



wwPDB EM Validation Summary Report ⓘ

Mar 28, 2026 – 04:20 PM UTC

PDB ID : 5TCR / pdb_00005tcr
EMDB ID : EMD-8400
Title : Atomic model of the Salmonella SPI-1 type III secretion injectisome basal body proteins InvG, PrgH, and PrgK
Authors : Worrall, L.J.; Hong, C.; Vuckovic, M.; Bergeron, J.R.C.; Huang, R.K.; Yu, Z.; Strynadka, N.C.J.
Deposited on : 2016-09-15
Resolution : 6.30 Å(reported)

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

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

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:

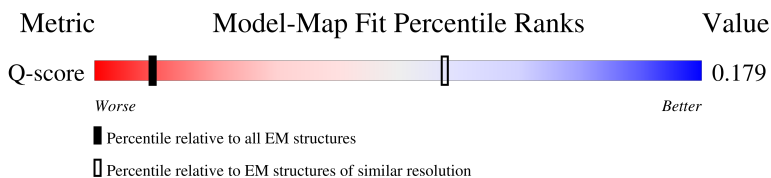
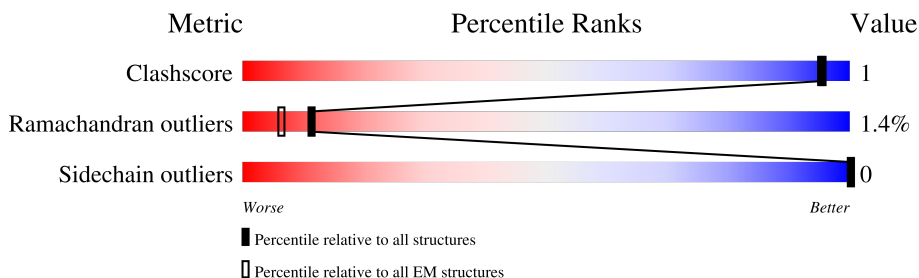
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDb archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & 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 6.30 Å.

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



Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	550 (5.80 - 6.80)

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

Mol	Chain	Length	Quality of chain
1	A	562	33% 78% 6% 16%
1	B	562	34% 78% 6% 16%
1	C	562	38% 78% 6% 16%
1	D	562	37% 78% 6% 16%

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Mol	Chain	Length	Quality of chain
1	E	562	36% 78% 6% 16%
1	F	562	36% 78% 6% 16%
1	G	562	34% 78% 6% 16%
1	H	562	36% 78% 6% 16%
1	I	562	37% 78% 6% 16%
1	J	562	36% 78% 6% 16%
1	K	562	36% 78% 6% 16%
1	L	562	40% 78% 6% 16%
1	M	562	38% 78% 6% 16%
1	N	562	35% 78% 6% 16%
1	O	562	34% 78% 6% 16%
2	1	235	73% 5% 22%
2	3	235	73% 5% 22%
2	5	235	73% 5% 22%
2	7	235	72% 5% 22%
2	9	235	72% 5% 22%
2	P	235	72% 5% 22%
2	R	235	73% 5% 22%
2	T	235	72% 5% 22%
2	V	235	5% 72% 5% 22%
2	X	235	73% 5% 22%
2	a	235	72% 5% 22%
2	c	235	5% 72% 5% 22%
2	e	235	7% 72% 5% 22%
2	g	235	5% 72% 5% 22%

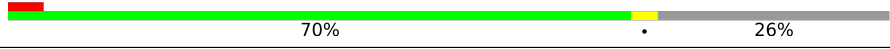
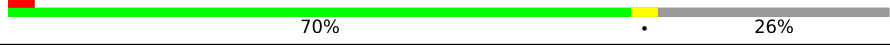
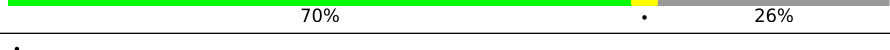
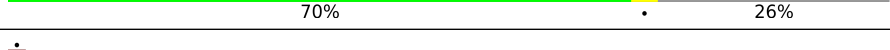
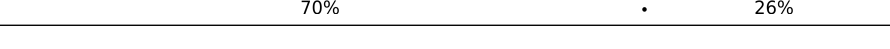
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Mol	Chain	Length	Quality of chain
2	i	235	
2	k	235	
2	m	235	
2	o	235	
2	q	235	
2	s	235	
2	u	235	
2	w	235	
2	y	235	
2	z	235	
3	0	263	
3	10	263	
3	2	263	
3	4	263	
3	6	263	
3	8	263	
3	Q	263	
3	S	263	
3	U	263	
3	W	263	
3	Y	263	
3	Z	263	
3	b	263	
3	d	263	
3	f	263	

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Mol	Chain	Length	Quality of chain
3	h	263	 70%26%
3	j	263	 70%26%
3	l	263	 70%26%
3	n	263	 70%26%
3	p	263	 70%26%
3	r	263	 70%26%
3	t	263	 70%26%
3	v	263	 70%26%
3	x	263	 70%26%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 128658 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 Protein InvG.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	474	Total	C	N	O	S	0	0
			3718	2357	645	704	12		
1	B	474	Total	C	N	O	S	0	0
			3718	2357	645	704	12		
1	C	474	Total	C	N	O	S	0	0
			3718	2357	645	704	12		
1	D	474	Total	C	N	O	S	0	0
			3718	2357	645	704	12		
1	E	474	Total	C	N	O	S	0	0
			3718	2357	645	704	12		
1	F	474	Total	C	N	O	S	0	0
			3718	2357	645	704	12		
1	G	474	Total	C	N	O	S	0	0
			3718	2357	645	704	12		
1	H	474	Total	C	N	O	S	0	0
			3718	2357	645	704	12		
1	I	474	Total	C	N	O	S	0	0
			3718	2357	645	704	12		
1	J	474	Total	C	N	O	S	0	0
			3718	2357	645	704	12		
1	K	474	Total	C	N	O	S	0	0
			3718	2357	645	704	12		
1	L	474	Total	C	N	O	S	0	0
			3718	2357	645	704	12		
1	M	474	Total	C	N	O	S	0	0
			3718	2357	645	704	12		
1	N	474	Total	C	N	O	S	0	0
			3718	2357	645	704	12		
1	O	474	Total	C	N	O	S	0	0
			3718	2357	645	704	12		

- Molecule 2 is a protein called Lipoprotein PrgK.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	P	184	Total 1437	C 905	N 250	O 279	S 3	0	0
2	R	184	Total 1437	C 905	N 250	O 279	S 3	0	0
2	T	184	Total 1437	C 905	N 250	O 279	S 3	0	0
2	V	184	Total 1437	C 905	N 250	O 279	S 3	0	0
2	X	184	Total 1437	C 905	N 250	O 279	S 3	0	0
2	a	184	Total 1437	C 905	N 250	O 279	S 3	0	0
2	c	184	Total 1437	C 905	N 250	O 279	S 3	0	0
2	e	184	Total 1437	C 905	N 250	O 279	S 3	0	0
2	g	184	Total 1437	C 905	N 250	O 279	S 3	0	0
2	i	184	Total 1437	C 905	N 250	O 279	S 3	0	0
2	k	184	Total 1437	C 905	N 250	O 279	S 3	0	0
2	m	184	Total 1437	C 905	N 250	O 279	S 3	0	0
2	o	184	Total 1437	C 905	N 250	O 279	S 3	0	0
2	q	184	Total 1437	C 905	N 250	O 279	S 3	0	0
2	s	184	Total 1437	C 905	N 250	O 279	S 3	0	0
2	u	184	Total 1437	C 905	N 250	O 279	S 3	0	0
2	w	184	Total 1437	C 905	N 250	O 279	S 3	0	0
2	y	184	Total 1437	C 905	N 250	O 279	S 3	0	0
2	z	184	Total 1437	C 905	N 250	O 279	S 3	0	0
2	1	184	Total 1437	C 905	N 250	O 279	S 3	0	0
2	3	184	Total 1437	C 905	N 250	O 279	S 3	0	0
2	5	184	Total 1437	C 905	N 250	O 279	S 3	0	0

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Mol	Chain	Residues	Atoms					AltConf	Trace
2	7	184	Total	C	N	O	S	0	0
			1437	905	250	279	3		
2	9	184	Total	C	N	O	S	0	0
			1437	905	250	279	3		

