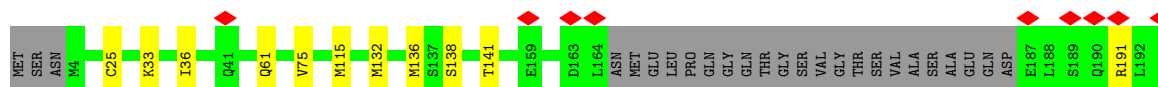
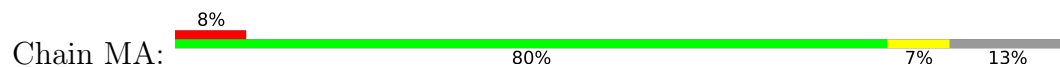
- Molecule 2: Charged multivesicular body protein 1b



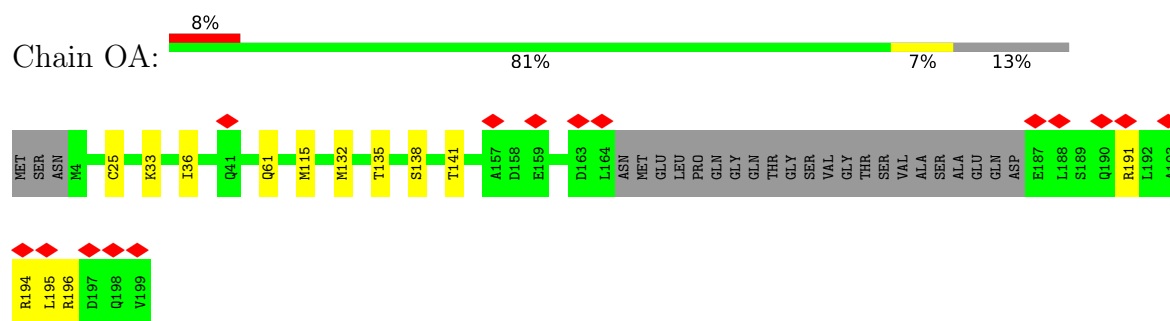
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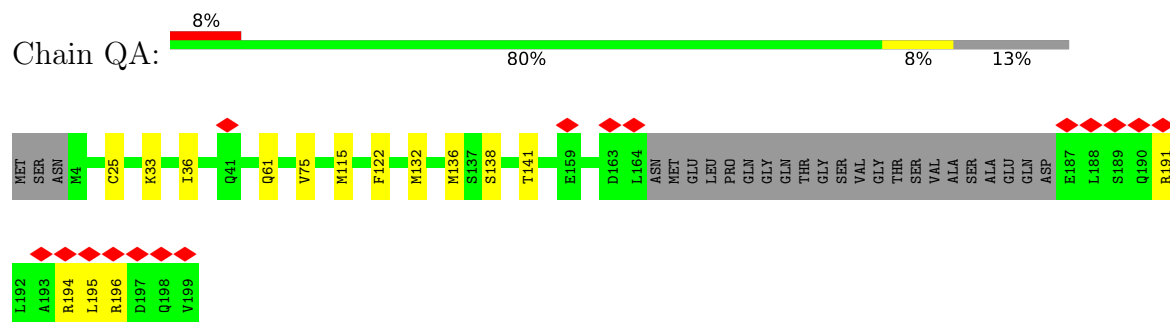
- Molecule 2: Charged multivesicular body protein 1b



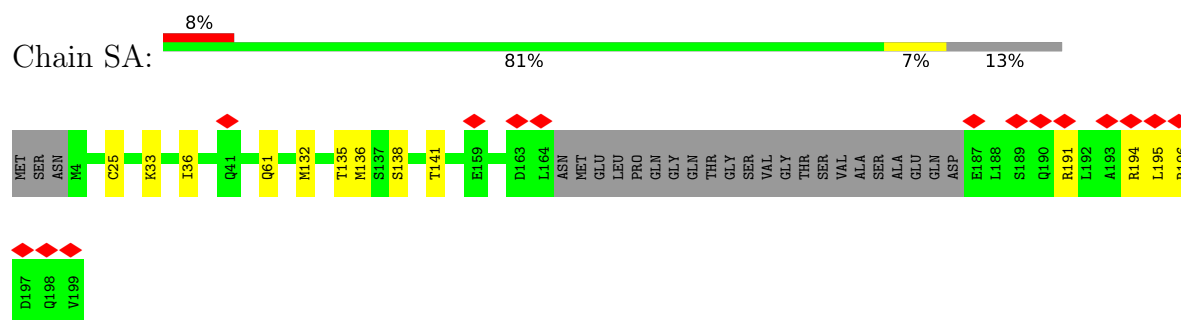
- Molecule 2: Charged multivesicular body protein 1b



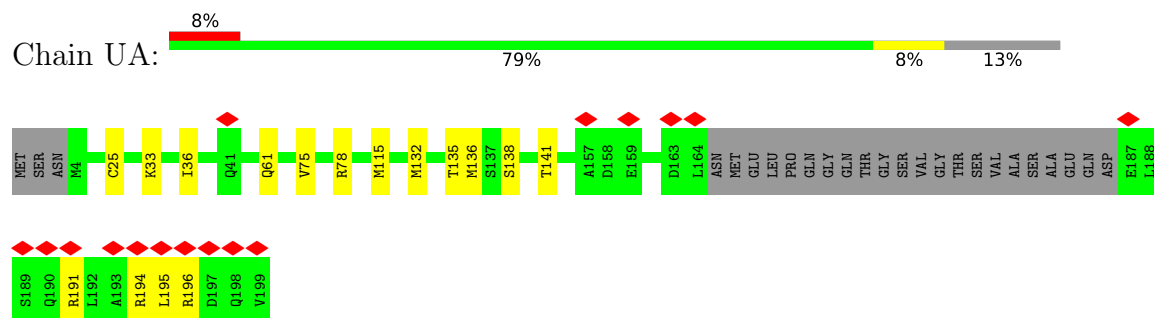
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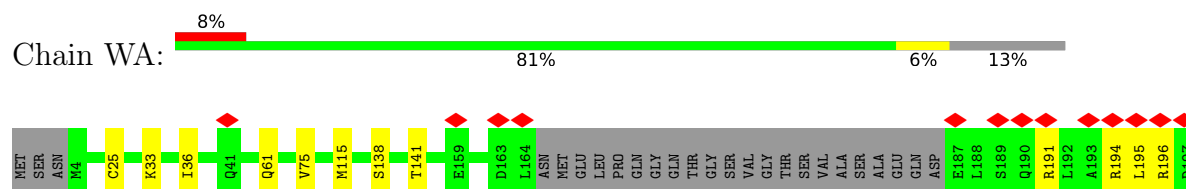
- Molecule 2: Charged multivesicular body protein 1b



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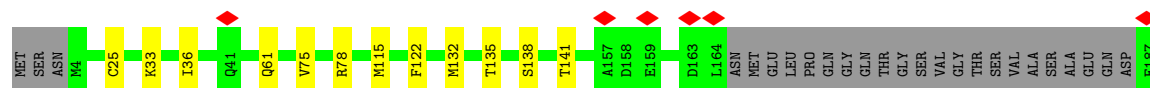
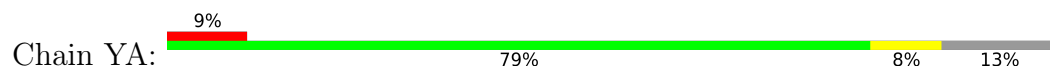


- Molecule 2: Charged multivesicular body protein 1b

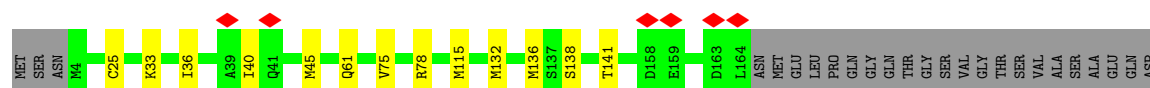
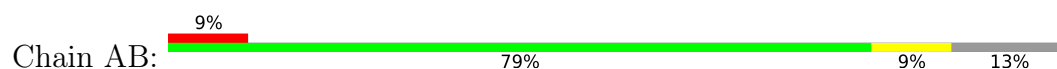




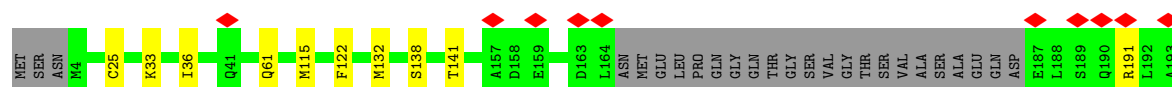
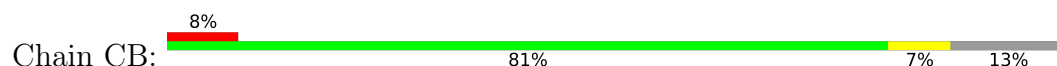
- Molecule 2: Charged multivesicular body protein 1b



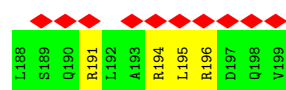
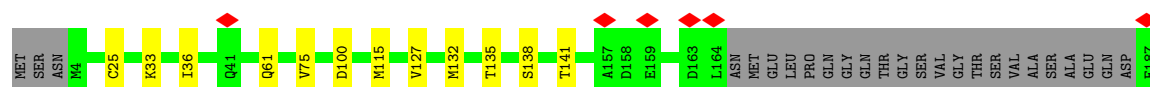
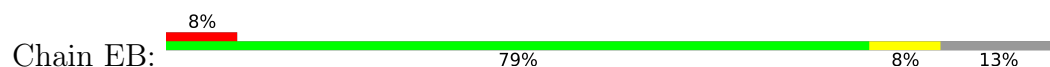
- Molecule 2: Charged multivesicular body protein 1b



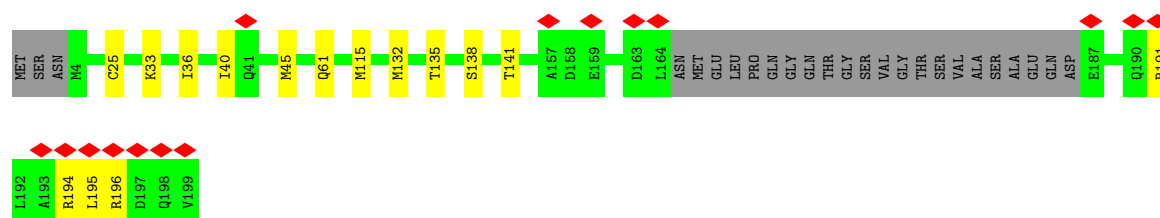
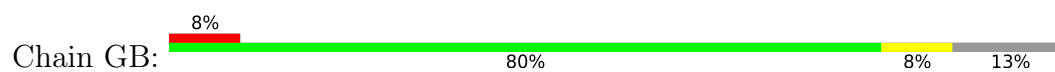
- Molecule 2: Charged multivesicular body protein 1b



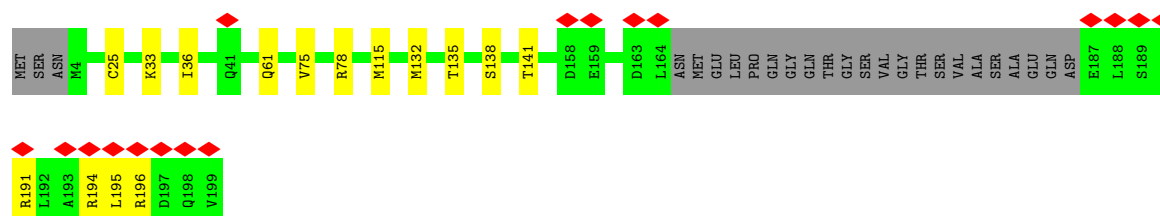
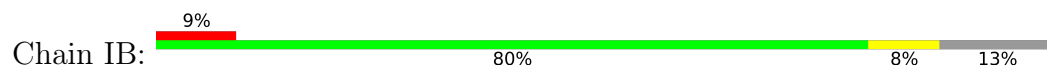
- Molecule 2: Charged multivesicular body protein 1b



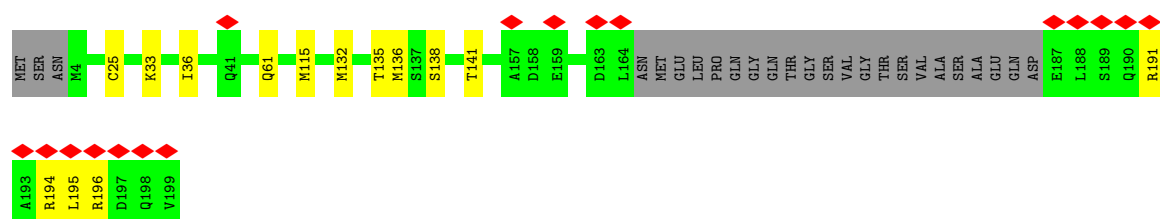
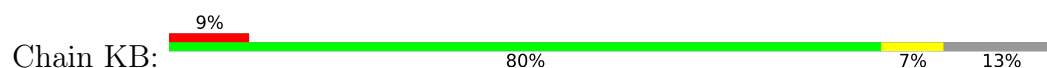
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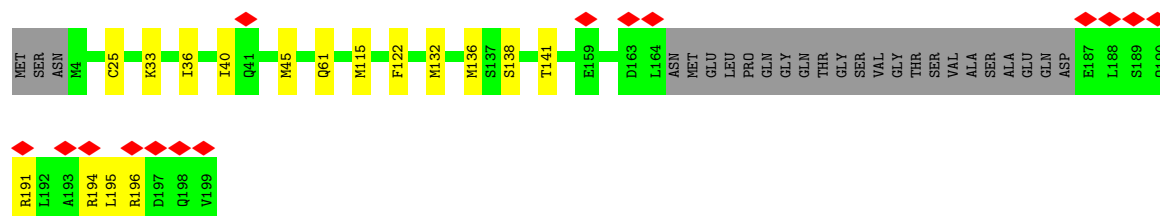
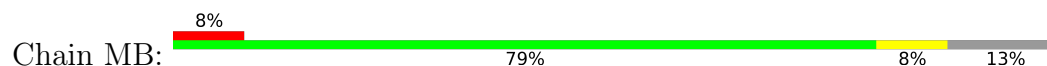
- Molecule 2: Charged multivesicular body protein 1b



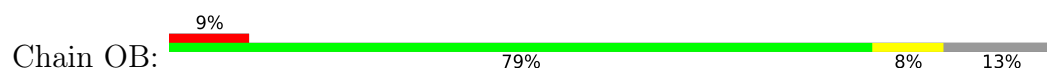
- Molecule 2: Charged multivesicular body protein 1b



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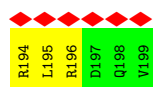
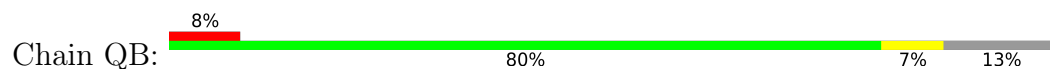


- Molecule 2: Charged multivesicular body protein 1b

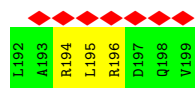
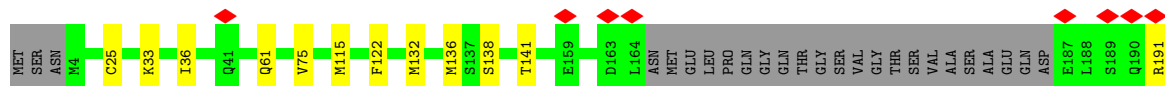
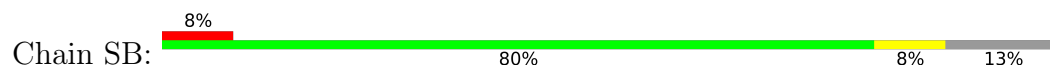




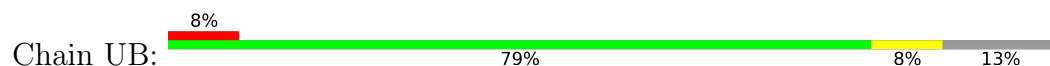
- Molecule 2: Charged multivesicular body protein 1b



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- Molecule 2: Charged multivesicular body protein 1b



4 Experimental information

Property	Value	Source
EM reconstruction method	HELICAL	Depositor
Imposed symmetry	HELICAL, twist=21.16°, rise=3.17 Å, axial sym=C1	Depositor
Number of segments used	101990	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{Å}^2$)	62	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	29000	Depositor
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.125	Depositor
Minimum map value	-0.063	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.008	Depositor
Recommended contour level	0.025	Depositor
Map size (Å)	335.2, 335.2, 335.2	wwPDB
Map dimensions	400, 400, 400	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.83800006, 0.83800006, 0.83800006	Depositor

5 Model quality

5.1 Standard geometry

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	A	0.25	0/1366	0.52	0/1838
1	BA	0.25	0/1366	0.52	0/1838
1	BB	0.26	0/1366	0.52	0/1838
1	C	0.25	0/1366	0.52	0/1838
1	DA	0.25	0/1366	0.52	0/1838
1	DB	0.26	0/1366	0.52	0/1838
1	E	0.25	0/1366	0.52	0/1838
1	FA	0.25	0/1366	0.52	0/1838
1	FB	0.25	0/1366	0.52	0/1838
1	G	0.25	0/1366	0.52	0/1838
1	HA	0.25	0/1366	0.52	0/1838
1	HB	0.25	0/1366	0.52	0/1838
1	I	0.25	0/1366	0.52	0/1838
1	JA	0.26	0/1366	0.52	0/1838
1	JB	0.25	0/1366	0.52	0/1838
1	K	0.25	0/1366	0.52	0/1838
1	LA	0.26	0/1366	0.52	0/1838
1	LB	0.26	0/1366	0.52	0/1838
1	M	0.25	0/1366	0.52	0/1838
1	NA	0.25	0/1366	0.52	0/1838
1	NB	0.25	0/1366	0.52	0/1838
1	O	0.25	0/1366	0.52	0/1838
1	PA	0.25	0/1366	0.52	0/1838
1	PB	0.26	0/1366	0.52	0/1838
1	Q	0.26	0/1366	0.52	0/1838
1	RA	0.26	0/1366	0.52	0/1838
1	RB	0.25	0/1366	0.52	0/1838
1	S	0.25	0/1366	0.52	0/1838
1	TA	0.25	0/1366	0.52	0/1838
1	TB	0.25	0/1366	0.52	0/1838
1	V	0.25	0/1366	0.52	0/1838
1	VA	0.25	0/1366	0.52	0/1838
1	X	0.25	0/1366	0.52	0/1838
1	XA	0.25	0/1366	0.52	0/1838

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	Z	0.25	0/1366	0.52	0/1838
1	ZA	0.25	0/1366	0.52	0/1838
2	AA	0.26	0/1360	0.48	0/1815
2	AB	0.26	0/1360	0.48	0/1815
2	B	0.26	0/1360	0.48	0/1815
2	CA	0.26	0/1360	0.48	0/1815
2	CB	0.26	0/1360	0.48	0/1815
2	D	0.26	0/1360	0.48	0/1815
2	EA	0.26	0/1360	0.48	0/1815
2	EB	0.26	0/1360	0.48	0/1815
2	F	0.26	0/1360	0.48	0/1815
2	GA	0.26	0/1360	0.48	0/1815
2	GB	0.26	0/1360	0.48	0/1815
2	H	0.26	0/1360	0.48	0/1815
2	IA	0.26	0/1360	0.48	0/1815
2	IB	0.26	0/1360	0.48	0/1815
2	J	0.26	0/1360	0.48	0/1815
2	KA	0.26	0/1360	0.48	0/1815
2	KB	0.26	0/1360	0.48	0/1815
2	L	0.26	0/1360	0.48	0/1815
2	MA	0.26	0/1360	0.48	0/1815
2	MB	0.26	0/1360	0.48	0/1815
2	N	0.26	0/1360	0.48	0/1815
2	OA	0.26	0/1360	0.48	0/1815
2	OB	0.26	0/1360	0.48	0/1815
2	P	0.26	0/1360	0.48	0/1815
2	QA	0.26	0/1360	0.48	0/1815
2	QB	0.26	0/1360	0.48	0/1815
2	R	0.26	0/1360	0.48	0/1815
2	SA	0.26	0/1360	0.48	0/1815
2	SB	0.26	0/1360	0.48	0/1815
2	T	0.26	0/1360	0.48	0/1815
2	UA	0.26	0/1360	0.48	0/1815
2	UB	0.26	0/1360	0.48	0/1815
2	W	0.26	0/1360	0.48	0/1815
2	WA	0.26	0/1360	0.48	0/1815
2	Y	0.26	0/1360	0.48	0/1815
2	YA	0.26	0/1360	0.48	0/1815
All	All	0.26	0/98136	0.50	0/131508

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	1349	0	1407	9	0
1	BA	1349	0	1407	8	0
1	BB	1349	0	1407	8	0
1	C	1349	0	1407	8	0
1	DA	1349	0	1407	8	0
1	DB	1349	0	1407	10	0
1	E	1349	0	1407	7	0
1	FA	1349	0	1407	7	0
1	FB	1349	0	1407	8	0
1	G	1349	0	1407	7	0
1	HA	1349	0	1407	8	0
1	HB	1349	0	1407	9	0
1	I	1349	0	1407	8	0
1	JA	1349	0	1407	9	0
1	JB	1349	0	1407	6	0
1	K	1349	0	1407	8	0
1	LA	1349	0	1407	7	0
1	LB	1349	0	1407	8	0
1	M	1349	0	1407	8	0
1	NA	1349	0	1407	8	0
1	NB	1349	0	1407	9	0
1	O	1349	0	1407	9	0
1	PA	1349	0	1407	8	0
1	PB	1349	0	1407	7	0
1	Q	1349	0	1407	7	0
1	RA	1349	0	1407	7	0
1	RB	1349	0	1407	9	0
1	S	1349	0	1407	8	0
1	TA	1349	0	1407	9	0
1	TB	1349	0	1407	9	0
1	V	1349	0	1407	13	0
1	VA	1349	0	1407	6	0
1	X	1349	0	1407	10	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	XA	1349	0	1407	8	0
1	Z	1349	0	1407	9	0
1	ZA	1349	0	1407	9	0
2	AA	1353	0	1385	10	0
2	AB	1353	0	1385	12	0
2	B	1353	0	1385	11	0
2	CA	1353	0	1385	13	0
2	CB	1353	0	1385	10	0
2	D	1353	0	1385	12	0
2	EA	1353	0	1385	10	0
2	EB	1353	0	1385	12	0
2	F	1353	0	1385	11	0
2	GA	1353	0	1385	12	0
2	GB	1353	0	1385	10	0
2	H	1353	0	1385	11	0
2	IA	1353	0	1385	10	0
2	IB	1353	0	1385	11	0
2	J	1353	0	1385	12	0
2	KA	1353	0	1385	10	0
2	KB	1353	0	1385	11	0
2	L	1353	0	1385	11	0
2	MA	1353	0	1385	11	0
2	MB	1353	0	1385	11	0
2	N	1353	0	1385	12	0
2	OA	1353	0	1385	9	0
2	OB	1353	0	1385	12	0
2	P	1353	0	1385	11	0
2	QA	1353	0	1385	11	0
2	QB	1353	0	1385	11	0
2	R	1353	0	1385	10	0
2	SA	1353	0	1385	9	0
2	SB	1353	0	1385	11	0
2	T	1353	0	1385	12	0
2	UA	1353	0	1385	12	0
2	UB	1353	0	1385	13	0
2	W	1353	0	1385	14	0
2	WA	1353	0	1385	9	0
2	Y	1353	0	1385	10	0
2	YA	1353	0	1385	12	0
All	All	97272	0	100512	510	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 3.