- Molecule 3 is a protein called Protein PrgH.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	Q	194	Total	C	N	O	S	0	0
			1600	1011	288	297	4		
3	S	194	Total	C	N	O	S	0	0
			1600	1011	288	297	4		
3	U	194	Total	C	N	O	S	0	0
			1600	1011	288	297	4		
3	W	194	Total	C	N	O	S	0	0
			1600	1011	288	297	4		
3	Y	194	Total	C	N	O	S	0	0
			1600	1011	288	297	4		
3	Z	194	Total	C	N	O	S	0	0
			1600	1011	288	297	4		
3	b	194	Total	C	N	O	S	0	0
			1600	1011	288	297	4		
3	d	194	Total	C	N	O	S	0	0
			1600	1011	288	297	4		
3	f	194	Total	C	N	O	S	0	0
			1600	1011	288	297	4		
3	h	194	Total	C	N	O	S	0	0
			1600	1011	288	297	4		
3	j	194	Total	C	N	O	S	0	0
			1600	1011	288	297	4		
3	l	194	Total	C	N	O	S	0	0
			1600	1011	288	297	4		
3	n	194	Total	C	N	O	S	0	0
			1600	1011	288	297	4		
3	p	194	Total	C	N	O	S	0	0
			1600	1011	288	297	4		
3	r	194	Total	C	N	O	S	0	0
			1600	1011	288	297	4		
3	t	194	Total	C	N	O	S	0	0
			1600	1011	288	297	4		
3	v	194	Total	C	N	O	S	0	0
			1600	1011	288	297	4		

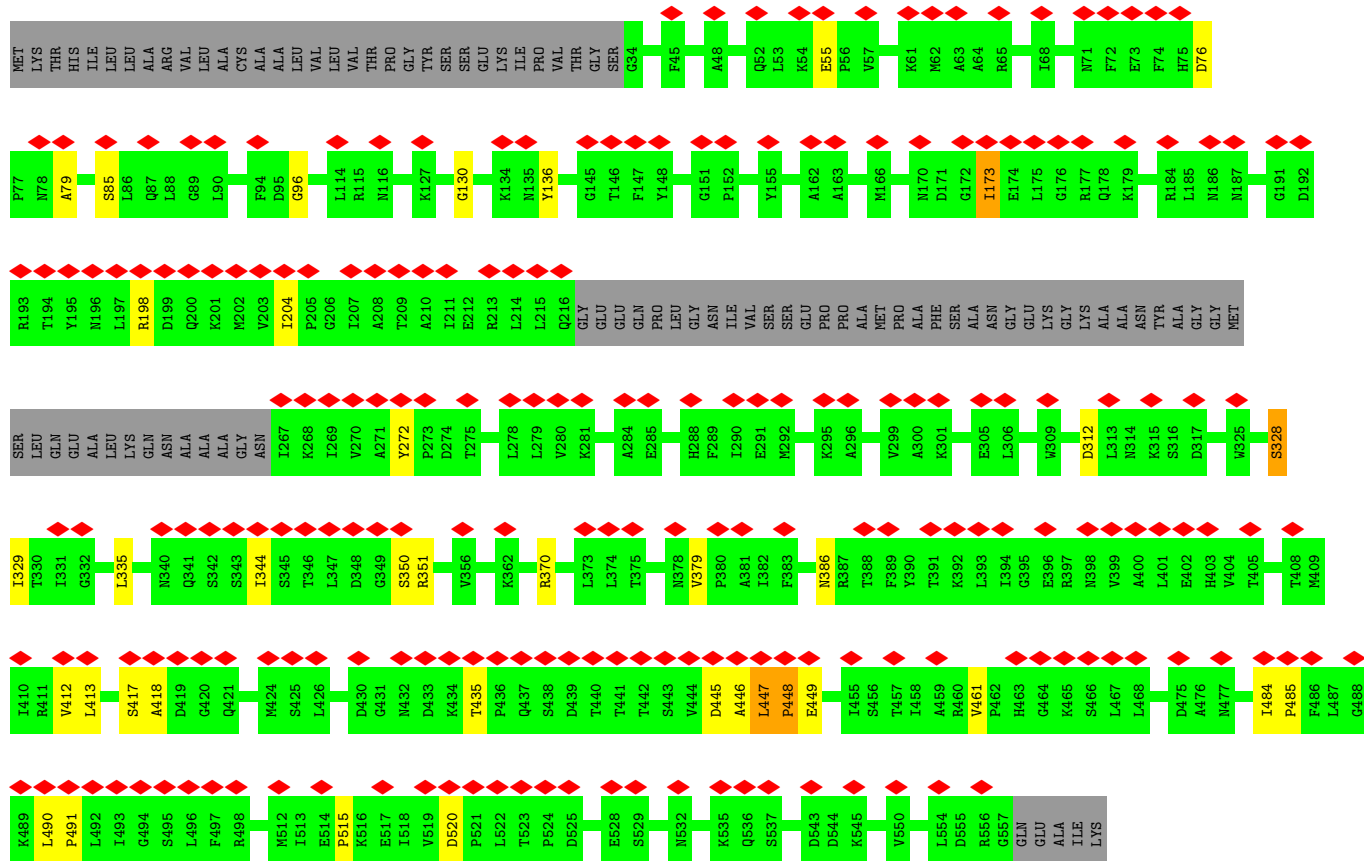
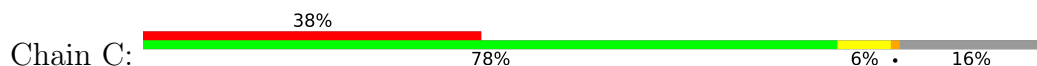
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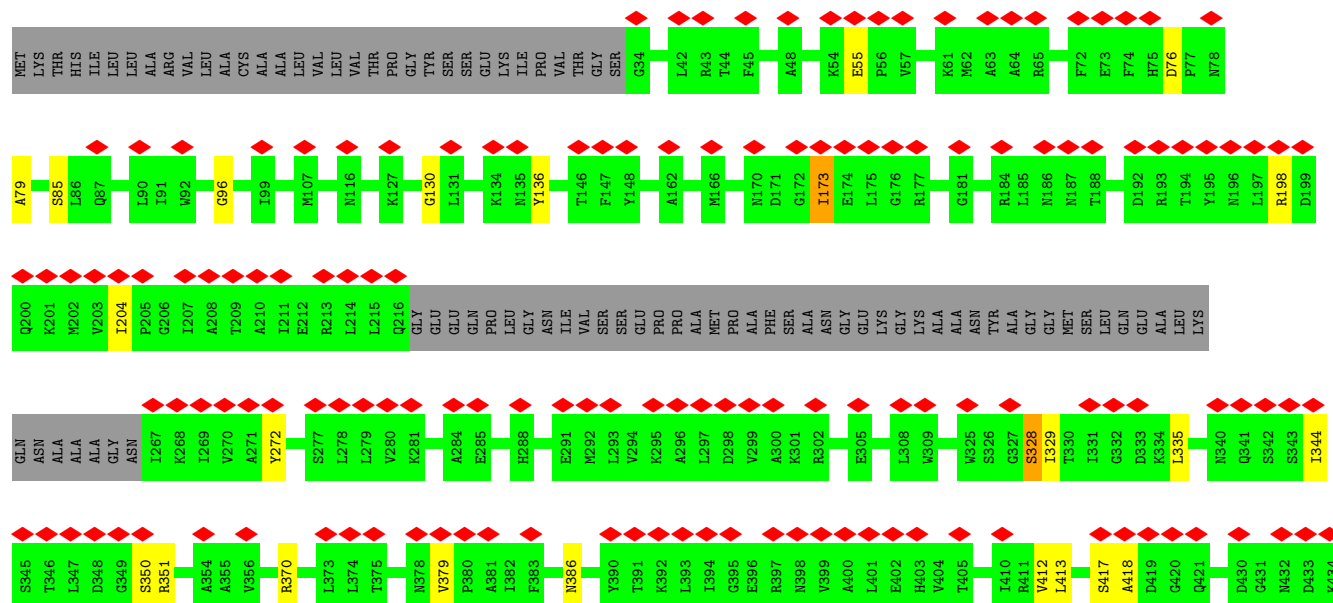
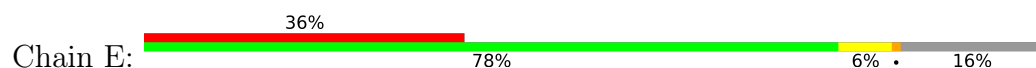
Mol	Chain	Residues	Atoms					AltConf	Trace
3	x	194	Total 1600	C 1011	N 288	O 297	S 4	0	0
3	0	194	Total 1600	C 1011	N 288	O 297	S 4	0	0
3	2	194	Total 1600	C 1011	N 288	O 297	S 4	0	0
3	4	194	Total 1600	C 1011	N 288	O 297	S 4	0	0
3	6	194	Total 1600	C 1011	N 288	O 297	S 4	0	0
3	8	194	Total 1600	C 1011	N 288	O 297	S 4	0	0
3	10	194	Total 1600	C 1011	N 288	O 297	S 4	0	0