All (510) 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:V:160:ILE:HG22	1:V:164:ARG:NH2	1.97	0.80
1:RA:160:ILE:HG22	1:RA:164:ARG:NH2	1.97	0.80
1:HB:160:ILE:HG22	1:HB:164:ARG:NH2	1.97	0.80
1:DA:160:ILE:HG22	1:DA:164:ARG:NH2	1.97	0.80
1:NA:160:ILE:HG22	1:NA:164:ARG:NH2	1.97	0.80
1:BB:160:ILE:HG22	1:BB:164:ARG:NH2	1.97	0.80
1:DB:160:ILE:HG22	1:DB:164:ARG:NH2	1.97	0.80
1:BA:160:ILE:HG22	1:BA:164:ARG:NH2	1.97	0.80
1:FA:160:ILE:HG22	1:FA:164:ARG:NH2	1.97	0.80
1:HA:160:ILE:HG22	1:HA:164:ARG:NH2	1.97	0.80
1:TA:160:ILE:HG22	1:TA:164:ARG:NH2	1.97	0.80
1:VA:160:ILE:HG22	1:VA:164:ARG:NH2	1.97	0.80
1:NB:160:ILE:HG22	1:NB:164:ARG:NH2	1.97	0.80
1:I:160:ILE:HG22	1:I:164:ARG:NH2	1.97	0.80
1:Q:160:ILE:HG22	1:Q:164:ARG:NH2	1.97	0.80
1:Z:160:ILE:HG22	1:Z:164:ARG:NH2	1.97	0.80
1:K:160:ILE:HG22	1:K:164:ARG:NH2	1.97	0.80
1:C:160:ILE:HG22	1:C:164:ARG:NH2	1.97	0.80
1:E:160:ILE:HG22	1:E:164:ARG:NH2	1.97	0.80
1:G:160:ILE:HG22	1:G:164:ARG:NH2	1.97	0.80
1:LA:160:ILE:HG22	1:LA:164:ARG:NH2	1.97	0.80
1:XA:160:ILE:HG22	1:XA:164:ARG:NH2	1.97	0.80
1:FB:160:ILE:HG22	1:FB:164:ARG:NH2	1.97	0.80
1:JB:160:ILE:HG22	1:JB:164:ARG:NH2	1.97	0.80
1:PB:160:ILE:HG22	1:PB:164:ARG:NH2	1.97	0.80
1:M:160:ILE:HG22	1:M:164:ARG:NH2	1.97	0.80
1:O:160:ILE:HG22	1:O:164:ARG:NH2	1.97	0.80
1:LB:160:ILE:HG22	1:LB:164:ARG:NH2	1.97	0.80
1:RB:160:ILE:HG22	1:RB:164:ARG:NH2	1.97	0.80
1:X:160:ILE:HG22	1:X:164:ARG:NH2	1.97	0.80
1:JA:160:ILE:HG22	1:JA:164:ARG:NH2	1.97	0.80
1:PA:160:ILE:HG22	1:PA:164:ARG:NH2	1.97	0.80
1:TB:160:ILE:HG22	1:TB:164:ARG:NH2	1.97	0.79
1:A:160:ILE:HG22	1:A:164:ARG:NH2	1.97	0.79
1:S:160:ILE:HG22	1:S:164:ARG:NH2	1.97	0.79
1:ZA:160:ILE:HG22	1:ZA:164:ARG:NH2	1.97	0.79
2:W:132:MET:HG2	2:EB:115:MET:HE3	1.65	0.77
1:VA:165:TYR:OH	2:WA:196:ARG:NH1	2.20	0.74
1:BA:165:TYR:OH	2:CA:196:ARG:NH1	2.20	0.74
1:DB:165:TYR:OH	2:EB:196:ARG:NH1	2.20	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:HB:165:TYR:OH	2:IB:196:ARG:NH1	2.20	0.74
1:K:165:TYR:OH	2:L:196:ARG:NH1	2.20	0.74
1:V:165:TYR:OH	2:W:196:ARG:NH1	2.20	0.74
1:Z:165:TYR:OH	2:AA:196:ARG:NH1	2.20	0.74
1:FA:165:TYR:OH	2:GA:196:ARG:NH1	2.20	0.74
1:NA:165:TYR:OH	2:OA:196:ARG:NH1	2.20	0.74
1:BB:165:TYR:OH	2:CB:196:ARG:NH1	2.20	0.74
1:NB:165:TYR:OH	2:OB:196:ARG:NH1	2.20	0.74
1:RA:165:TYR:OH	2:SA:196:ARG:NH1	2.20	0.74
1:FB:165:TYR:OH	2:GB:196:ARG:NH1	2.20	0.74
1:LB:165:TYR:OH	2:MB:196:ARG:NH1	2.20	0.74
1:O:165:TYR:OH	2:P:196:ARG:NH1	2.20	0.74
1:A:165:TYR:OH	2:B:196:ARG:NH1	2.20	0.74
1:RB:165:TYR:OH	2:SB:196:ARG:NH1	2.20	0.74
1:JA:165:TYR:OH	2:KA:196:ARG:NH1	2.20	0.74
1:M:165:TYR:OH	2:N:196:ARG:NH1	2.20	0.73
1:TB:165:TYR:OH	2:UB:196:ARG:NH1	2.20	0.73
1:S:165:TYR:OH	2:T:196:ARG:NH1	2.20	0.73
1:ZA:165:TYR:OH	2:AB:196:ARG:NH1	2.20	0.73
1:JB:165:TYR:OH	2:KB:196:ARG:NH1	2.20	0.73
1:E:165:TYR:OH	2:F:196:ARG:NH1	2.20	0.73
1:XA:165:TYR:OH	2:YA:196:ARG:NH1	2.20	0.73
1:C:165:TYR:OH	2:D:196:ARG:NH1	2.20	0.73
1:PB:165:TYR:OH	2:QB:196:ARG:NH1	2.20	0.73
1:X:165:TYR:OH	2:Y:196:ARG:NH1	2.20	0.73
1:TA:165:TYR:OH	2:UA:196:ARG:NH1	2.20	0.73
1:G:165:TYR:OH	2:H:196:ARG:NH1	2.20	0.73
1:PA:165:TYR:OH	2:QA:196:ARG:NH1	2.20	0.73
1:I:165:TYR:OH	2:J:196:ARG:NH1	2.20	0.72
1:LA:165:TYR:OH	2:MA:196:ARG:NH1	2.20	0.72
1:Q:165:TYR:OH	2:R:196:ARG:NH1	2.20	0.72
1:HA:165:TYR:OH	2:IA:196:ARG:NH1	2.20	0.72
1:DA:165:TYR:OH	2:EA:196:ARG:NH1	2.20	0.72
2:Y:132:MET:HG2	2:UA:115:MET:HE3	1.76	0.68
2:SB:115:MET:HE3	2:UB:132:MET:HG2	1.76	0.68
2:MA:115:MET:HE3	2:UA:132:MET:HG2	1.75	0.67
2:N:132:MET:HG2	2:P:115:MET:HE3	1.77	0.67
2:L:115:MET:HE3	2:CA:132:MET:HG2	1.77	0.67
2:H:132:MET:HG2	2:J:115:MET:HE3	1.77	0.66
2:EA:132:MET:HG2	2:QA:115:MET:HE3	1.78	0.66
2:YA:115:MET:HE3	2:AB:132:MET:HG2	1.76	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:WA:115:MET:HE3	2:IB:132:MET:HG2	1.78	0.66
2:B:132:MET:HG2	2:AB:115:MET:HE3	1.77	0.66
2:D:115:MET:HE3	2:IA:132:MET:HG2	1.79	0.65
2:T:132:MET:HG2	2:QB:115:MET:HE3	1.79	0.65
2:D:132:MET:HG2	2:KB:115:MET:HE3	1.78	0.64
2:B:115:MET:HE3	2:MB:132:MET:HG2	1.78	0.64
2:SA:132:MET:HG2	2:CB:115:MET:HE3	1.79	0.64
2:J:132:MET:HG2	2:R:115:MET:HE3	1.79	0.63
2:EA:115:MET:HE3	2:MA:132:MET:HG2	1.80	0.63
2:KA:115:MET:HE3	2:GB:132:MET:HG2	1.80	0.63
2:Y:115:MET:HE3	2:QB:132:MET:HG2	1.80	0.63
2:F:132:MET:HG2	2:UB:115:MET:HE3	1.80	0.63
2:P:132:MET:HG2	2:AA:115:MET:HE3	1.79	0.63
2:GA:115:MET:HE3	2:OA:132:MET:HG2	1.79	0.63
2:N:115:MET:HE3	2:KB:132:MET:HG2	1.81	0.63
2:L:132:MET:HG2	2:MB:115:MET:HE3	1.80	0.62
2:CA:115:MET:HE3	2:EB:132:MET:HG2	1.81	0.62
2:AA:132:MET:HG2	2:IB:115:MET:HE3	1.81	0.61
2:H:115:MET:HE3	2:YA:132:MET:HG2	1.81	0.61
2:T:115:MET:HE3	2:KA:132:MET:HG2	1.82	0.61
1:I:160:ILE:CG2	1:I:164:ARG:NH2	2.65	0.60
1:Q:160:ILE:CG2	1:Q:164:ARG:NH2	2.64	0.60
2:GA:132:MET:HG2	2:GB:115:MET:HE3	1.82	0.60
1:LA:160:ILE:CG2	1:LA:164:ARG:NH2	2.64	0.60
1:TA:160:ILE:CG2	1:TA:164:ARG:NH2	2.65	0.60
1:G:160:ILE:CG2	1:G:164:ARG:NH2	2.64	0.60
2:R:132:MET:HG2	2:IA:115:MET:HE3	1.82	0.60
2:F:115:MET:HE3	2:QA:132:MET:HG2	1.82	0.60
1:X:160:ILE:CG2	1:X:164:ARG:NH2	2.65	0.60
1:DA:160:ILE:CG2	1:DA:164:ARG:NH2	2.64	0.60
1:HA:160:ILE:CG2	1:HA:164:ARG:NH2	2.65	0.60
2:OB:115:MET:HE3	2:SB:132:MET:HG2	1.82	0.60
1:V:160:ILE:CG2	1:V:164:ARG:NH2	2.65	0.60
1:HB:14:ASN:ND2	1:HB:116:GLU:OE1	2.35	0.60
1:NB:14:ASN:ND2	1:NB:116:GLU:OE1	2.35	0.60
1:K:14:ASN:ND2	1:K:116:GLU:OE1	2.35	0.60
1:O:14:ASN:ND2	1:O:116:GLU:OE1	2.35	0.60
1:V:14:ASN:ND2	1:V:116:GLU:OE1	2.35	0.60
1:Z:14:ASN:ND2	1:Z:116:GLU:OE1	2.35	0.60
1:BA:14:ASN:ND2	1:BA:116:GLU:OE1	2.35	0.60
1:RA:14:ASN:ND2	1:RA:116:GLU:OE1	2.35	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:RA:160:ILE:CG2	1:RA:164:ARG:NH2	2.65	0.60
1:VA:14:ASN:ND2	1:VA:116:GLU:OE1	2.35	0.60
1:BB:14:ASN:ND2	1:BB:116:GLU:OE1	2.35	0.60
1:LB:14:ASN:ND2	1:LB:116:GLU:OE1	2.35	0.60
1:A:14:ASN:ND2	1:A:116:GLU:OE1	2.35	0.59
1:JA:14:ASN:ND2	1:JA:116:GLU:OE1	2.35	0.59
1:NA:14:ASN:ND2	1:NA:116:GLU:OE1	2.35	0.59
1:BB:160:ILE:CG2	1:BB:164:ARG:NH2	2.65	0.59
1:DB:14:ASN:ND2	1:DB:116:GLU:OE1	2.35	0.59
1:DB:160:ILE:CG2	1:DB:164:ARG:NH2	2.65	0.59
1:FB:14:ASN:ND2	1:FB:116:GLU:OE1	2.35	0.59
1:HB:160:ILE:CG2	1:HB:164:ARG:NH2	2.65	0.59
1:NB:160:ILE:CG2	1:NB:164:ARG:NH2	2.65	0.59
1:RB:14:ASN:ND2	1:RB:116:GLU:OE1	2.35	0.59
1:Z:160:ILE:CG2	1:Z:164:ARG:NH2	2.65	0.59
1:FA:14:ASN:ND2	1:FA:116:GLU:OE1	2.35	0.59
1:M:14:ASN:ND2	1:M:116:GLU:OE1	2.35	0.59
1:BA:160:ILE:CG2	1:BA:164:ARG:NH2	2.65	0.59
1:PA:160:ILE:CG2	1:PA:164:ARG:NH2	2.65	0.59
1:VA:160:ILE:CG2	1:VA:164:ARG:NH2	2.65	0.59
1:C:160:ILE:CG2	1:C:164:ARG:NH2	2.65	0.59
1:NA:160:ILE:CG2	1:NA:164:ARG:NH2	2.65	0.59
1:RB:160:ILE:CG2	1:RB:164:ARG:NH2	2.65	0.59
1:TB:14:ASN:ND2	1:TB:116:GLU:OE1	2.35	0.59
1:S:14:ASN:ND2	1:S:116:GLU:OE1	2.35	0.59
1:M:160:ILE:CG2	1:M:164:ARG:NH2	2.64	0.59
1:O:160:ILE:CG2	1:O:164:ARG:NH2	2.65	0.59
1:ZA:14:ASN:ND2	1:ZA:116:GLU:OE1	2.35	0.59
1:JB:14:ASN:ND2	1:JB:116:GLU:OE1	2.35	0.59
1:E:14:ASN:ND2	1:E:116:GLU:OE1	2.35	0.59
1:FA:160:ILE:CG2	1:FA:164:ARG:NH2	2.64	0.59
1:JA:160:ILE:CG2	1:JA:164:ARG:NH2	2.64	0.59
1:XA:14:ASN:ND2	1:XA:116:GLU:OE1	2.35	0.59
1:FB:160:ILE:CG2	1:FB:164:ARG:NH2	2.65	0.59
1:LB:160:ILE:CG2	1:LB:164:ARG:NH2	2.64	0.59
1:TB:160:ILE:CG2	1:TB:164:ARG:NH2	2.65	0.59
1:A:160:ILE:CG2	1:A:164:ARG:NH2	2.65	0.59
1:C:14:ASN:ND2	1:C:116:GLU:OE1	2.35	0.59
1:K:160:ILE:CG2	1:K:164:ARG:NH2	2.65	0.59
1:PA:14:ASN:ND2	1:PA:116:GLU:OE1	2.35	0.59
1:PB:14:ASN:ND2	1:PB:116:GLU:OE1	2.35	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:14:ASN:ND2	1:G:116:GLU:OE1	2.35	0.59
1:X:14:ASN:ND2	1:X:116:GLU:OE1	2.35	0.58
1:HA:14:ASN:ND2	1:HA:116:GLU:OE1	2.35	0.58
1:S:160:ILE:CG2	1:S:164:ARG:NH2	2.65	0.58
2:W:115:MET:HE3	2:OB:132:MET:HG2	1.83	0.58
1:DA:14:ASN:ND2	1:DA:116:GLU:OE1	2.35	0.58
1:TA:14:ASN:ND2	1:TA:116:GLU:OE1	2.35	0.58
1:I:14:ASN:ND2	1:I:116:GLU:OE1	2.35	0.58
1:Q:14:ASN:ND2	1:Q:116:GLU:OE1	2.35	0.58
1:LA:14:ASN:ND2	1:LA:116:GLU:OE1	2.35	0.58
1:ZA:160:ILE:CG2	1:ZA:164:ARG:NH2	2.64	0.58
1:E:160:ILE:CG2	1:E:164:ARG:NH2	2.65	0.58
1:JB:160:ILE:CG2	1:JB:164:ARG:NH2	2.65	0.58
1:PB:160:ILE:CG2	1:PB:164:ARG:NH2	2.65	0.58
2:OA:115:MET:HE3	2:CB:132:MET:HG2	1.84	0.57
1:XA:160:ILE:CG2	1:XA:164:ARG:NH2	2.65	0.57
2:P:33:LYS:HA	2:P:36:ILE:HD12	1.87	0.57
2:T:33:LYS:HA	2:T:36:ILE:HD12	1.87	0.57
2:D:33:LYS:HA	2:D:36:ILE:HD12	1.87	0.57
2:J:33:LYS:HA	2:J:36:ILE:HD12	1.87	0.57
2:AB:33:LYS:HA	2:AB:36:ILE:HD12	1.87	0.57
2:MA:33:LYS:HA	2:MA:36:ILE:HD12	1.87	0.57
2:QA:33:LYS:HA	2:QA:36:ILE:HD12	1.87	0.57
2:UA:33:LYS:HA	2:UA:36:ILE:HD12	1.87	0.57
2:QB:33:LYS:HA	2:QB:36:ILE:HD12	1.87	0.57
2:SB:33:LYS:HA	2:SB:36:ILE:HD12	1.87	0.57
2:F:33:LYS:HA	2:F:36:ILE:HD12	1.87	0.57
2:GB:33:LYS:HA	2:GB:36:ILE:HD12	1.87	0.57
2:KB:33:LYS:HA	2:KB:36:ILE:HD12	1.87	0.57
2:J:75:VAL:HG21	2:AB:136:MET:HG2	1.86	0.57
2:R:33:LYS:HA	2:R:36:ILE:HD12	1.87	0.57
2:YA:33:LYS:HA	2:YA:36:ILE:HD12	1.87	0.57
2:MB:33:LYS:HA	2:MB:36:ILE:HD12	1.87	0.57
2:CB:33:LYS:HA	2:CB:36:ILE:HD12	1.87	0.57
2:EB:33:LYS:HA	2:EB:36:ILE:HD12	1.87	0.57
2:WA:33:LYS:HA	2:WA:36:ILE:HD12	1.87	0.56
2:AA:33:LYS:HA	2:AA:36:ILE:HD12	1.87	0.56
2:OB:33:LYS:HA	2:OB:36:ILE:HD12	1.87	0.56
2:L:33:LYS:HA	2:L:36:ILE:HD12	1.87	0.56
2:GA:33:LYS:HA	2:GA:36:ILE:HD12	1.87	0.56
2:L:75:VAL:HG21	2:W:136:MET:HG2	1.86	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:N:33:LYS:HA	2:N:36:ILE:HD12	1.87	0.56
2:SA:33:LYS:HA	2:SA:36:ILE:HD12	1.87	0.56
2:IB:33:LYS:HA	2:IB:36:ILE:HD12	1.87	0.56
2:H:33:LYS:HA	2:H:36:ILE:HD12	1.87	0.56
2:W:33:LYS:HA	2:W:36:ILE:HD12	1.87	0.56
2:UB:33:LYS:HA	2:UB:36:ILE:HD12	1.87	0.56
2:CA:33:LYS:HA	2:CA:36:ILE:HD12	1.87	0.56
2:IA:33:LYS:HA	2:IA:36:ILE:HD12	1.87	0.56
2:OA:33:LYS:HA	2:OA:36:ILE:HD12	1.87	0.56
2:Y:33:LYS:HA	2:Y:36:ILE:HD12	1.87	0.56
2:KA:33:LYS:HA	2:KA:36:ILE:HD12	1.87	0.56
2:EA:33:LYS:HA	2:EA:36:ILE:HD12	1.87	0.55
2:B:33:LYS:HA	2:B:36:ILE:HD12	1.87	0.55
2:W:75:VAL:HG21	2:UB:136:MET:HG2	1.90	0.53
1:NA:54:ILE:HD12	1:BB:78:LEU:HD23	1.90	0.52
1:E:54:ILE:HD12	1:PA:78:LEU:HD23	1.93	0.51
2:CA:115:MET:HB3	2:EB:135:THR:HG21	1.93	0.50
1:V:21:ARG:HD3	2:EB:127:VAL:HG21	1.94	0.49
1:V:78:LEU:HD23	1:DB:54:ILE:HD12	1.94	0.49
2:B:75:VAL:HG21	2:CA:136:MET:HG2	1.95	0.48
2:L:136:MET:HG2	2:AB:75:VAL:HG21	1.93	0.48
1:V:77:ASP:OD1	1:DB:58:HIS:NE2	2.46	0.48
1:XA:54:ILE:HD12	1:ZA:78:LEU:HD23	1.95	0.48
2:T:136:MET:HG2	2:UA:75:VAL:HG21	1.96	0.48
1:O:78:LEU:HD23	1:Z:54:ILE:HD12	1.96	0.48
1:FA:78:LEU:HD23	1:FB:54:ILE:HD12	1.95	0.47
1:V:73:GLU:HG2	1:DB:51:ARG:NH1	2.29	0.47
2:Y:136:MET:HG2	2:EA:75:VAL:HG21	1.97	0.47
2:EB:75:VAL:HG21	2:SB:136:MET:HG2	1.96	0.47
1:V:50:GLU:HG2	1:NB:74:LEU:HD21	1.97	0.47
1:NB:54:ILE:HD12	1:RB:78:LEU:HD23	1.97	0.47
2:CA:75:VAL:HG21	2:OB:136:MET:HG2	1.96	0.46
2:B:136:MET:HG2	2:H:75:VAL:HG21	1.98	0.46
2:P:136:MET:HG2	2:WA:75:VAL:HG21	1.97	0.46
2:H:136:MET:HG2	2:IA:75:VAL:HG21	1.98	0.46
2:N:136:MET:HG2	2:IB:75:VAL:HG21	1.97	0.46
2:QA:75:VAL:HG21	2:UA:136:MET:HG2	1.97	0.46
1:X:78:LEU:HD23	1:TA:54:ILE:HD12	1.97	0.45
2:EA:25:CYS:SG	2:EA:61:GLN:NE2	2.90	0.45
2:AB:78:ARG:HE	2:MB:122:PHE:HB3	1.82	0.45
2:SB:25:CYS:SG	2:SB:61:GLN:NE2	2.90	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:T:25:CYS:SG	2:T:61:GLN:NE2	2.90	0.45
2:P:25:CYS:SG	2:P:61:GLN:NE2	2.90	0.45
2:IA:25:CYS:SG	2:IA:61:GLN:NE2	2.90	0.45
2:QA:122:PHE:HB3	2:UB:78:ARG:HE	1.81	0.45
2:AB:25:CYS:SG	2:AB:61:GLN:NE2	2.90	0.45
2:QB:25:CYS:SG	2:QB:61:GLN:NE2	2.90	0.45
2:J:25:CYS:SG	2:J:61:GLN:NE2	2.90	0.45
2:N:25:CYS:SG	2:N:61:GLN:NE2	2.90	0.45
2:N:75:VAL:HG21	2:IA:136:MET:HG2	1.99	0.45
2:UA:25:CYS:SG	2:UA:61:GLN:NE2	2.90	0.45
2:YA:25:CYS:SG	2:YA:61:GLN:NE2	2.90	0.45
2:D:78:ARG:HE	2:R:122:PHE:HB3	1.82	0.45
2:D:136:MET:HG2	2:P:75:VAL:HG21	1.98	0.45
2:Y:75:VAL:HG21	2:KA:136:MET:HG2	1.99	0.45
2:MA:115:MET:HB3	2:UA:135:THR:HG21	1.98	0.45
2:WA:25:CYS:SG	2:WA:61:GLN:NE2	2.90	0.45
2:CB:25:CYS:SG	2:CB:61:GLN:NE2	2.90	0.45
2:EB:25:CYS:SG	2:EB:61:GLN:NE2	2.90	0.45
2:UB:25:CYS:SG	2:UB:61:GLN:NE2	2.90	0.45
2:F:25:CYS:SG	2:F:61:GLN:NE2	2.90	0.45
2:R:25:CYS:SG	2:R:61:GLN:NE2	2.90	0.45
2:W:78:ARG:HE	2:SB:122:PHE:HB3	1.81	0.45
2:MA:25:CYS:SG	2:MA:61:GLN:NE2	2.90	0.45
2:KB:25:CYS:SG	2:KB:61:GLN:NE2	2.90	0.45
2:D:25:CYS:SG	2:D:61:GLN:NE2	2.90	0.45
1:BA:113:GLU:HG3	1:BA:114:VAL:HG22	1.99	0.45
2:QA:25:CYS:SG	2:QA:61:GLN:NE2	2.90	0.45
1:VA:113:GLU:HG3	1:VA:114:VAL:HG22	1.99	0.45
2:GB:25:CYS:SG	2:GB:61:GLN:NE2	2.90	0.45
2:IB:25:CYS:SG	2:IB:61:GLN:NE2	2.90	0.45
2:MB:25:CYS:SG	2:MB:61:GLN:NE2	2.90	0.45
2:Y:25:CYS:SG	2:Y:61:GLN:NE2	2.90	0.45
1:NA:113:GLU:HG3	1:NA:114:VAL:HG22	1.99	0.45
2:SA:25:CYS:SG	2:SA:61:GLN:NE2	2.90	0.45
2:YA:75:VAL:HG21	2:MB:136:MET:HG2	1.98	0.45
2:H:25:CYS:SG	2:H:61:GLN:NE2	2.90	0.44
1:K:113:GLU:HG3	1:K:114:VAL:HG22	1.99	0.44
2:L:25:CYS:SG	2:L:61:GLN:NE2	2.90	0.44
1:V:113:GLU:HG3	1:V:114:VAL:HG22	1.99	0.44
2:W:25:CYS:SG	2:W:61:GLN:NE2	2.90	0.44
2:QA:136:MET:HG2	2:SB:75:VAL:HG21	1.98	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BB:113:GLU:HG3	1:BB:114:VAL:HG22	1.99	0.44
1:DB:113:GLU:HG3	1:DB:114:VAL:HG22	1.99	0.44
1:HB:113:GLU:HG3	1:HB:114:VAL:HG22	1.99	0.44
1:LB:113:GLU:HG3	1:LB:114:VAL:HG22	1.99	0.44
2:OB:25:CYS:SG	2:OB:61:GLN:NE2	2.90	0.44
2:AA:25:CYS:SG	2:AA:61:GLN:NE2	2.90	0.44
1:DA:54:ILE:HD12	1:LA:78:LEU:HD23	1.98	0.44
1:FA:113:GLU:HG3	1:FA:114:VAL:HG22	1.99	0.44
2:GA:25:CYS:SG	2:GA:61:GLN:NE2	2.90	0.44
1:RA:113:GLU:HG3	1:RA:114:VAL:HG22	1.99	0.44
1:FB:113:GLU:HG3	1:FB:114:VAL:HG22	1.99	0.44
1:Z:113:GLU:HG3	1:Z:114:VAL:HG22	1.99	0.44
1:Q:113:GLU:HG3	1:Q:114:VAL:HG22	1.99	0.44
1:NB:113:GLU:HG3	1:NB:114:VAL:HG22	1.99	0.44
1:A:113:GLU:HG3	1:A:114:VAL:HG22	1.99	0.44
1:S:113:GLU:HG3	1:S:114:VAL:HG22	1.99	0.44
2:B:25:CYS:SG	2:B:61:GLN:NE2	2.90	0.44
1:C:54:ILE:HD12	1:HA:78:LEU:HD23	1.99	0.44
2:CA:25:CYS:SG	2:CA:61:GLN:NE2	2.90	0.44
1:DA:113:GLU:HG3	1:DA:114:VAL:HG22	1.99	0.44
1:JA:113:GLU:HG3	1:JA:114:VAL:HG22	1.99	0.44
1:LA:113:GLU:HG3	1:LA:114:VAL:HG22	1.99	0.44
1:RB:113:GLU:HG3	1:RB:114:VAL:HG22	1.99	0.44
1:O:113:GLU:HG3	1:O:114:VAL:HG22	1.99	0.44
1:Z:78:LEU:HD23	1:HB:54:ILE:HD12	2.00	0.44
1:HA:113:GLU:HG3	1:HA:114:VAL:HG22	1.99	0.44
2:KA:25:CYS:SG	2:KA:61:GLN:NE2	2.90	0.44
2:MA:75:VAL:HG21	2:QB:136:MET:HG2	1.99	0.44
2:OA:25:CYS:SG	2:OA:61:GLN:NE2	2.90	0.44
1:ZA:113:GLU:HG3	1:ZA:114:VAL:HG22	1.99	0.44
1:JB:113:GLU:HG3	1:JB:114:VAL:HG22	1.99	0.44
1:E:113:GLU:HG3	1:E:114:VAL:HG22	1.99	0.44
1:X:113:GLU:HG3	1:X:114:VAL:HG22	1.99	0.44
1:G:113:GLU:HG3	1:G:114:VAL:HG22	1.99	0.44
2:D:135:THR:HG21	2:KB:115:MET:HB3	1.99	0.43
2:J:135:THR:HG21	2:R:115:MET:HB3	2.00	0.43
1:XA:113:GLU:HG3	1:XA:114:VAL:HG22	1.99	0.43
1:PB:113:GLU:HG3	1:PB:114:VAL:HG22	1.99	0.43
1:RB:54:ILE:HD12	1:TB:78:LEU:HD23	2.00	0.43
1:TB:113:GLU:HG3	1:TB:114:VAL:HG22	1.99	0.43
1:M:113:GLU:HG3	1:M:114:VAL:HG22	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:GA:115:MET:HB3	2:OA:135:THR:HG21	1.99	0.43
1:C:113:GLU:HG3	1:C:114:VAL:HG22	1.99	0.43
1:FB:125:CYS:O	1:FB:129:SER:OG	2.36	0.43
1:LB:125:CYS:O	1:LB:129:SER:OG	2.36	0.43
2:F:75:VAL:HG21	2:MA:136:MET:HG2	2.00	0.43
2:IA:138:SER:HA	2:IA:141:THR:HG23	2.01	0.43
2:QB:138:SER:HA	2:QB:141:THR:HG23	2.01	0.43
1:I:113:GLU:HG3	1:I:114:VAL:HG22	1.99	0.43
2:J:191:ARG:HG2	2:J:194:ARG:HH21	1.84	0.43
2:EA:138:SER:HA	2:EA:141:THR:HG23	2.01	0.43
1:PA:113:GLU:HG3	1:PA:114:VAL:HG22	1.99	0.43
1:TA:113:GLU:HG3	1:TA:114:VAL:HG22	1.99	0.43
2:IA:191:ARG:HG2	2:IA:194:ARG:HH21	1.84	0.43
2:QA:138:SER:HA	2:QA:141:THR:HG23	2.01	0.43
2:QA:191:ARG:HG2	2:QA:194:ARG:HH21	1.84	0.43
2:UA:191:ARG:HG2	2:UA:194:ARG:HH21	1.84	0.43
2:YA:138:SER:HA	2:YA:141:THR:HG23	2.01	0.43
2:D:191:ARG:HG2	2:D:194:ARG:HH21	1.84	0.43
1:E:164:ARG:HD3	2:F:195:LEU:HD11	2.01	0.43
2:J:78:ARG:HE	2:YA:122:PHE:HB3	1.84	0.43
2:EA:136:MET:HG2	2:UB:75:VAL:HG21	2.00	0.43
2:EA:191:ARG:HG2	2:EA:194:ARG:HH21	1.84	0.43
2:KB:191:ARG:HG2	2:KB:194:ARG:HH21	1.84	0.43
1:RB:164:ARG:HD3	2:SB:195:LEU:HD11	2.01	0.43
2:D:138:SER:HA	2:D:141:THR:HG23	2.01	0.42
2:F:191:ARG:HG2	2:F:194:ARG:HH21	1.84	0.42
1:O:164:ARG:HD3	2:P:195:LEU:HD11	2.01	0.42
1:JB:164:ARG:HD3	2:KB:195:LEU:HD11	2.01	0.42
1:S:54:ILE:HD12	1:JA:78:LEU:HD23	2.00	0.42
2:D:75:VAL:HG21	2:J:136:MET:HG2	2.00	0.42
2:H:138:SER:HA	2:H:141:THR:HG23	2.01	0.42
2:R:138:SER:HA	2:R:141:THR:HG23	2.01	0.42
2:Y:138:SER:HA	2:Y:141:THR:HG23	2.01	0.42
2:AA:75:VAL:HG21	2:KB:136:MET:HG2	2.00	0.42
2:MA:191:ARG:HG2	2:MA:194:ARG:HH21	1.84	0.42
1:VA:164:ARG:HD3	2:WA:195:LEU:HD11	2.01	0.42
1:BB:164:ARG:HD3	2:CB:195:LEU:HD11	2.01	0.42
1:DB:164:ARG:HD3	2:EB:195:LEU:HD11	2.01	0.42
2:B:122:PHE:HB3	2:YA:78:ARG:HE	1.83	0.42
2:R:191:ARG:HG2	2:R:194:ARG:HH21	1.84	0.42
2:W:122:PHE:HB3	2:CA:78:ARG:HE	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:MA:138:SER:HA	2:MA:141:THR:HG23	2.01	0.42
2:YA:191:ARG:HG2	2:YA:194:ARG:HH21	1.84	0.42
1:NB:164:ARG:HD3	2:OB:195:LEU:HD11	2.01	0.42
2:QB:191:ARG:HG2	2:QB:194:ARG:HH21	1.84	0.42
2:W:115:MET:HB3	2:OB:135:THR:HG21	2.00	0.42
1:Z:164:ARG:HD3	2:AA:195:LEU:HD11	2.01	0.42
2:SB:138:SER:HA	2:SB:141:THR:HG23	2.01	0.42
2:T:138:SER:HA	2:T:141:THR:HG23	2.01	0.42
2:N:138:SER:HA	2:N:141:THR:HG23	2.01	0.42
2:P:138:SER:HA	2:P:141:THR:HG23	2.01	0.42
2:GB:138:SER:HA	2:GB:141:THR:HG23	2.01	0.42
1:A:125:CYS:O	1:A:129:SER:OG	2.36	0.42
1:C:164:ARG:HD3	2:D:195:LEU:HD11	2.01	0.42
1:JA:125:CYS:O	1:JA:129:SER:OG	2.36	0.42
2:AB:138:SER:HA	2:AB:141:THR:HG23	2.01	0.42
2:MB:138:SER:HA	2:MB:141:THR:HG23	2.01	0.42
2:UB:138:SER:HA	2:UB:141:THR:HG23	2.01	0.42
2:H:191:ARG:HG2	2:H:194:ARG:HH21	1.84	0.42
2:L:191:ARG:HG2	2:L:194:ARG:HH21	1.84	0.42
2:CA:191:ARG:HG2	2:CA:194:ARG:HH21	1.84	0.42
2:GA:75:VAL:HG21	2:SA:136:MET:HG2	2.00	0.42
2:KA:115:MET:HB3	2:GB:135:THR:HG21	2.01	0.42
1:PA:164:ARG:HD3	2:QA:195:LEU:HD11	2.01	0.42
2:SA:135:THR:HG21	2:CB:115:MET:HB3	2.01	0.42
2:WA:115:MET:HB3	2:IB:135:THR:HG21	2.01	0.42
1:HB:164:ARG:HD3	2:IB:195:LEU:HD11	2.01	0.42
2:T:135:THR:HG21	2:QB:115:MET:HB3	2.01	0.42
2:T:191:ARG:HG2	2:T:194:ARG:HH21	1.84	0.42
1:M:78:LEU:HD23	1:O:54:ILE:HD12	2.02	0.42
2:N:191:ARG:HG2	2:N:194:ARG:HH21	1.84	0.42
1:Q:164:ARG:HD3	2:R:195:LEU:HD11	2.01	0.42
2:GA:191:ARG:HG2	2:GA:194:ARG:HH21	1.84	0.42
1:LA:164:ARG:HD3	2:MA:195:LEU:HD11	2.01	0.42
2:WA:191:ARG:HG2	2:WA:194:ARG:HH21	1.84	0.42
1:XA:125:CYS:O	1:XA:129:SER:OG	2.36	0.42
2:AB:191:ARG:HG2	2:AB:194:ARG:HH21	1.84	0.42
2:CB:191:ARG:HG2	2:CB:194:ARG:HH21	1.84	0.42
2:EB:191:ARG:HG2	2:EB:194:ARG:HH21	1.84	0.42
2:L:115:MET:HB3	2:CA:135:THR:HG21	2.02	0.42
1:V:164:ARG:HD3	2:W:195:LEU:HD11	2.01	0.42
1:JA:160:ILE:CG2	1:JA:164:ARG:HH22	2.33	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:OA:191:ARG:HG2	2:OA:194:ARG:HH21	1.84	0.42
1:A:160:ILE:CG2	1:A:164:ARG:HH22	2.33	0.42
1:V:125:CYS:O	1:V:129:SER:OG	2.36	0.42
1:X:50:GLU:HG2	1:PB:74:LEU:HD21	2.01	0.42
2:CB:138:SER:HA	2:CB:141:THR:HG23	2.01	0.42
1:HB:125:CYS:O	1:HB:129:SER:OG	2.36	0.42
1:TB:125:CYS:O	1:TB:129:SER:OG	2.36	0.42
1:TB:160:ILE:CG2	1:TB:164:ARG:HH22	2.33	0.42
2:UB:191:ARG:HG2	2:UB:194:ARG:HH21	1.84	0.42
2:F:136:MET:HG2	2:OB:75:VAL:HG21	2.01	0.41
2:L:138:SER:HA	2:L:141:THR:HG23	2.01	0.41
1:M:160:ILE:CG2	1:M:164:ARG:HH22	2.33	0.41
1:O:160:ILE:CG2	1:O:164:ARG:HH22	2.33	0.41
2:Y:191:ARG:HG2	2:Y:194:ARG:HH21	1.84	0.41
2:AA:191:ARG:HG2	2:AA:194:ARG:HH21	1.84	0.41
1:RA:164:ARG:HD3	2:SA:195:LEU:HD11	2.01	0.41
2:WA:138:SER:HA	2:WA:141:THR:HG23	2.01	0.41
2:EB:138:SER:HA	2:EB:141:THR:HG23	2.01	0.41
1:RB:160:ILE:CG2	1:RB:164:ARG:HH22	2.33	0.41
2:SB:191:ARG:HG2	2:SB:194:ARG:HH21	1.84	0.41
2:B:191:ARG:HG2	2:B:194:ARG:HH21	1.84	0.41
1:G:160:ILE:CG2	1:G:164:ARG:HH22	2.33	0.41
2:H:115:MET:HB3	2:YA:135:THR:HG21	2.02	0.41
1:K:54:ILE:HD12	1:BA:78:LEU:HD23	2.03	0.41
2:P:191:ARG:HG2	2:P:194:ARG:HH21	1.84	0.41
2:KA:191:ARG:HG2	2:KA:194:ARG:HH21	1.84	0.41
2:UA:138:SER:HA	2:UA:141:THR:HG23	2.01	0.41
1:LB:164:ARG:HD3	2:MB:195:LEU:HD11	2.01	0.41
2:OB:191:ARG:HG2	2:OB:194:ARG:HH21	1.84	0.41
2:J:138:SER:HA	2:J:141:THR:HG23	2.01	0.41
1:K:164:ARG:HD3	2:L:195:LEU:HD11	2.01	0.41
1:X:160:ILE:CG2	1:X:164:ARG:HH22	2.33	0.41
1:BA:164:ARG:HD3	2:CA:195:LEU:HD11	2.01	0.41
2:GA:78:ARG:HE	2:CB:122:PHE:HB3	1.85	0.41
2:GA:138:SER:HA	2:GA:141:THR:HG23	2.01	0.41
1:NA:164:ARG:HD3	2:OA:195:LEU:HD11	2.01	0.41
1:FB:164:ARG:HD3	2:GB:195:LEU:HD11	2.01	0.41
1:PB:164:ARG:HD3	2:QB:195:LEU:HD11	2.01	0.41
1:S:164:ARG:HD3	2:T:195:LEU:HD11	2.01	0.41
2:F:138:SER:HA	2:F:141:THR:HG23	2.01	0.41
1:DA:164:ARG:HD3	2:EA:195:LEU:HD11	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:XA:164:ARG:HD3	2:YA:195:LEU:HD11	2.01	0.41
1:ZA:164:ARG:HD3	2:AB:195:LEU:HD11	2.01	0.41
1:G:164:ARG:HD3	2:H:195:LEU:HD11	2.01	0.41
1:M:164:ARG:HD3	2:N:195:LEU:HD11	2.01	0.41
2:P:122:PHE:HB3	2:IB:78:ARG:HE	1.85	0.41
1:DA:160:ILE:CG2	1:DA:164:ARG:HH22	2.33	0.41
1:FA:164:ARG:HD3	2:GA:195:LEU:HD11	2.01	0.41
1:HA:164:ARG:HD3	2:IA:195:LEU:HD11	2.01	0.41
2:SA:138:SER:HA	2:SA:141:THR:HG23	2.01	0.41
2:UA:78:ARG:HE	2:QB:122:PHE:HB3	1.85	0.41
1:BB:160:ILE:CG2	1:BB:164:ARG:HH22	2.33	0.41
1:DB:160:ILE:CG2	1:DB:164:ARG:HH22	2.33	0.41
2:IB:138:SER:HA	2:IB:141:THR:HG23	2.01	0.41
1:TB:164:ARG:HD3	2:UB:195:LEU:HD11	2.01	0.41
1:A:74:LEU:HD21	1:ZA:50:GLU:HG2	2.02	0.41
1:A:164:ARG:HD3	2:B:195:LEU:HD11	2.01	0.41
2:B:138:SER:HA	2:B:141:THR:HG23	2.01	0.41
2:F:135:THR:HG21	2:UB:115:MET:HB3	2.02	0.41
1:K:78:LEU:HD23	1:LB:54:ILE:HD12	2.02	0.41
1:HA:160:ILE:CG2	1:HA:164:ARG:HH22	2.33	0.41
1:JA:164:ARG:HD3	2:KA:195:LEU:HD11	2.01	0.41
2:KA:138:SER:HA	2:KA:141:THR:HG23	2.01	0.41
1:ZA:160:ILE:CG2	1:ZA:164:ARG:HH22	2.33	0.41
1:S:160:ILE:CG2	1:S:164:ARG:HH22	2.33	0.41
1:C:160:ILE:CG2	1:C:164:ARG:HH22	2.33	0.41
1:I:164:ARG:HD3	2:J:195:LEU:HD11	2.01	0.41
2:W:121:GLN:HE22	2:EB:100:ASP:HB2	1.86	0.41
2:W:138:SER:HA	2:W:141:THR:HG23	2.01	0.41
1:X:164:ARG:HD3	2:Y:195:LEU:HD11	2.01	0.41
2:SA:191:ARG:HG2	2:SA:194:ARG:HH21	1.84	0.41
2:KB:138:SER:HA	2:KB:141:THR:HG23	2.01	0.41
1:I:160:ILE:CG2	1:I:164:ARG:HH22	2.33	0.41
2:W:191:ARG:HG2	2:W:194:ARG:HH21	1.84	0.41
1:I:78:LEU:HD23	1:Q:54:ILE:HD12	2.02	0.41
1:X:73:GLU:HG2	1:TA:51:ARG:NH1	2.36	0.41
2:AA:138:SER:HA	2:AA:141:THR:HG23	2.01	0.41
1:BA:160:ILE:CG2	1:BA:164:ARG:HH22	2.33	0.41
2:CA:138:SER:HA	2:CA:141:THR:HG23	2.01	0.41
1:NA:160:ILE:CG2	1:NA:164:ARG:HH22	2.33	0.41
1:PA:160:ILE:CG2	1:PA:164:ARG:HH22	2.33	0.41
1:TA:160:ILE:CG2	1:TA:164:ARG:HH22	2.33	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:TA:164:ARG:HD3	2:UA:195:LEU:HD11	2.01	0.41
2:IB:191:ARG:HG2	2:IB:194:ARG:HH21	1.84	0.41
2:OB:138:SER:HA	2:OB:141:THR:HG23	2.01	0.41
2:OA:138:SER:HA	2:OA:141:THR:HG23	2.01	0.41
1:V:160:ILE:CG2	1:V:164:ARG:HH22	2.33	0.40
2:OB:78:ARG:HE	2:UB:122:PHE:HB3	1.87	0.40
2:N:115:MET:HB3	2:KB:135:THR:HG21	2.03	0.40
1:Z:160:ILE:CG2	1:Z:164:ARG:HH22	2.33	0.40
1:RA:160:ILE:CG2	1:RA:164:ARG:HH22	2.33	0.40
2:T:75:VAL:HG21	2:GA:136:MET:HG2	2.02	0.40
2:GB:191:ARG:HG2	2:GB:194:ARG:HH21	1.84	0.40
1:HB:160:ILE:CG2	1:HB:164:ARG:HH22	2.33	0.40
2:GB:40:ILE:HG12	2:GB:45:MET:HE1	2.04	0.40
2:MB:40:ILE:HG12	2:MB:45:MET:HE1	2.04	0.40
2:MB:191:ARG:HG2	2:MB:194:ARG:HH21	1.84	0.40
1:NB:160:ILE:CG2	1:NB:164:ARG:HH22	2.33	0.40
2:T:40:ILE:HG12	2:T:45:MET:HE1	2.04	0.40
2:N:122:PHE:HB3	2:AA:78:ARG:HE	1.87	0.40
2:AB:40:ILE:HG12	2:AB:45:MET:HE1	2.04	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	163/366 (44%)	161 (99%)	2 (1%)	0	100	100
1	BA	163/366 (44%)	161 (99%)	2 (1%)	0	100	100
1	BB	163/366 (44%)	161 (99%)	2 (1%)	0	100	100
1	C	163/366 (44%)	161 (99%)	2 (1%)	0	100	100
1	DA	163/366 (44%)	161 (99%)	2 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	DB	163/366 (44%)	161 (99%)	2 (1%)	0	100	100
1	E	163/366 (44%)	161 (99%)	2 (1%)	0	100	100
1	FA	163/366 (44%)	161 (99%)	2 (1%)	0	100	100
1	FB	163/366 (44%)	161 (99%)	2 (1%)	0	100	100
1	G	163/366 (44%)	161 (99%)	2 (1%)	0	100	100
1	HA	163/366 (44%)	161 (99%)	2 (1%)	0	100	100
1	HB	163/366 (44%)	161 (99%)	2 (1%)	0	100	100
1	I	163/366 (44%)	161 (99%)	2 (1%)	0	100	100
1	JA	163/366 (44%)	161 (99%)	2 (1%)	0	100	100
1	JB	163/366 (44%)	161 (99%)	2 (1%)	0	100	100
1	K	163/366 (44%)	161 (99%)	2 (1%)	0	100	100
1	LA	163/366 (44%)	161 (99%)	2 (1%)	0	100	100
1	LB	163/366 (44%)	161 (99%)	2 (1%)	0	100	100
1	M	163/366 (44%)	161 (99%)	2 (1%)	0	100	100
1	NA	163/366 (44%)	161 (99%)	2 (1%)	0	100	100
1	NB	163/366 (44%)	161 (99%)	2 (1%)	0	100	100
1	O	163/366 (44%)	161 (99%)	2 (1%)	0	100	100
1	PA	163/366 (44%)	161 (99%)	2 (1%)	0	100	100
1	PB	163/366 (44%)	161 (99%)	2 (1%)	0	100	100
1	Q	163/366 (44%)	161 (99%)	2 (1%)	0	100	100
1	RA	163/366 (44%)	161 (99%)	2 (1%)	0	100	100
1	RB	163/366 (44%)	161 (99%)	2 (1%)	0	100	100
1	S	163/366 (44%)	161 (99%)	2 (1%)	0	100	100
1	TA	163/366 (44%)	161 (99%)	2 (1%)	0	100	100
1	TB	163/366 (44%)	161 (99%)	2 (1%)	0	100	100
1	V	163/366 (44%)	161 (99%)	2 (1%)	0	100	100
1	VA	163/366 (44%)	161 (99%)	2 (1%)	0	100	100
1	X	163/366 (44%)	161 (99%)	2 (1%)	0	100	100
1	XA	163/366 (44%)	161 (99%)	2 (1%)	0	100	100
1	Z	163/366 (44%)	161 (99%)	2 (1%)	0	100	100
1	ZA	163/366 (44%)	161 (99%)	2 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	AA	170/199 (85%)	169 (99%)	1 (1%)	0	100	100
2	AB	170/199 (85%)	169 (99%)	1 (1%)	0	100	100
2	B	170/199 (85%)	169 (99%)	1 (1%)	0	100	100
2	CA	170/199 (85%)	169 (99%)	1 (1%)	0	100	100
2	CB	170/199 (85%)	169 (99%)	1 (1%)	0	100	100
2	D	170/199 (85%)	169 (99%)	1 (1%)	0	100	100
2	EA	170/199 (85%)	169 (99%)	1 (1%)	0	100	100
2	EB	170/199 (85%)	169 (99%)	1 (1%)	0	100	100
2	F	170/199 (85%)	169 (99%)	1 (1%)	0	100	100
2	GA	170/199 (85%)	169 (99%)	1 (1%)	0	100	100
2	GB	170/199 (85%)	169 (99%)	1 (1%)	0	100	100
2	H	170/199 (85%)	169 (99%)	1 (1%)	0	100	100
2	IA	170/199 (85%)	169 (99%)	1 (1%)	0	100	100
2	IB	170/199 (85%)	169 (99%)	1 (1%)	0	100	100
2	J	170/199 (85%)	169 (99%)	1 (1%)	0	100	100
2	KA	170/199 (85%)	169 (99%)	1 (1%)	0	100	100
2	KB	170/199 (85%)	169 (99%)	1 (1%)	0	100	100
2	L	170/199 (85%)	169 (99%)	1 (1%)	0	100	100
2	MA	170/199 (85%)	169 (99%)	1 (1%)	0	100	100
2	MB	170/199 (85%)	169 (99%)	1 (1%)	0	100	100
2	N	170/199 (85%)	169 (99%)	1 (1%)	0	100	100
2	OA	170/199 (85%)	169 (99%)	1 (1%)	0	100	100
2	OB	170/199 (85%)	169 (99%)	1 (1%)	0	100	100
2	P	170/199 (85%)	169 (99%)	1 (1%)	0	100	100
2	QA	170/199 (85%)	169 (99%)	1 (1%)	0	100	100
2	QB	170/199 (85%)	169 (99%)	1 (1%)	0	100	100
2	R	170/199 (85%)	169 (99%)	1 (1%)	0	100	100
2	SA	170/199 (85%)	169 (99%)	1 (1%)	0	100	100
2	SB	170/199 (85%)	169 (99%)	1 (1%)	0	100	100
2	T	170/199 (85%)	169 (99%)	1 (1%)	0	100	100
2	UA	170/199 (85%)	169 (99%)	1 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	UB	170/199 (85%)	169 (99%)	1 (1%)	0	100	100
2	W	170/199 (85%)	169 (99%)	1 (1%)	0	100	100
2	WA	170/199 (85%)	169 (99%)	1 (1%)	0	100	100
2	Y	170/199 (85%)	169 (99%)	1 (1%)	0	100	100
2	YA	170/199 (85%)	169 (99%)	1 (1%)	0	100	100
All	All	11988/20340 (59%)	11880 (99%)	108 (1%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	144/309 (47%)	143 (99%)	1 (1%)	76	92
1	BA	144/309 (47%)	143 (99%)	1 (1%)	76	92
1	BB	144/309 (47%)	143 (99%)	1 (1%)	76	92
1	C	144/309 (47%)	143 (99%)	1 (1%)	76	92
1	DA	144/309 (47%)	143 (99%)	1 (1%)	76	92
1	DB	144/309 (47%)	143 (99%)	1 (1%)	76	92
1	E	144/309 (47%)	143 (99%)	1 (1%)	76	92
1	FA	144/309 (47%)	143 (99%)	1 (1%)	76	92
1	FB	144/309 (47%)	143 (99%)	1 (1%)	76	92
1	G	144/309 (47%)	143 (99%)	1 (1%)	76	92
1	HA	144/309 (47%)	143 (99%)	1 (1%)	76	92
1	HB	144/309 (47%)	143 (99%)	1 (1%)	76	92
1	I	144/309 (47%)	143 (99%)	1 (1%)	76	92
1	JA	144/309 (47%)	143 (99%)	1 (1%)	76	92
1	JB	144/309 (47%)	143 (99%)	1 (1%)	76	92
1	K	144/309 (47%)	143 (99%)	1 (1%)	76	92