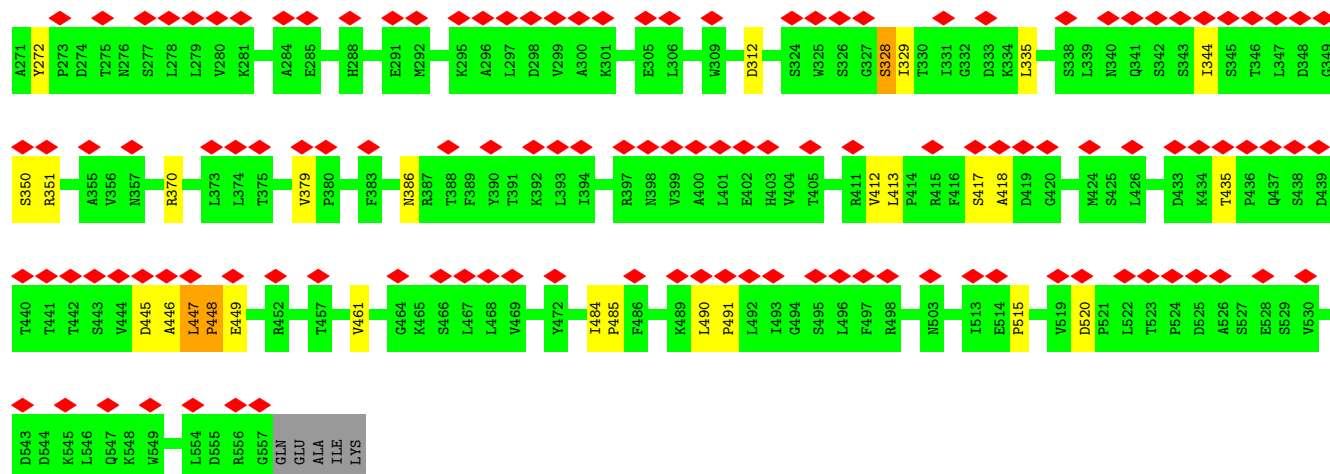


• Molecule 1: Protein InvG

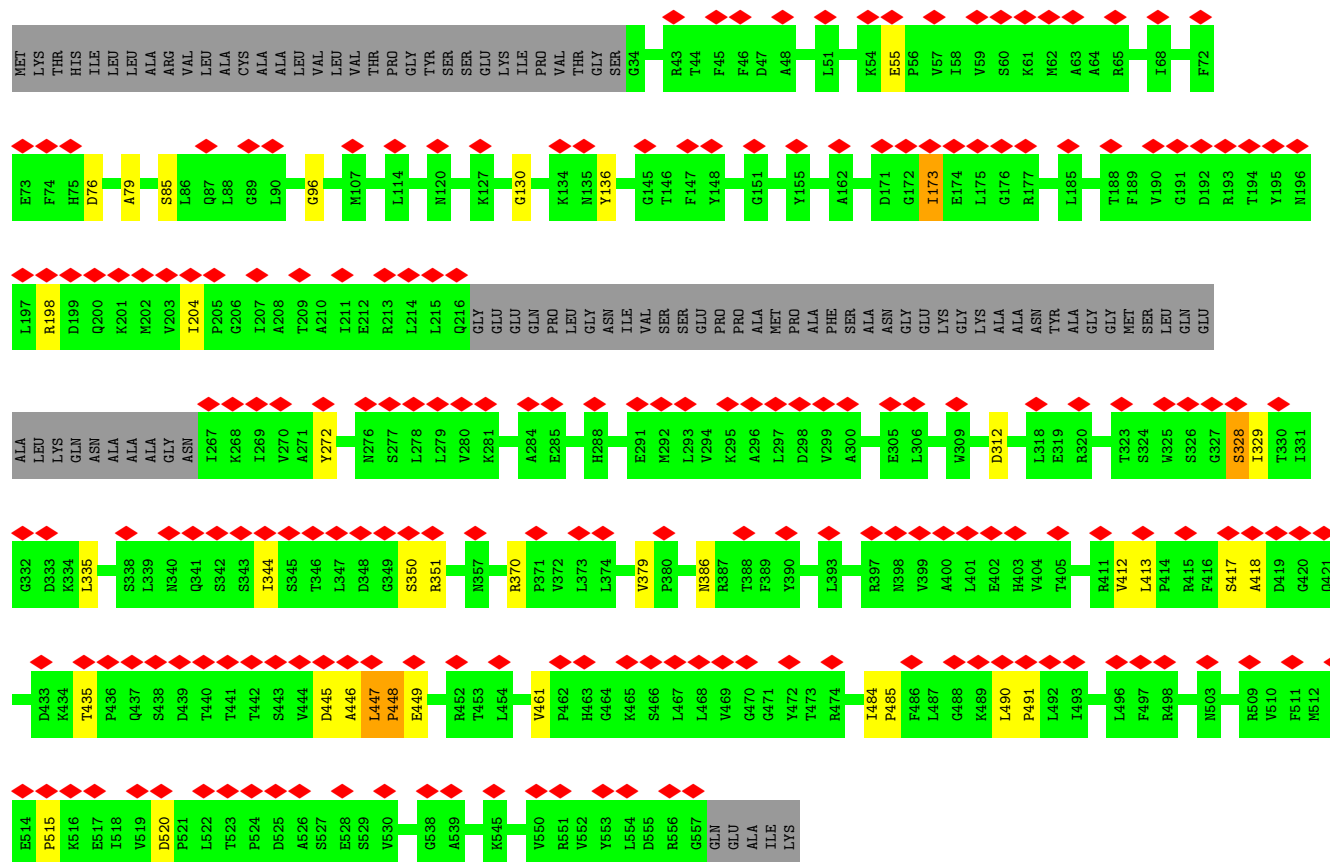
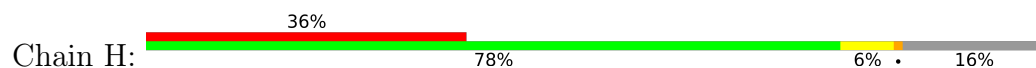


Chain D:

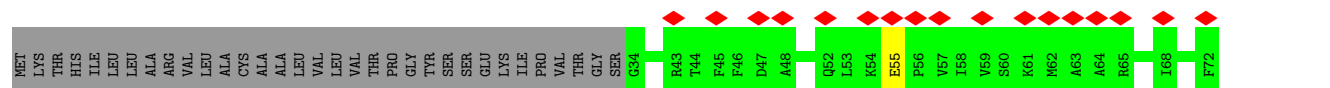
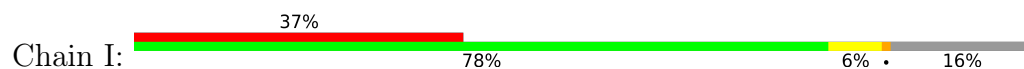