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	LA	144/309 (47%)	143 (99%)	1 (1%)	76	92
1	LB	144/309 (47%)	143 (99%)	1 (1%)	76	92
1	M	144/309 (47%)	143 (99%)	1 (1%)	76	92
1	NA	144/309 (47%)	143 (99%)	1 (1%)	76	92
1	NB	144/309 (47%)	143 (99%)	1 (1%)	76	92
1	O	144/309 (47%)	143 (99%)	1 (1%)	76	92
1	PA	144/309 (47%)	143 (99%)	1 (1%)	76	92
1	PB	144/309 (47%)	143 (99%)	1 (1%)	76	92
1	Q	144/309 (47%)	143 (99%)	1 (1%)	76	92
1	RA	144/309 (47%)	143 (99%)	1 (1%)	76	92
1	RB	144/309 (47%)	143 (99%)	1 (1%)	76	92
1	S	144/309 (47%)	143 (99%)	1 (1%)	76	92
1	TA	144/309 (47%)	143 (99%)	1 (1%)	76	92
1	TB	144/309 (47%)	143 (99%)	1 (1%)	76	92
1	V	144/309 (47%)	143 (99%)	1 (1%)	76	92
1	VA	144/309 (47%)	143 (99%)	1 (1%)	76	92
1	X	144/309 (47%)	143 (99%)	1 (1%)	76	92
1	XA	144/309 (47%)	143 (99%)	1 (1%)	76	92
1	Z	144/309 (47%)	143 (99%)	1 (1%)	76	92
1	ZA	144/309 (47%)	143 (99%)	1 (1%)	76	92
2	AA	146/169 (86%)	146 (100%)	0	100	100
2	AB	146/169 (86%)	146 (100%)	0	100	100
2	B	146/169 (86%)	146 (100%)	0	100	100
2	CA	146/169 (86%)	146 (100%)	0	100	100
2	CB	146/169 (86%)	146 (100%)	0	100	100
2	D	146/169 (86%)	146 (100%)	0	100	100
2	EA	146/169 (86%)	146 (100%)	0	100	100
2	EB	146/169 (86%)	146 (100%)	0	100	100
2	F	146/169 (86%)	146 (100%)	0	100	100
2	GA	146/169 (86%)	146 (100%)	0	100	100
2	GB	146/169 (86%)	146 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	H	146/169 (86%)	146 (100%)	0	100	100
2	IA	146/169 (86%)	146 (100%)	0	100	100
2	IB	146/169 (86%)	146 (100%)	0	100	100
2	J	146/169 (86%)	146 (100%)	0	100	100
2	KA	146/169 (86%)	146 (100%)	0	100	100
2	KB	146/169 (86%)	146 (100%)	0	100	100
2	L	146/169 (86%)	146 (100%)	0	100	100
2	MA	146/169 (86%)	146 (100%)	0	100	100
2	MB	146/169 (86%)	146 (100%)	0	100	100
2	N	146/169 (86%)	146 (100%)	0	100	100
2	OA	146/169 (86%)	146 (100%)	0	100	100
2	OB	146/169 (86%)	146 (100%)	0	100	100
2	P	146/169 (86%)	146 (100%)	0	100	100
2	QA	146/169 (86%)	146 (100%)	0	100	100
2	QB	146/169 (86%)	146 (100%)	0	100	100
2	R	146/169 (86%)	146 (100%)	0	100	100
2	SA	146/169 (86%)	146 (100%)	0	100	100
2	SB	146/169 (86%)	146 (100%)	0	100	100
2	T	146/169 (86%)	146 (100%)	0	100	100
2	UA	146/169 (86%)	146 (100%)	0	100	100
2	UB	146/169 (86%)	146 (100%)	0	100	100
2	W	146/169 (86%)	146 (100%)	0	100	100
2	WA	146/169 (86%)	146 (100%)	0	100	100
2	Y	146/169 (86%)	146 (100%)	0	100	100
2	YA	146/169 (86%)	146 (100%)	0	100	100
All	All	10440/17208 (61%)	10404 (100%)	36 (0%)	84	96