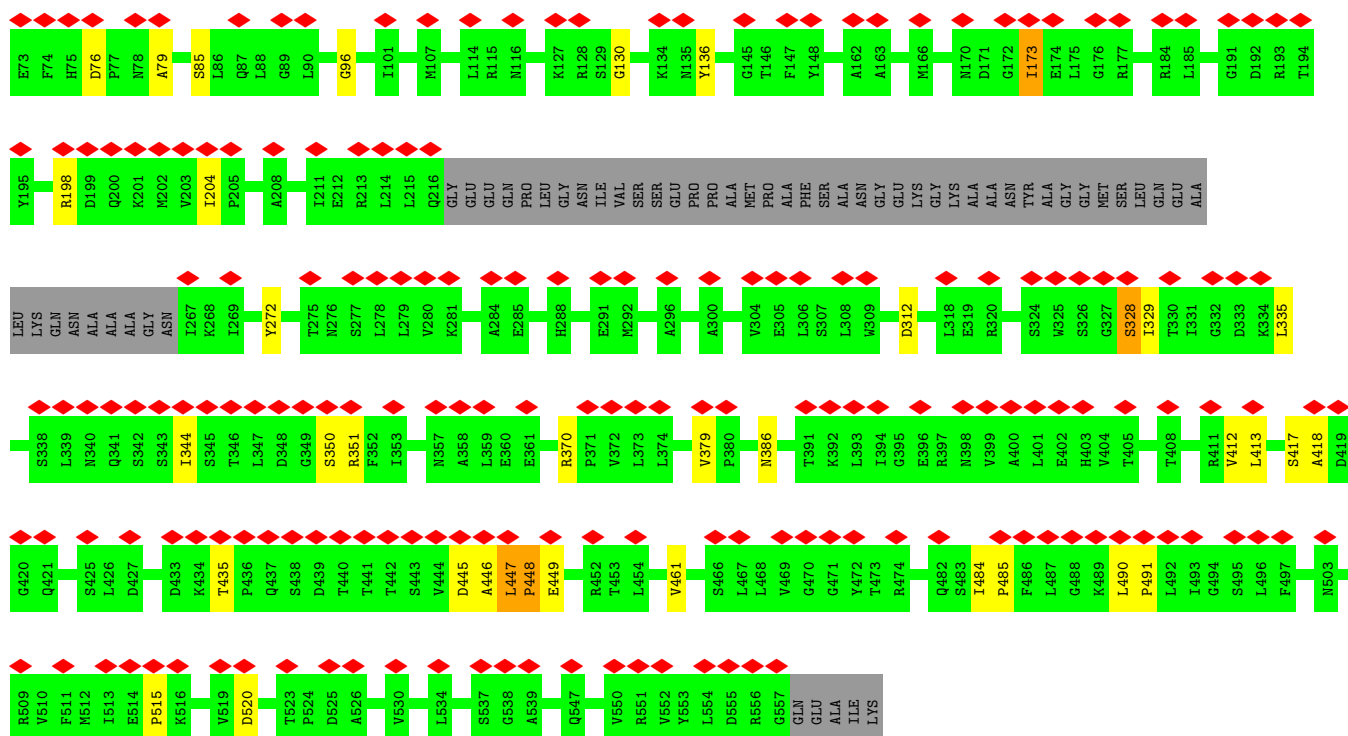


• Molecule 1: Protein InvG

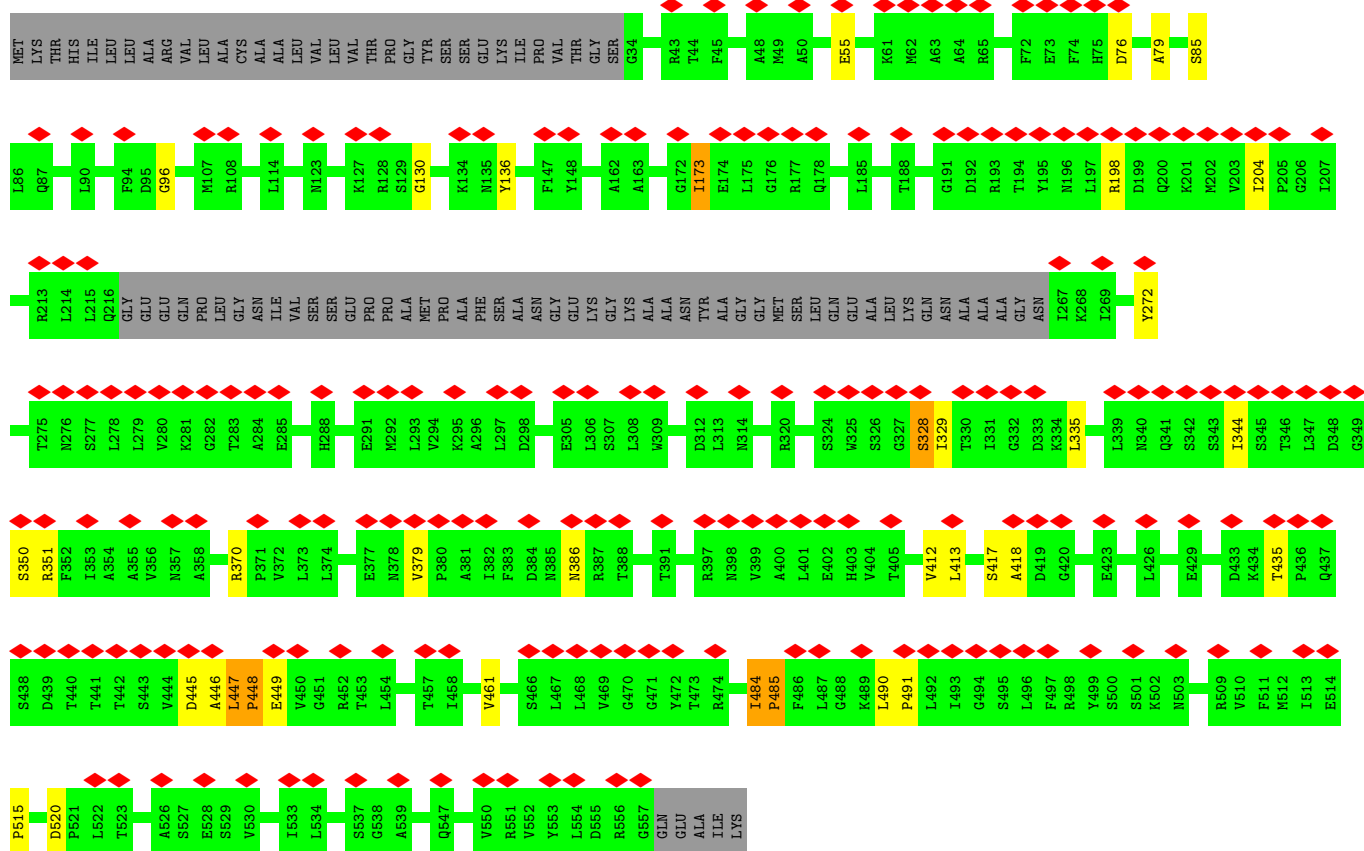
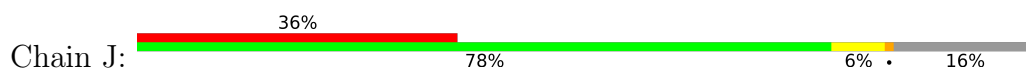