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	S	165	TYR
1	A	165	TYR
1	C	165	TYR

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Mol	Chain	Res	Type
1	E	165	TYR
1	G	165	TYR
1	I	165	TYR
1	K	165	TYR
1	M	165	TYR
1	O	165	TYR
1	Q	165	TYR
1	V	165	TYR
1	X	165	TYR
1	Z	165	TYR
1	BA	165	TYR
1	DA	165	TYR
1	FA	165	TYR
1	HA	165	TYR
1	JA	165	TYR
1	LA	165	TYR
1	NA	165	TYR
1	PA	165	TYR
1	RA	165	TYR
1	TA	165	TYR
1	VA	165	TYR
1	XA	165	TYR
1	ZA	165	TYR
1	BB	165	TYR
1	DB	165	TYR
1	FB	165	TYR
1	HB	165	TYR
1	JB	165	TYR
1	LB	165	TYR
1	NB	165	TYR
1	PB	165	TYR
1	RB	165	TYR
1	TB	165	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (198) such sidechains are listed below:

Mol	Chain	Res	Type
1	S	14	ASN
2	T	44	ASN
2	T	58	GLN
2	T	61	GLN
2	T	121	GLN

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Mol	Chain	Res	Type
1	A	14	ASN
2	B	41	GLN
2	B	44	ASN
2	B	58	GLN
2	B	61	GLN
2	B	121	GLN
1	C	14	ASN
2	D	44	ASN
2	D	58	GLN
2	D	61	GLN
2	D	121	GLN
1	E	14	ASN
2	F	44	ASN
2	F	58	GLN
2	F	61	GLN
2	F	121	GLN
1	G	14	ASN
2	H	41	GLN
2	H	44	ASN
2	H	58	GLN
2	H	61	GLN
2	H	121	GLN
1	I	14	ASN
2	J	44	ASN
2	J	58	GLN
2	J	61	GLN
2	J	121	GLN
1	K	14	ASN
2	L	41	GLN
2	L	44	ASN
2	L	58	GLN
2	L	61	GLN
2	L	121	GLN
1	M	14	ASN
2	N	41	GLN
2	N	44	ASN
2	N	58	GLN
2	N	61	GLN
2	N	121	GLN
1	O	14	ASN
2	P	44	ASN
2	P	58	GLN

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Mol	Chain	Res	Type
2	P	61	GLN
2	P	121	GLN
1	Q	14	ASN
2	R	44	ASN
2	R	58	GLN
2	R	61	GLN
2	R	121	GLN
1	V	14	ASN
2	W	41	GLN
2	W	44	ASN
2	W	58	GLN
2	W	61	GLN
2	W	121	GLN
1	X	14	ASN
2	Y	41	GLN
2	Y	44	ASN
2	Y	58	GLN
2	Y	61	GLN
2	Y	121	GLN
1	Z	14	ASN
2	AA	41	GLN
2	AA	44	ASN
2	AA	58	GLN
2	AA	61	GLN
2	AA	121	GLN
1	BA	14	ASN
2	CA	44	ASN
2	CA	58	GLN
2	CA	61	GLN
2	CA	121	GLN
1	DA	14	ASN
2	EA	41	GLN
2	EA	44	ASN
2	EA	58	GLN
2	EA	61	GLN
2	EA	121	GLN
1	FA	14	ASN
2	GA	44	ASN
2	GA	58	GLN
2	GA	61	GLN
2	GA	121	GLN
1	HA	14	ASN

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Mol	Chain	Res	Type
2	IA	41	GLN
2	IA	44	ASN
2	IA	58	GLN
2	IA	61	GLN
2	IA	121	GLN
1	JA	14	ASN
2	KA	41	GLN
2	KA	44	ASN
2	KA	58	GLN
2	KA	61	GLN
2	KA	121	GLN
1	LA	14	ASN
2	MA	44	ASN
2	MA	58	GLN
2	MA	61	GLN
2	MA	121	GLN
1	NA	14	ASN
2	OA	44	ASN
2	OA	58	GLN
2	OA	61	GLN
2	OA	121	GLN
1	PA	14	ASN
2	QA	44	ASN
2	QA	58	GLN
2	QA	61	GLN
2	QA	121	GLN
1	RA	14	ASN
2	SA	44	ASN
2	SA	58	GLN
2	SA	61	GLN
2	SA	121	GLN
1	TA	14	ASN
2	UA	41	GLN
2	UA	44	ASN
2	UA	58	GLN
2	UA	61	GLN
2	UA	121	GLN
1	VA	14	ASN
2	WA	41	GLN
2	WA	44	ASN
2	WA	58	GLN
2	WA	61	GLN

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Mol	Chain	Res	Type
2	WA	107	ASN
1	XA	14	ASN
2	YA	41	GLN
2	YA	44	ASN
2	YA	58	GLN
2	YA	61	GLN
2	YA	121	GLN
1	ZA	14	ASN
2	AB	44	ASN
2	AB	58	GLN
2	AB	61	GLN
2	AB	121	GLN
1	BB	14	ASN
2	CB	44	ASN
2	CB	58	GLN
2	CB	61	GLN
2	CB	121	GLN
1	DB	14	ASN
2	EB	44	ASN
2	EB	58	GLN
2	EB	61	GLN
2	EB	121	GLN
1	FB	14	ASN
2	GB	41	GLN
2	GB	44	ASN
2	GB	58	GLN
2	GB	61	GLN
2	GB	121	GLN
1	HB	14	ASN
2	IB	41	GLN
2	IB	44	ASN
2	IB	58	GLN
2	IB	61	GLN
2	IB	121	GLN
1	JB	14	ASN
2	KB	41	GLN
2	KB	44	ASN
2	KB	58	GLN
2	KB	61	GLN
2	KB	121	GLN
1	LB	14	ASN
2	MB	44	ASN

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Mol	Chain	Res	Type
2	MB	58	GLN
2	MB	61	GLN
2	MB	121	GLN
1	NB	14	ASN
2	OB	44	ASN
2	OB	58	GLN
2	OB	61	GLN
2	OB	121	GLN
1	PB	14	ASN
2	QB	41	GLN
2	QB	44	ASN
2	QB	58	GLN
2	QB	61	GLN
2	QB	121	GLN
1	RB	14	ASN
2	SB	41	GLN
2	SB	44	ASN
2	SB	58	GLN
2	SB	61	GLN
2	SB	121	GLN
1	TB	14	ASN
2	UB	44	ASN
2	UB	58	GLN
2	UB	61	GLN
2	UB	121	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

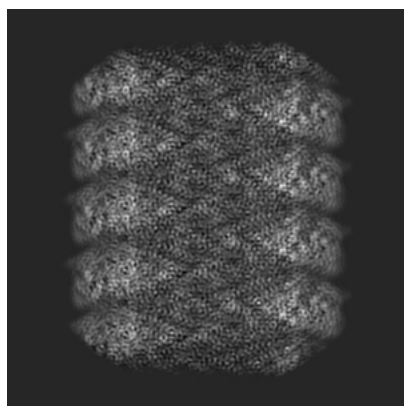
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-9005. These allow visual inspection of the internal detail of the map and identification of artifacts.

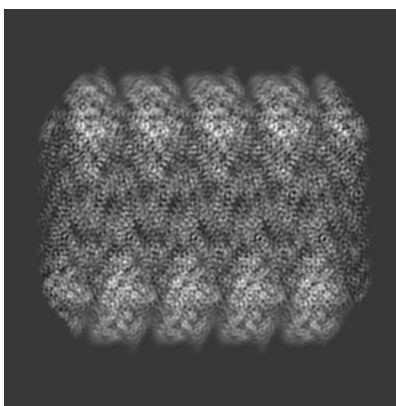
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

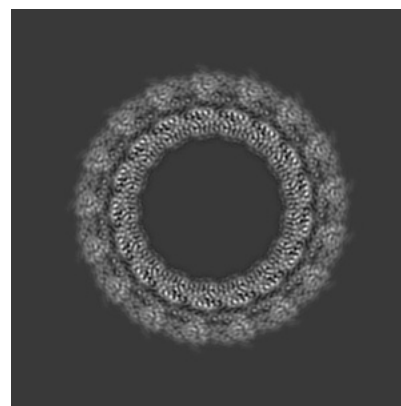
6.1.1 Primary map



X

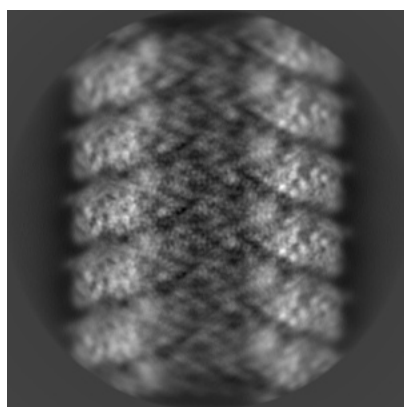


Y

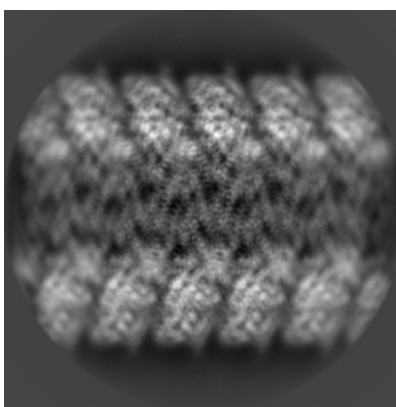


Z

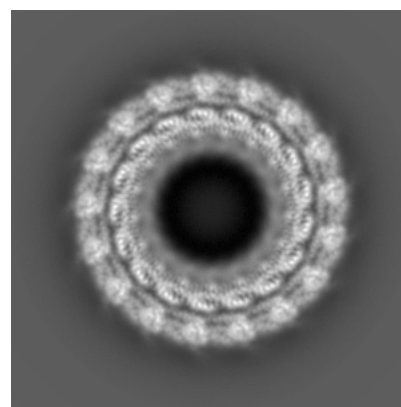
6.1.2 Raw map



X



Y

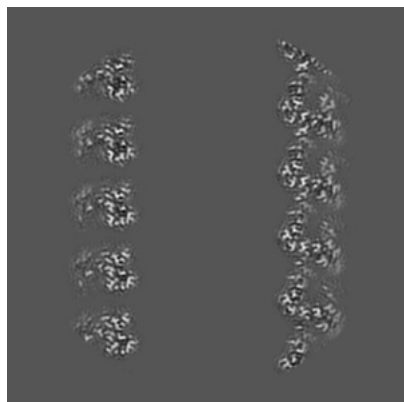


Z

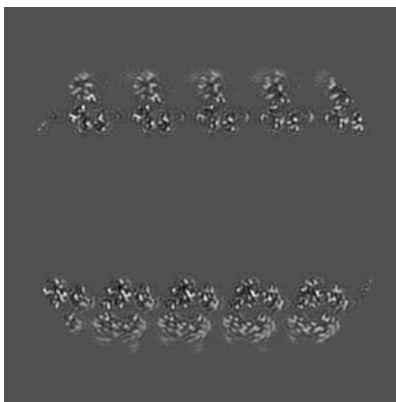
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

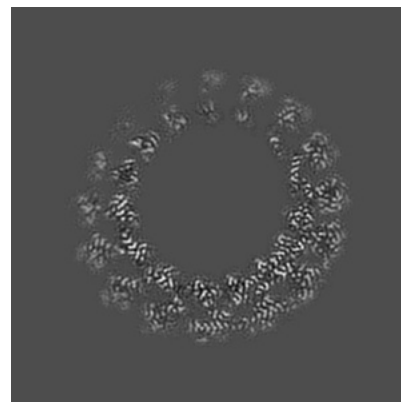
6.2.1 Primary map



X Index: 200

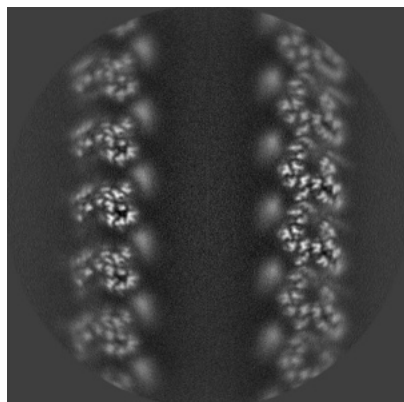


Y Index: 200

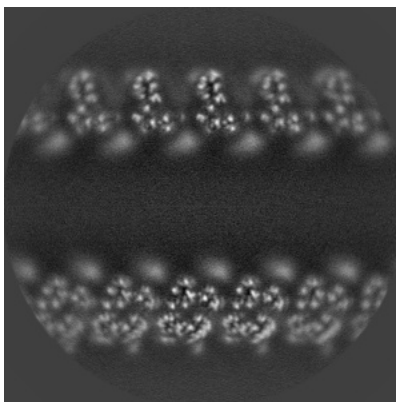


Z Index: 200

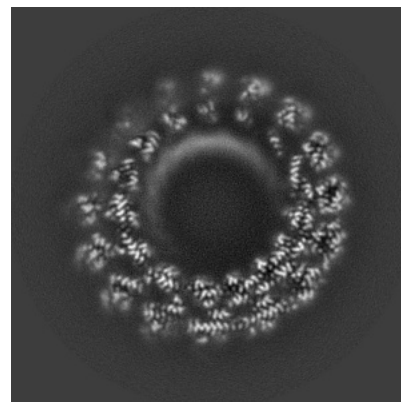
6.2.2 Raw map



X Index: 200



Y Index: 200

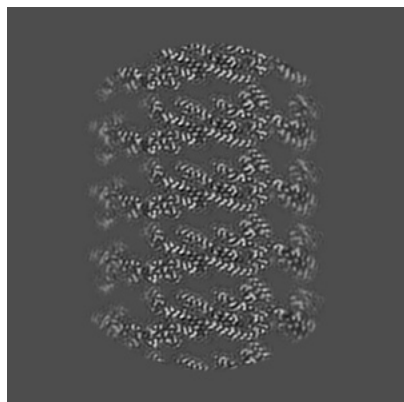


Z Index: 200

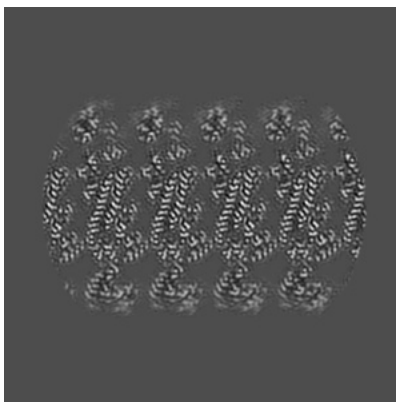
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

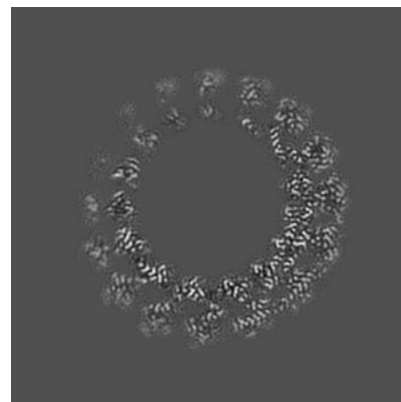
6.3.1 Primary map



X Index: 282

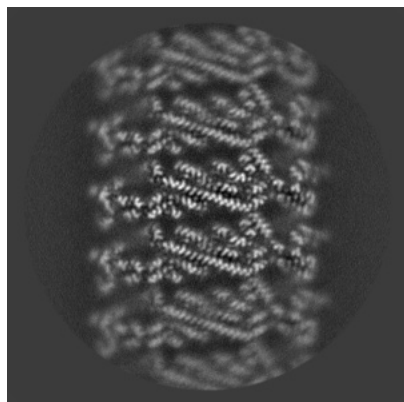


Y Index: 286

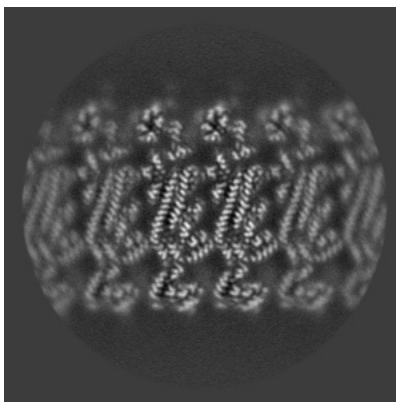


Z Index: 204

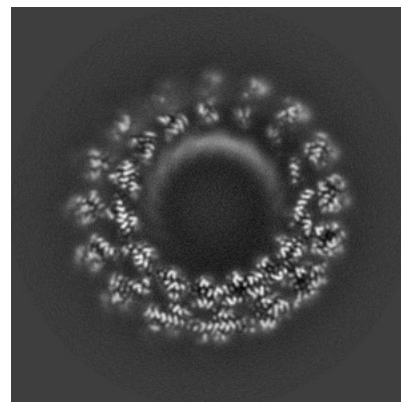
6.3.2 Raw map



X Index: 283



Y Index: 286

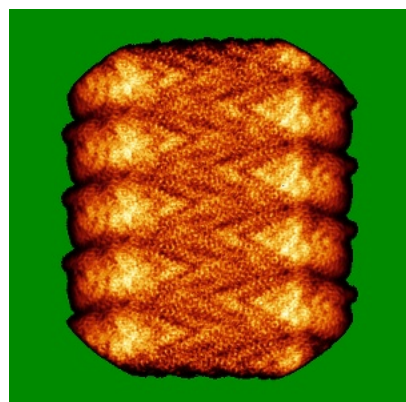


Z Index: 197

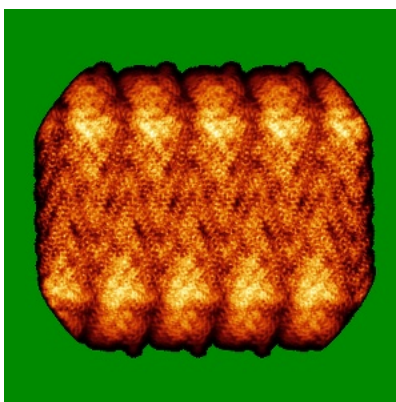
The images above show the largest variance slices of the map in three orthogonal directions.