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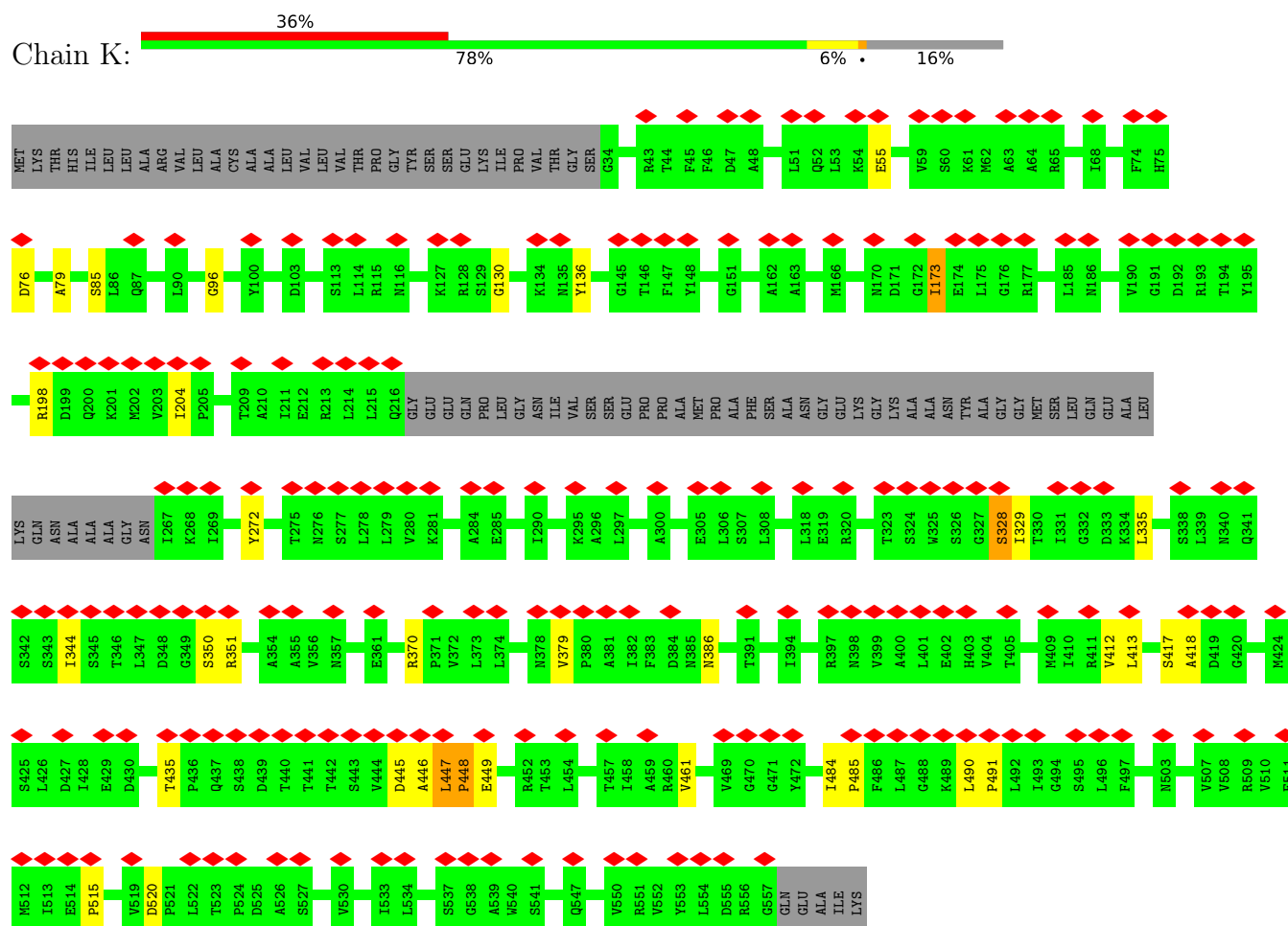




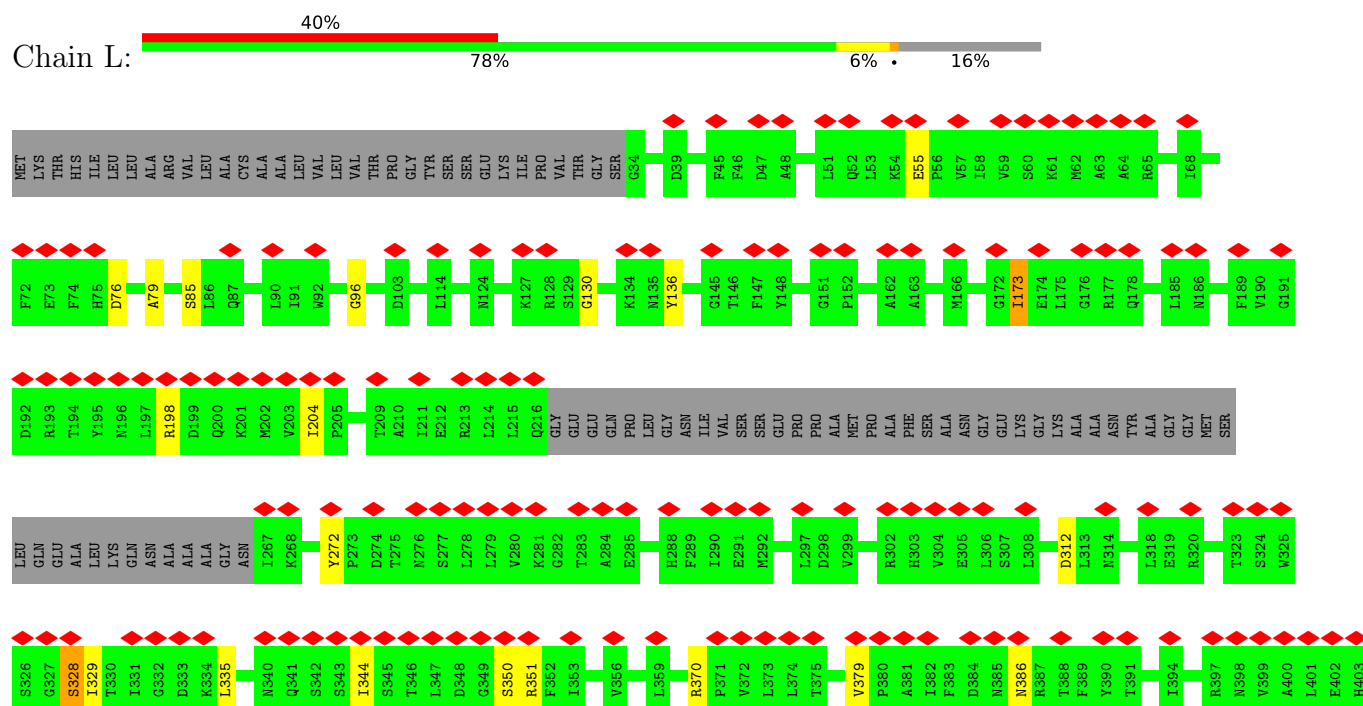
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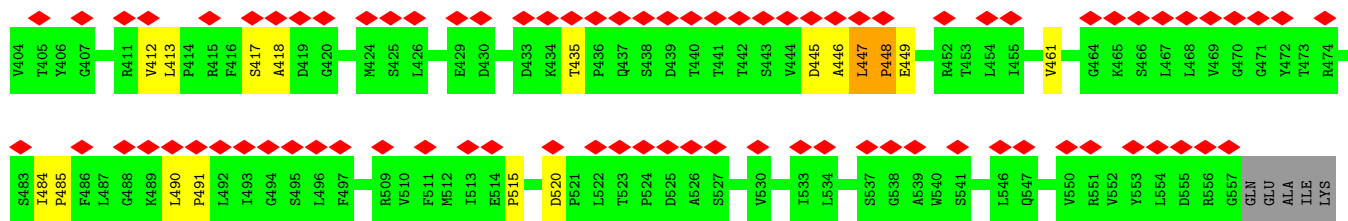


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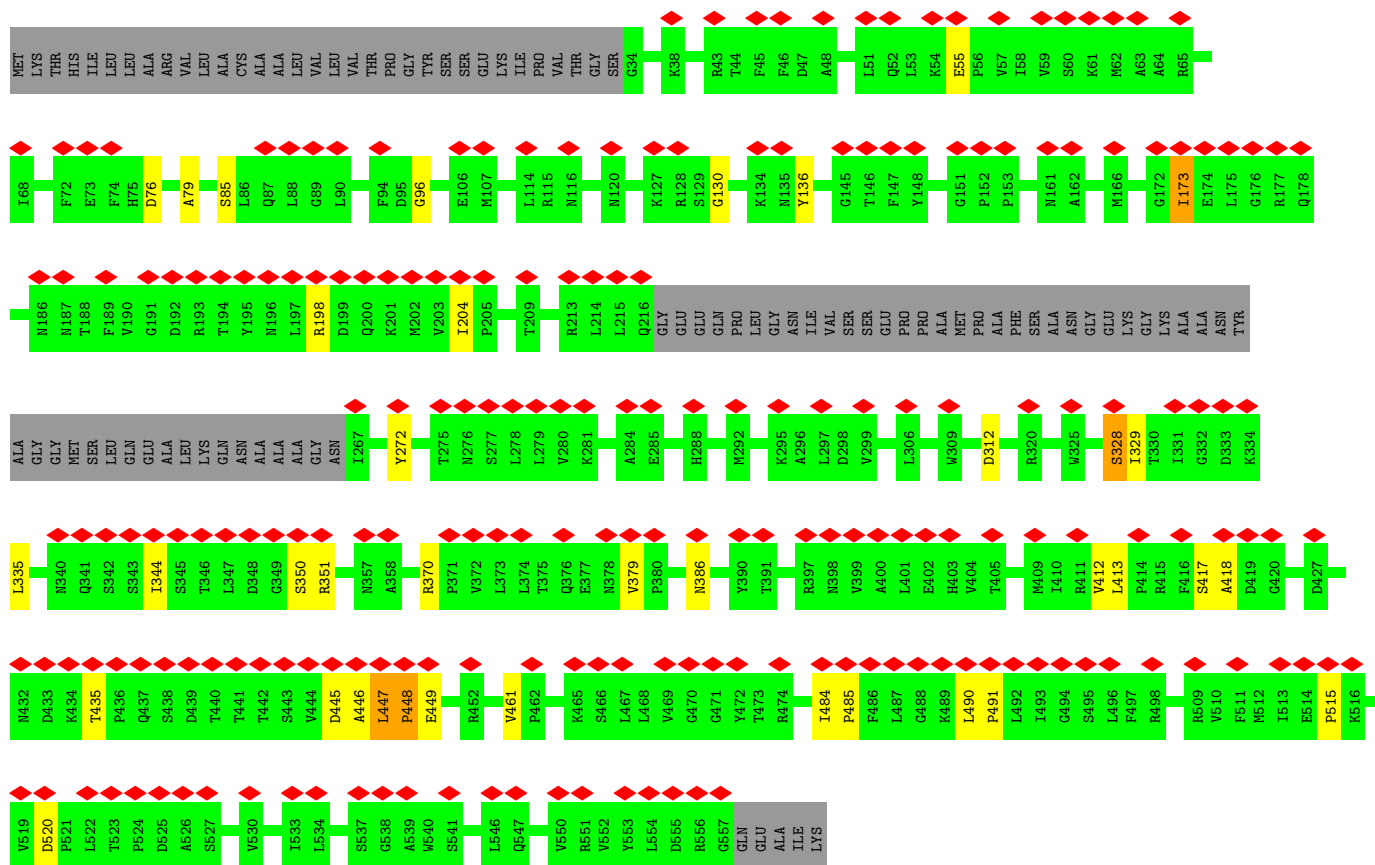
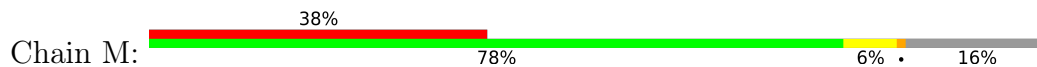