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

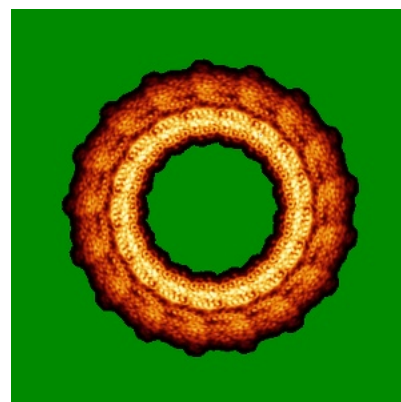
6.4.1 Primary map



X

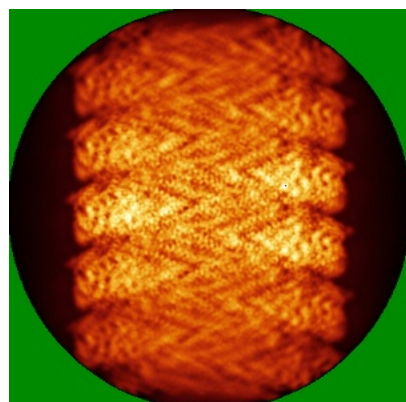


Y

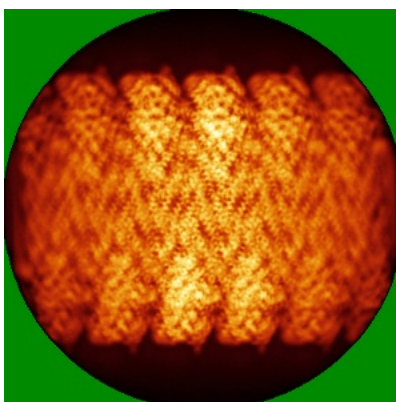


Z

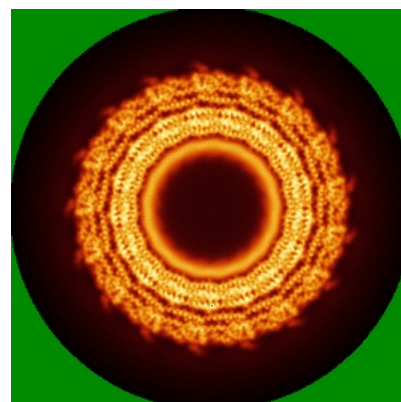
6.4.2 Raw map



X



Y

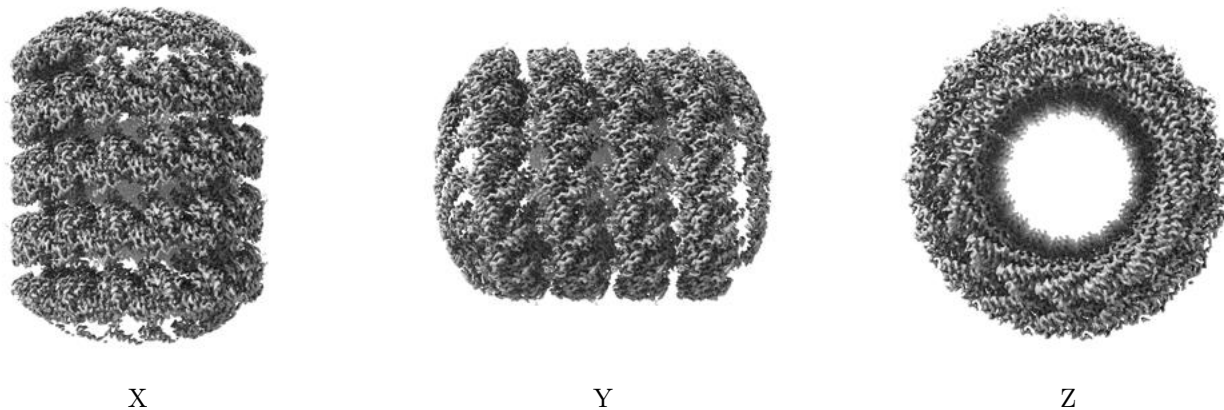


Z

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

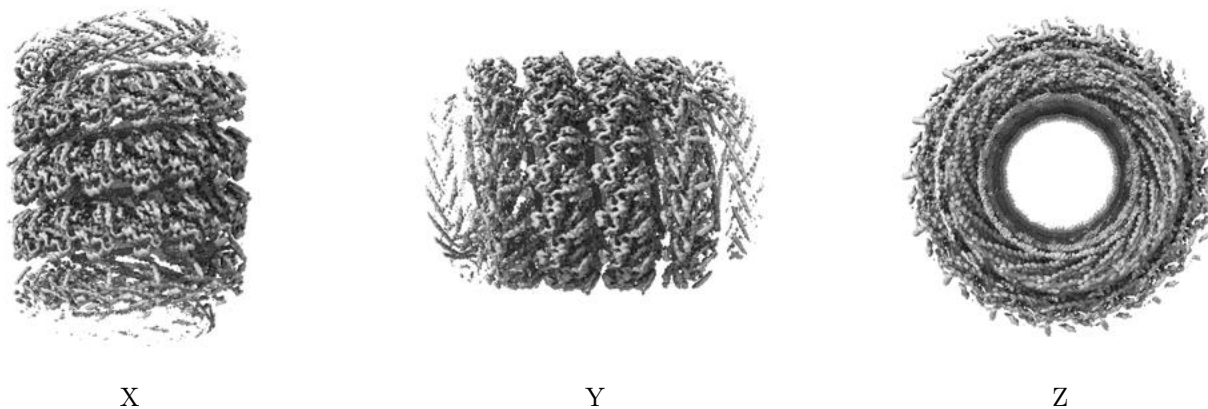
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.025. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



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

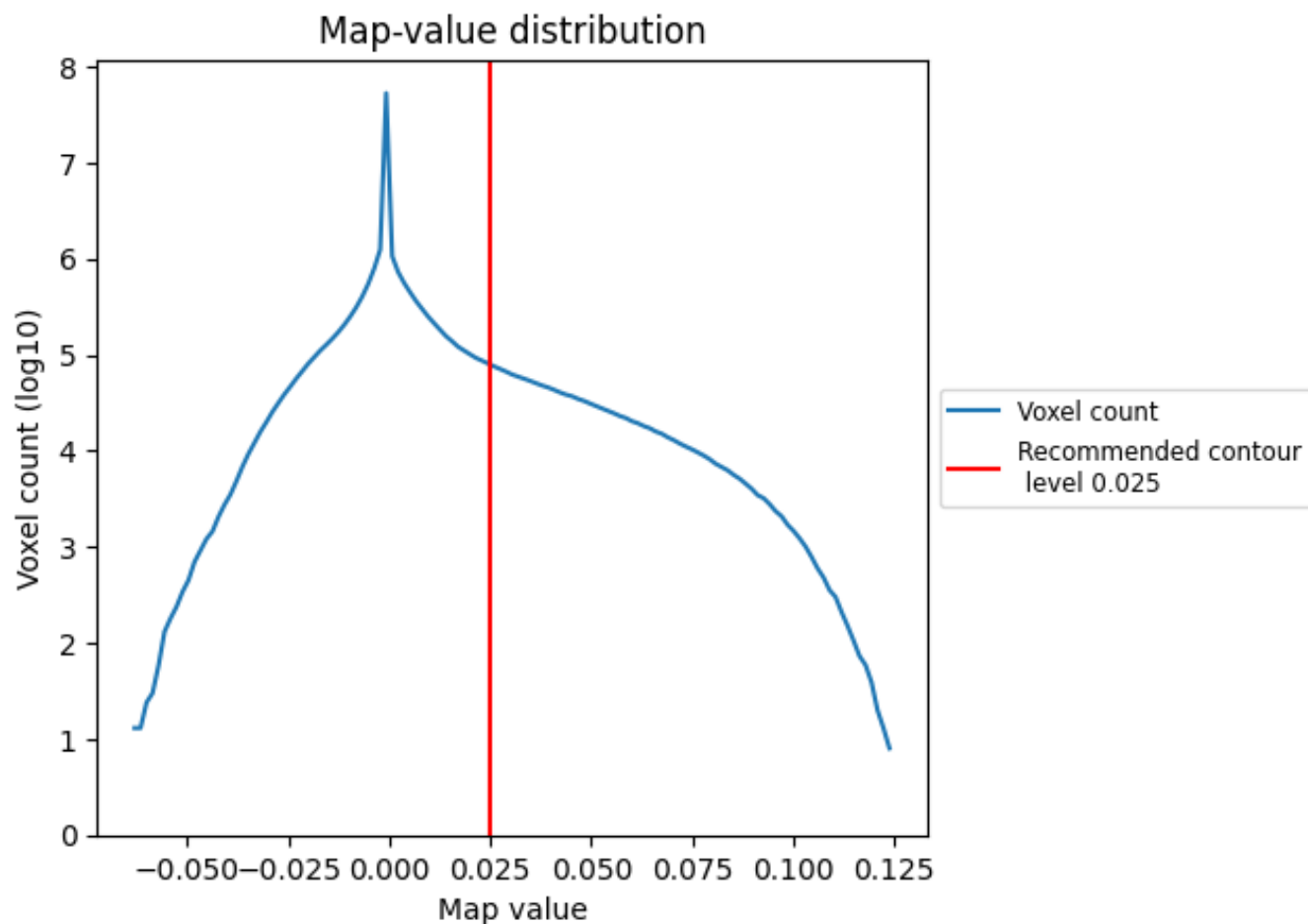
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

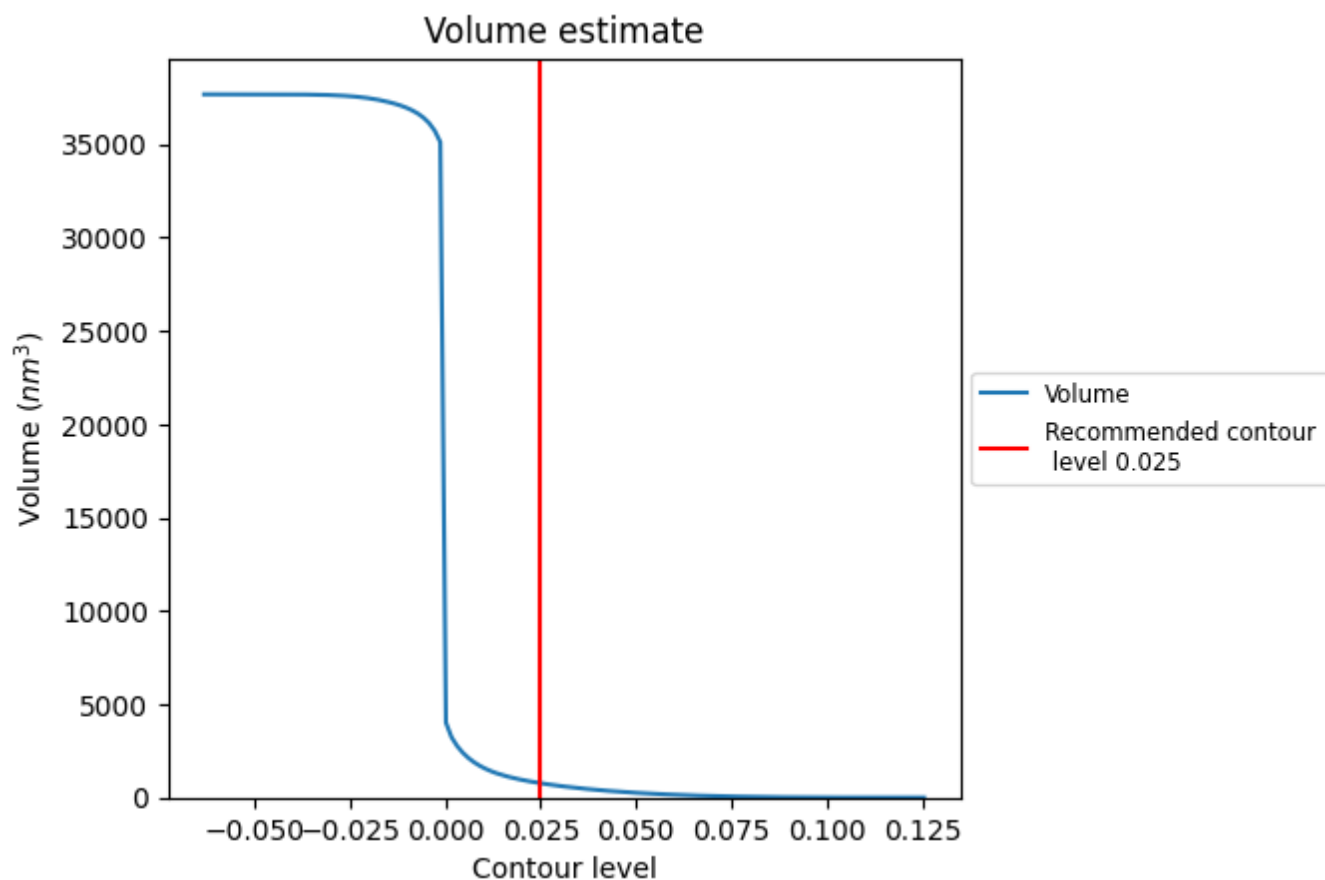
This section contains the results of statistical analysis of the map.

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.

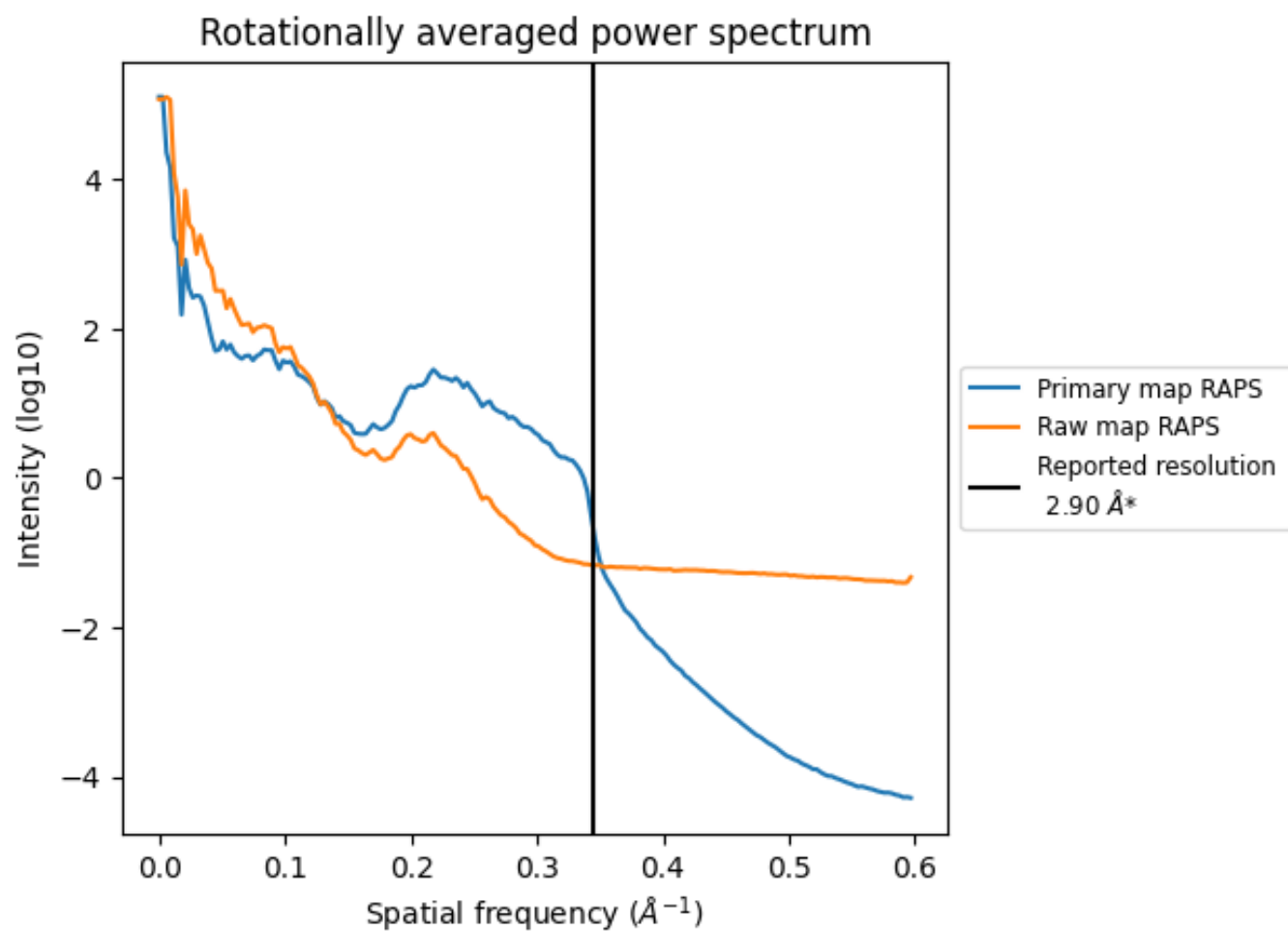
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 769 nm³; this corresponds to an approximate mass of 695 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 ⓘ

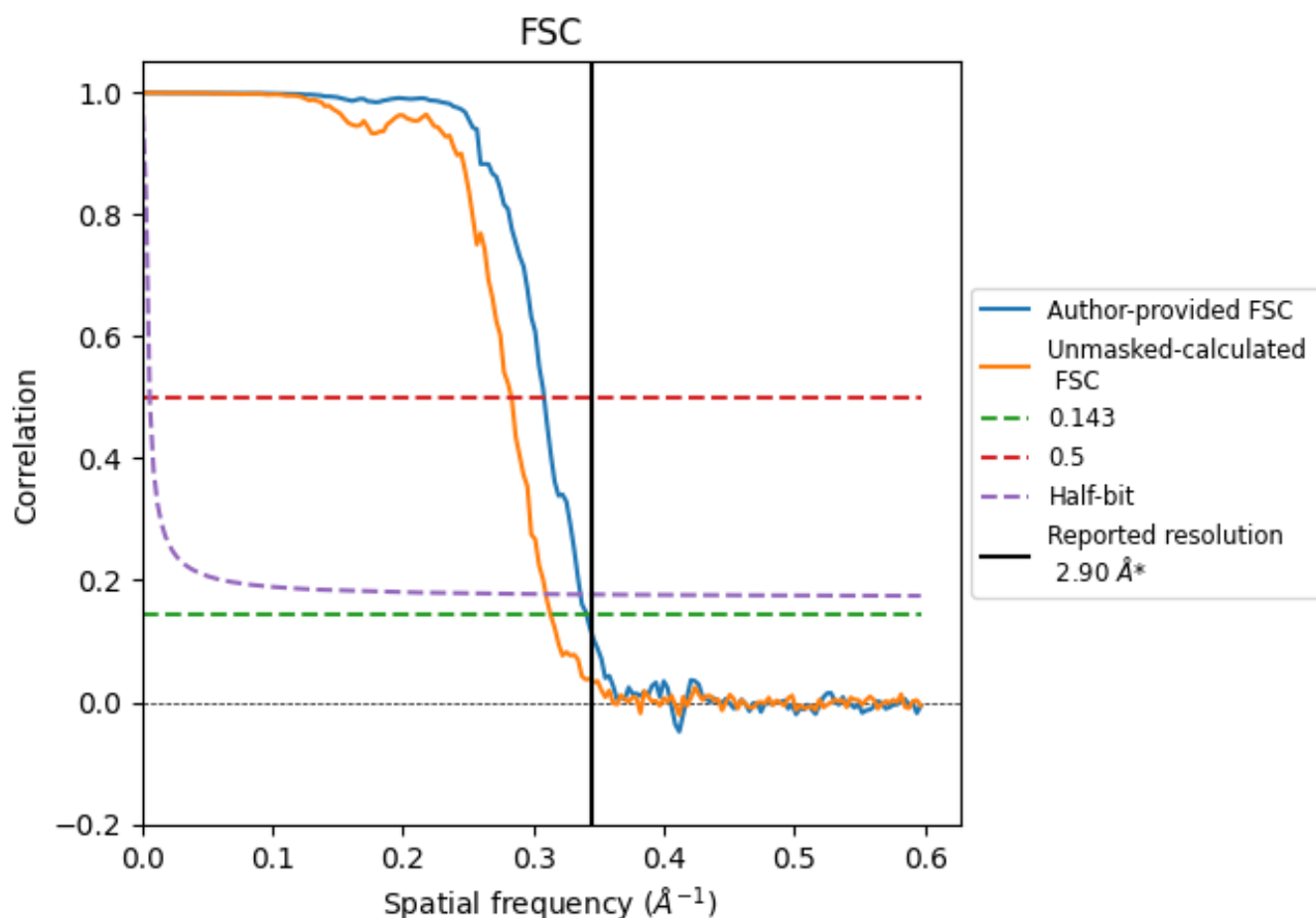


*Reported resolution corresponds to spatial frequency of 0.345 Å⁻¹

8 Fourier-Shell correlation [i](#)

Fourier-Shell Correlation (FSC) is the most commonly used method to estimate the resolution of single-particle and subtomogram-averaged maps. The shape of the curve depends on the imposed symmetry, mask and whether or not the two 3D reconstructions used were processed from a common reference. The reported resolution is shown as a black line. A curve is displayed for the half-bit criterion in addition to lines showing the 0.143 gold standard cut-off and 0.5 cut-off.

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.345 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

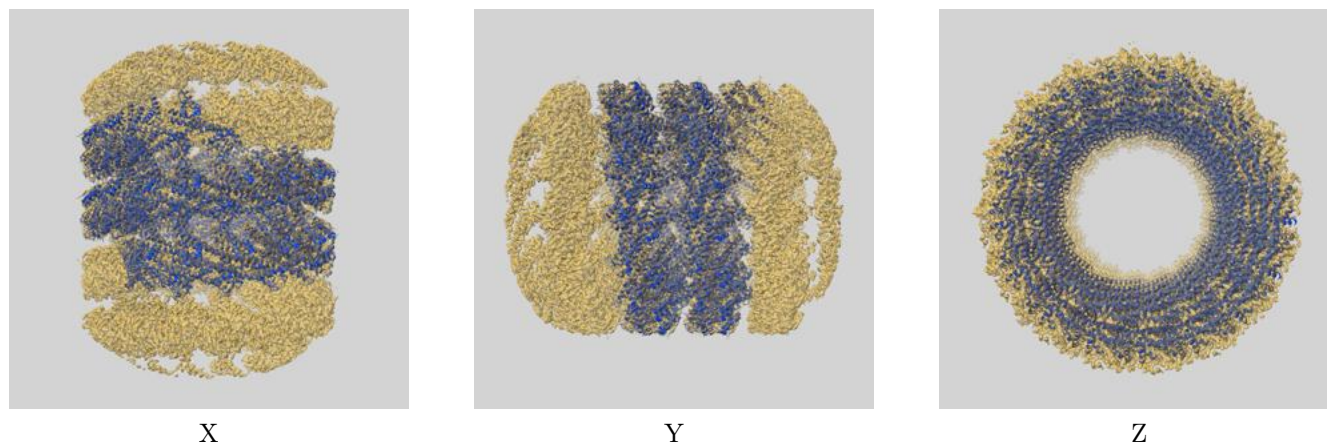
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.90	-	-
Author-provided FSC curve	2.94	3.25	2.98
Unmasked-calculated*	3.19	3.53	3.23

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 3.19 differs from the reported value 2.9 by more than 10 %

9 Map-model fit [i](#)

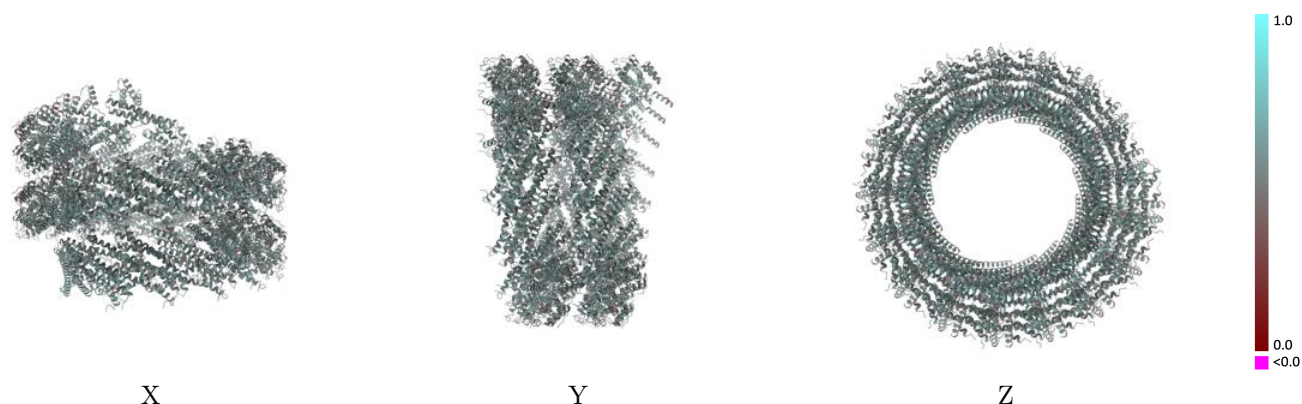
This section contains information regarding the fit between EMDB map EMD-9005 and PDB model 6E8G. Per-residue inclusion information can be found in section 3 on page 11.

9.1 Map-model overlay [i](#)



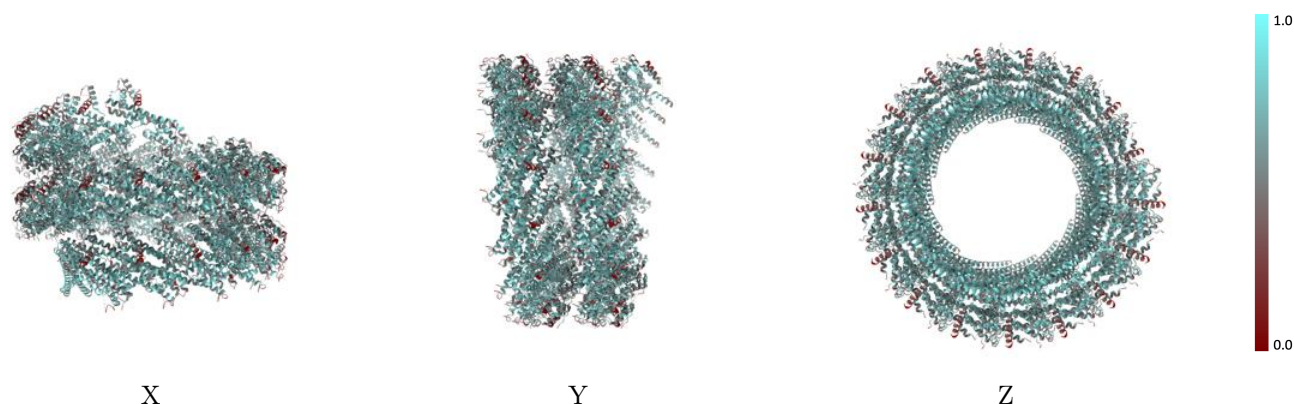
The images above show the 3D surface view of the map at the recommended contour level 0.025 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



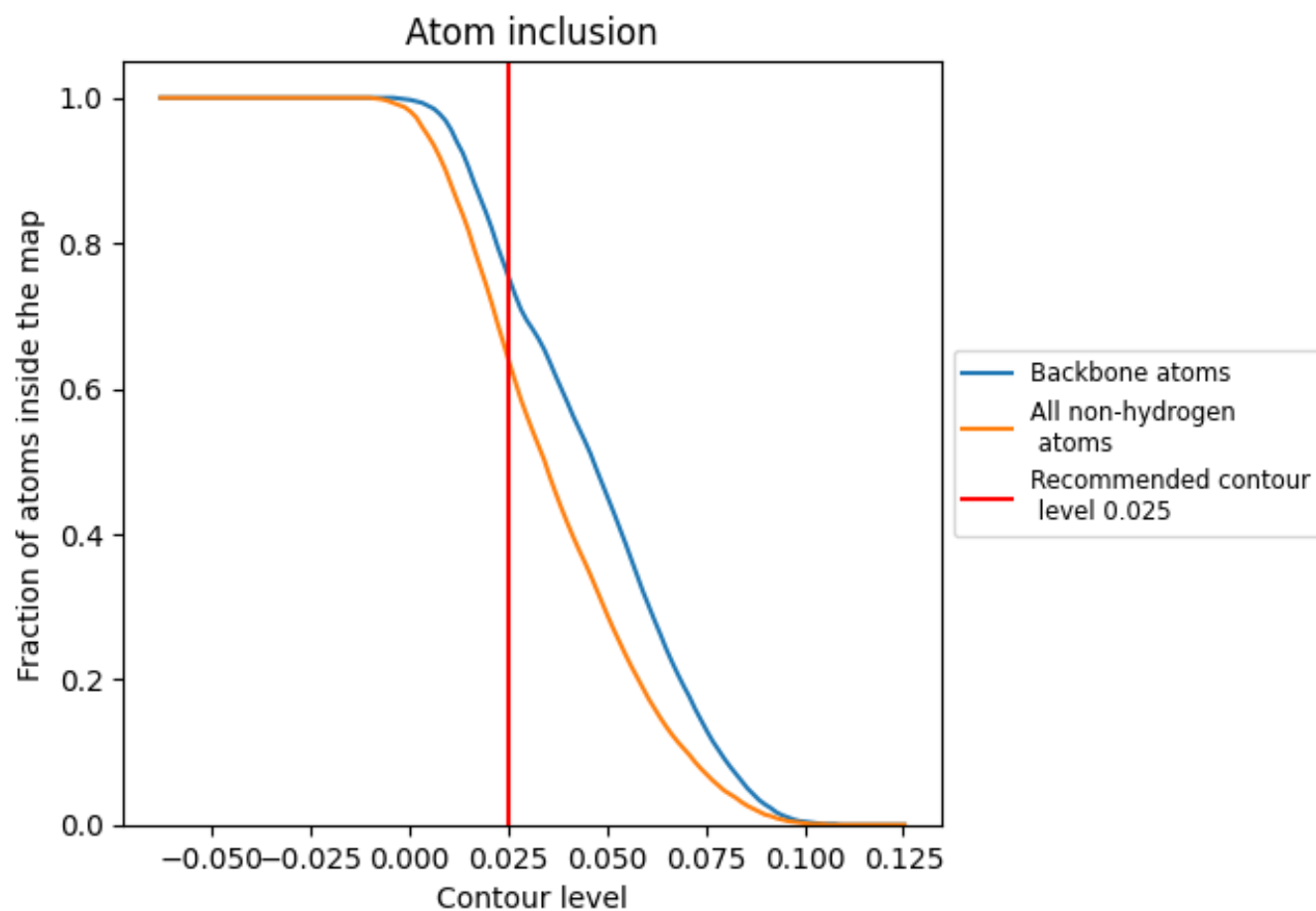
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.025).

9.4 Atom inclusion [i](#)



At the recommended contour level, 76% of all backbone atoms, 64% 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 (0.025) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.6400	 0.5370
A	 0.6160	 0.5320
AA	 0.6640	 0.5450
AB	 0.6700	 0.5470
B	 0.6680	 0.5460
BA	 0.6230	 0.5310
BB	 0.6190	 0.5290
C	 0.6190	 0.5310
CA	 0.6640	 0.5450
CB	 0.6580	 0.5420
D	 0.6610	 0.5480
DA	 0.6200	 0.5300
DB	 0.6210	 0.5250
E	 0.6130	 0.5300
EA	 0.6640	 0.5460
EB	 0.6700	 0.5420
F	 0.6610	 0.5450
FA	 0.6150	 0.5310
FB	 0.6100	 0.5230
G	 0.6200	 0.5310
GA	 0.6610	 0.5440
GB	 0.6580	 0.5410
H	 0.6670	 0.5500
HA	 0.6130	 0.5300
HB	 0.6200	 0.5300
I	 0.6160	 0.5310
IA	 0.6670	 0.5480
IB	 0.6590	 0.5450
J	 0.6710	 0.5480
JA	 0.6160	 0.5310
JB	 0.6200	 0.5300
K	 0.6140	 0.5310
KA	 0.6660	 0.5460
KB	 0.6550	 0.5440
L	 0.6610	 0.5450



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Chain	Atom inclusion	Q-score
LA	 0.6180	 0.5330
LB	 0.6200	 0.5310
M	 0.6130	 0.5310
MA	 0.6670	 0.5480
MB	 0.6620	 0.5450
N	 0.6530	 0.5430
NA	 0.6130	 0.5210
NB	 0.6170	 0.5300
O	 0.6190	 0.5310
OA	 0.6540	 0.5410
OB	 0.6610	 0.5470
P	 0.6610	 0.5440
PA	 0.6170	 0.5310
PB	 0.6120	 0.5280
Q	 0.6120	 0.5290
QA	 0.6650	 0.5470
QB	 0.6620	 0.5470
R	 0.6640	 0.5470
RA	 0.6170	 0.5320
RB	 0.6230	 0.5300
S	 0.6160	 0.5300
SA	 0.6640	 0.5440
SB	 0.6640	 0.5450
T	 0.6670	 0.5470
TA	 0.6170	 0.5300
TB	 0.6160	 0.5310
UA	 0.6690	 0.5470
UB	 0.6570	 0.5430
V	 0.6180	 0.5230
VA	 0.6170	 0.5300
W	 0.6560	 0.5380
WA	 0.6620	 0.5440
X	 0.6120	 0.5320
XA	 0.6190	 0.5280
Y	 0.6670	 0.5500
YA	 0.6620	 0.5470
Z	 0.6210	 0.5290
ZA	 0.6160	 0.5310