• Molecule 1: Protein InvG

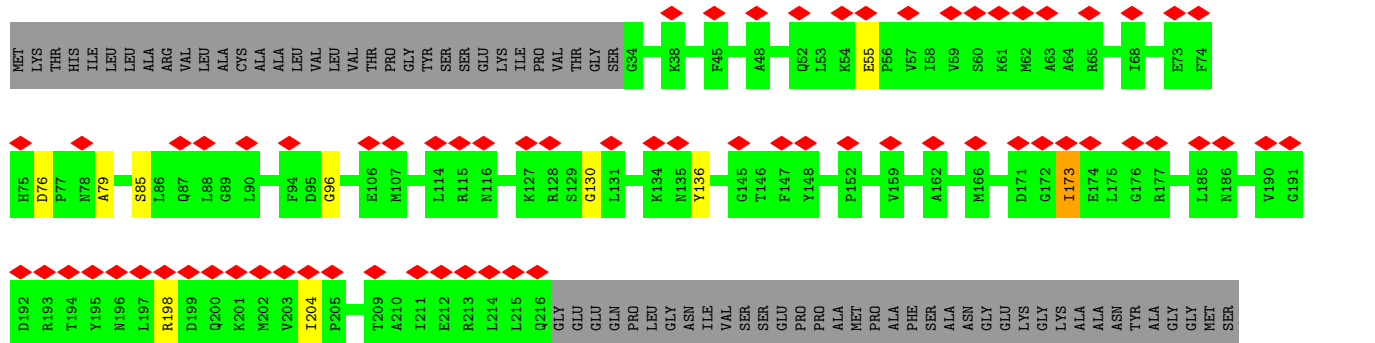
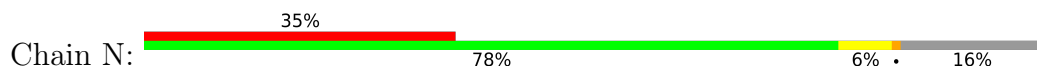


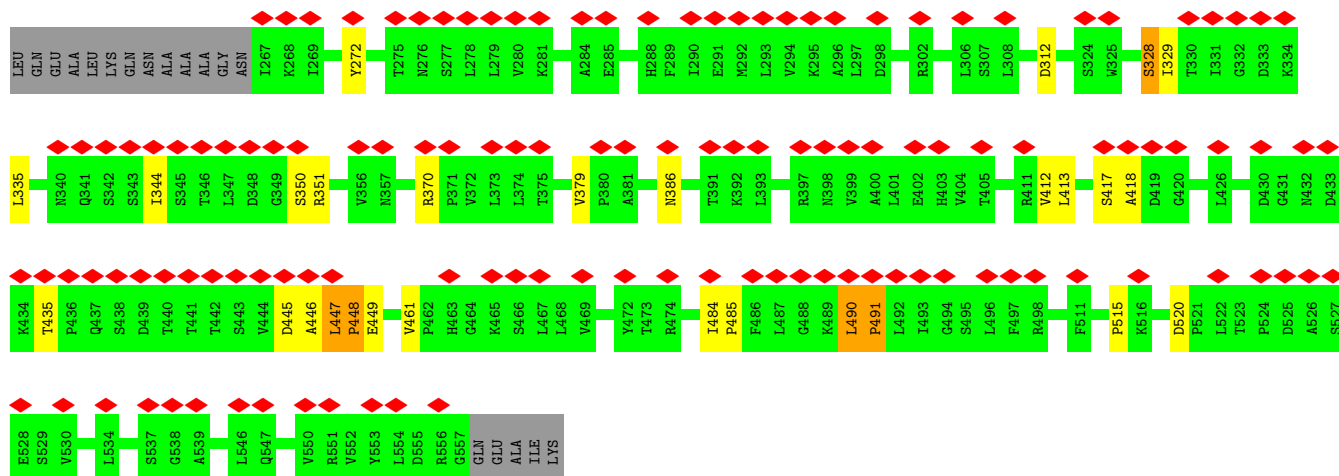


• Molecule 1: Protein InvG

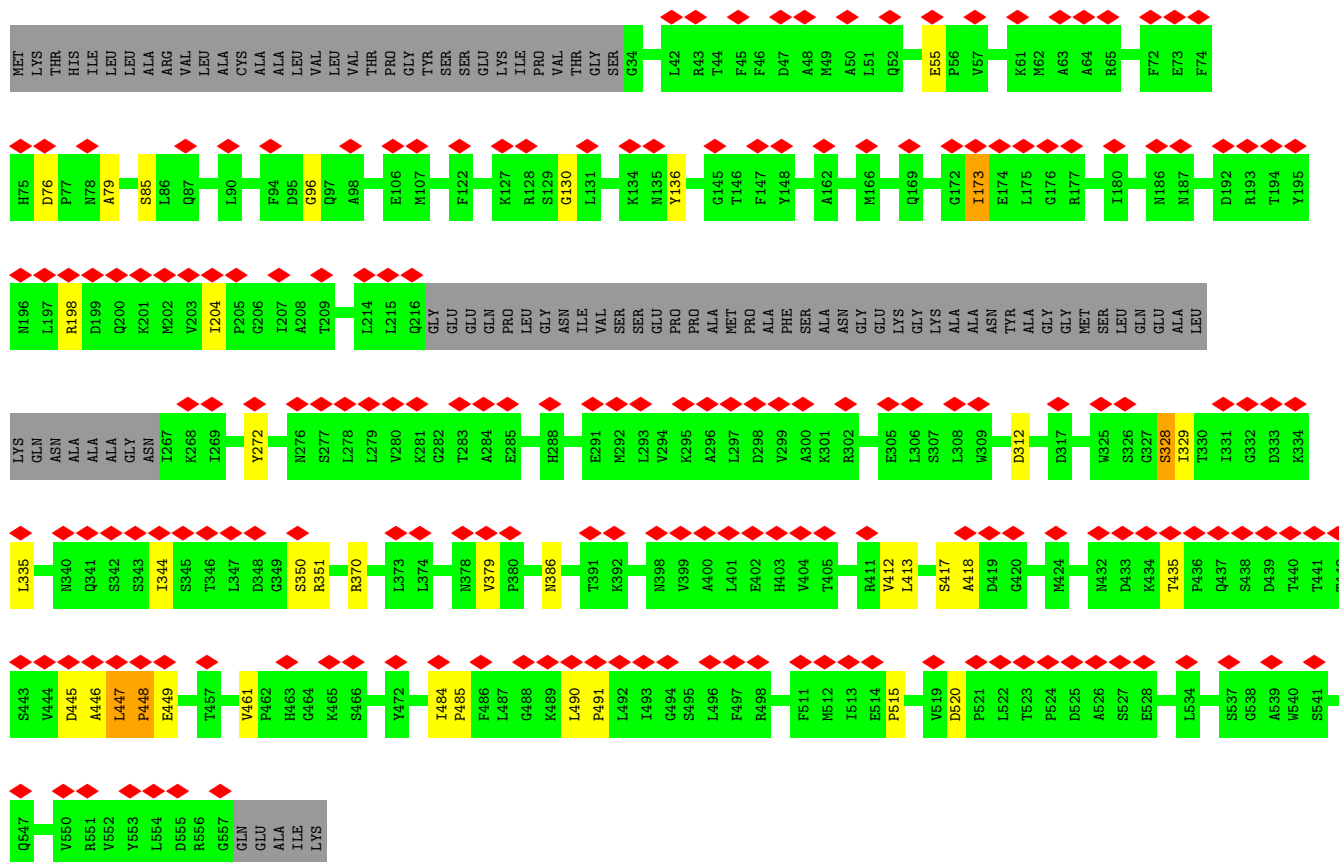
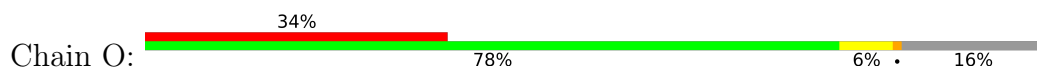


• Molecule 1: Protein InvG



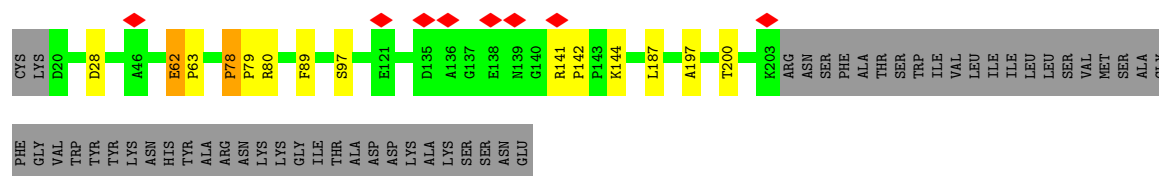


• Molecule 1: Protein InvG



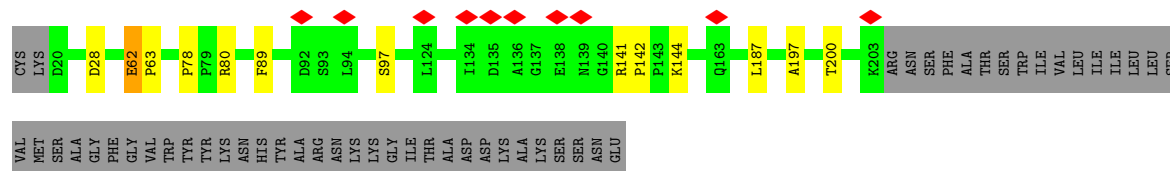
• Molecule 2: Lipoprotein PrgK





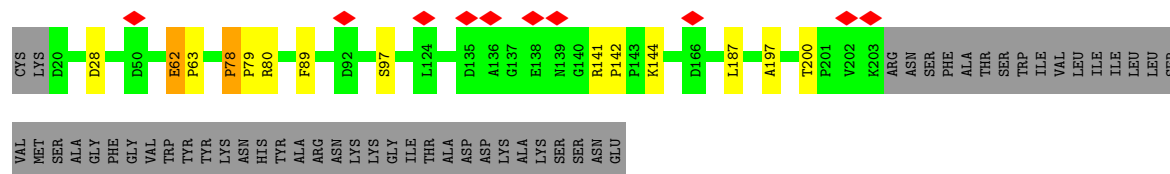
- Molecule 2: Lipoprotein PrgK

Chain R: 73% 5% 22%



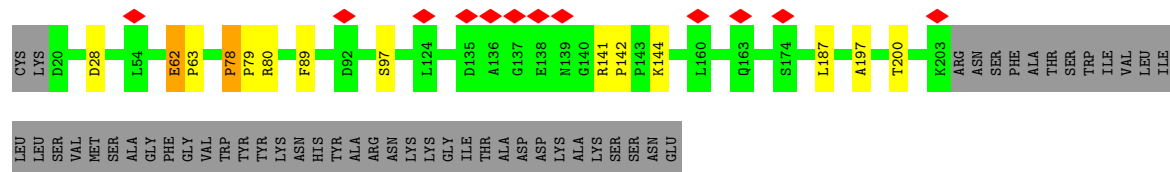
- Molecule 2: Lipoprotein PrgK

Chain T: 72% 5% 22%



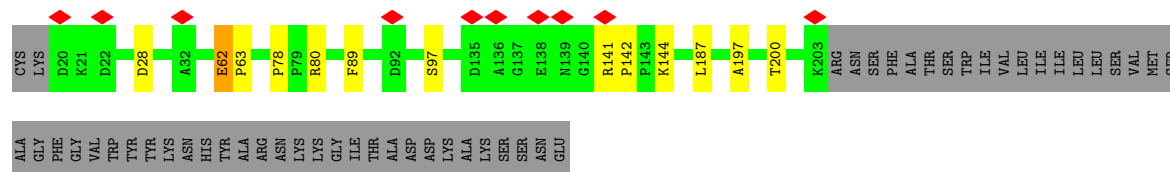
- Molecule 2: Lipoprotein PrgK

Chain V: 5% 72% 5% 22%



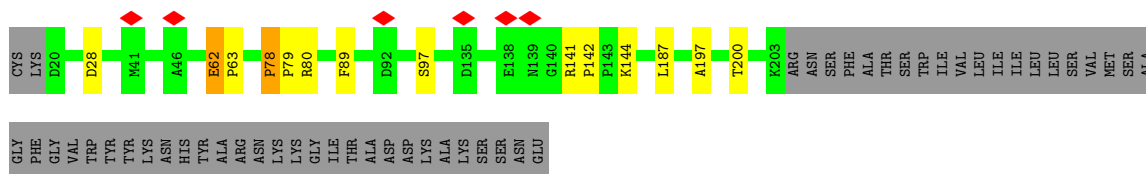
- Molecule 2: Lipoprotein PrgK

Chain X: 73% 5% 22%

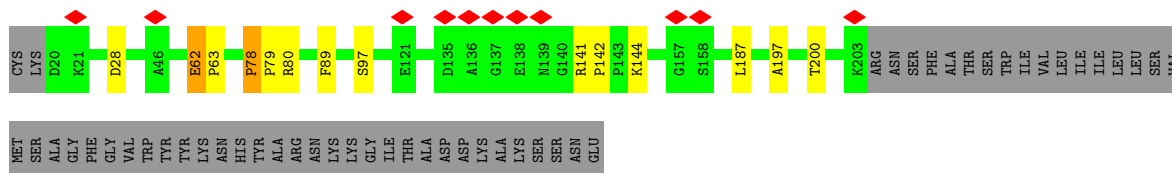


- Molecule 2: Lipoprotein PrgK

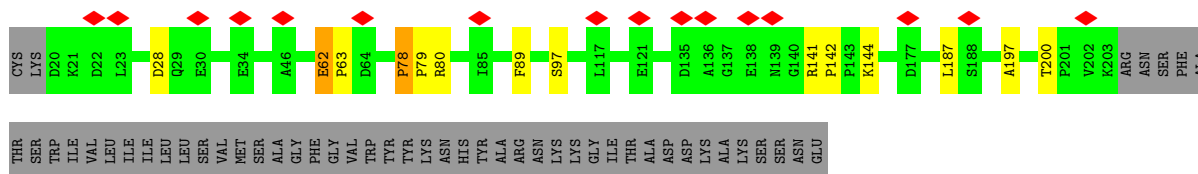
Chain a: 72% 5% 22%



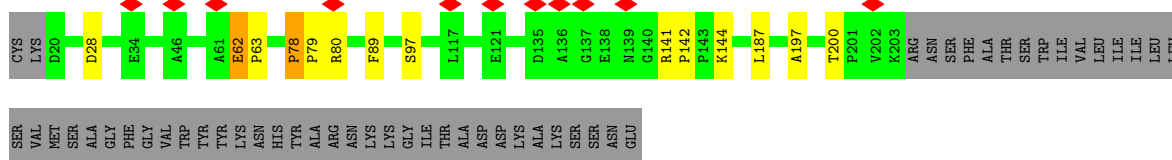
- Molecule 2: Lipoprotein PrgK



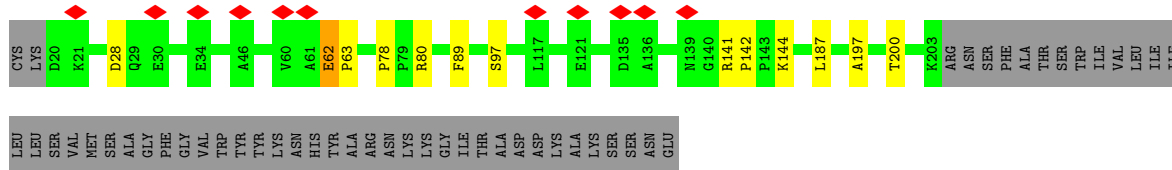
- Molecule 2: Lipoprotein PrgK



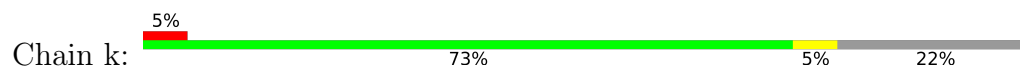
- Molecule 2: Lipoprotein PrgK

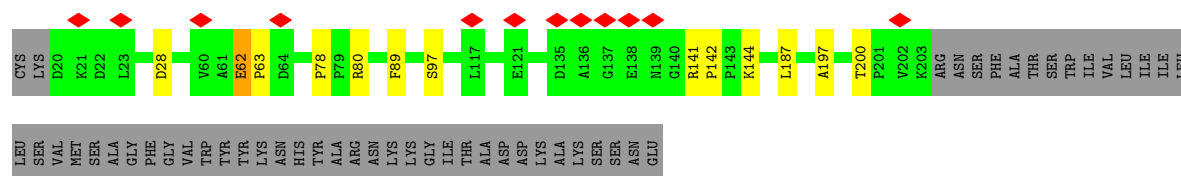


- Molecule 2: Lipoprotein PrgK

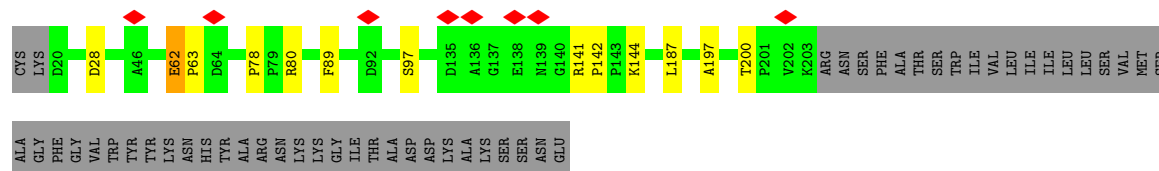


- Molecule 2: Lipoprotein PrgK

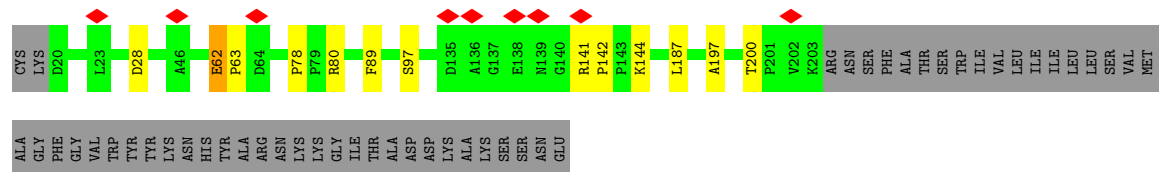




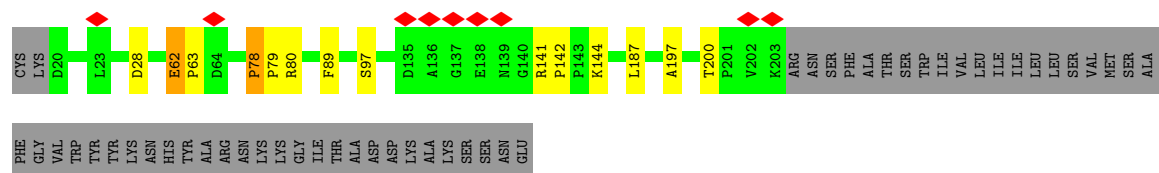
• Molecule 2: Lipoprotein PrgK



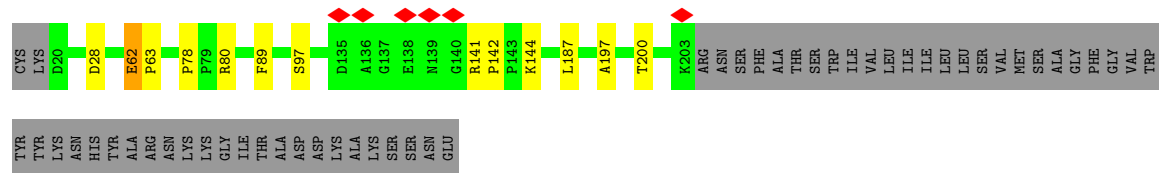
• Molecule 2: Lipoprotein PrgK



• Molecule 2: Lipoprotein PrgK

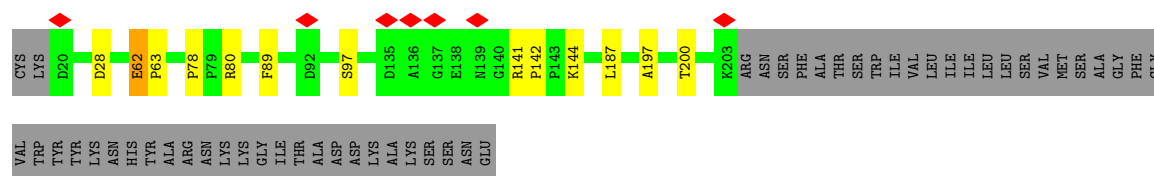


• Molecule 2: Lipoprotein PrgK

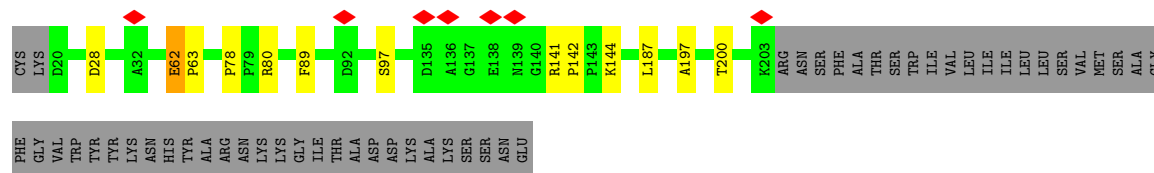


• Molecule 2: Lipoprotein PrgK

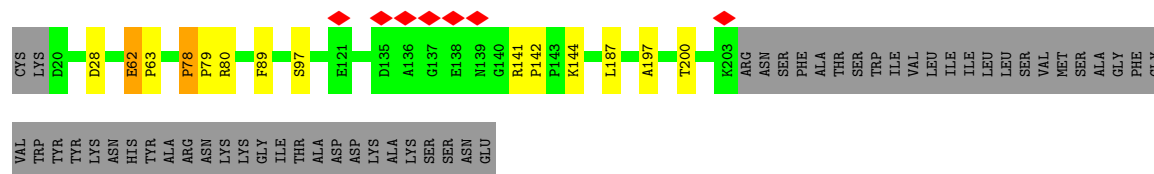




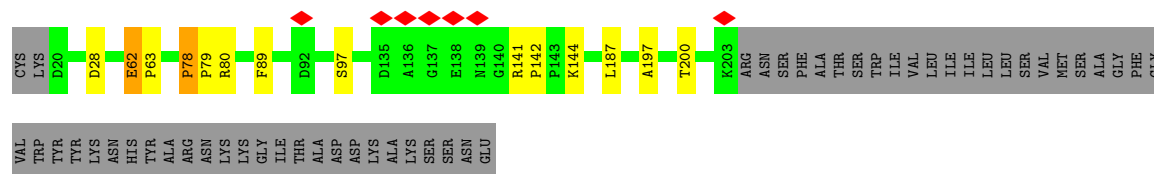
• Molecule 2: Lipoprotein PrgK



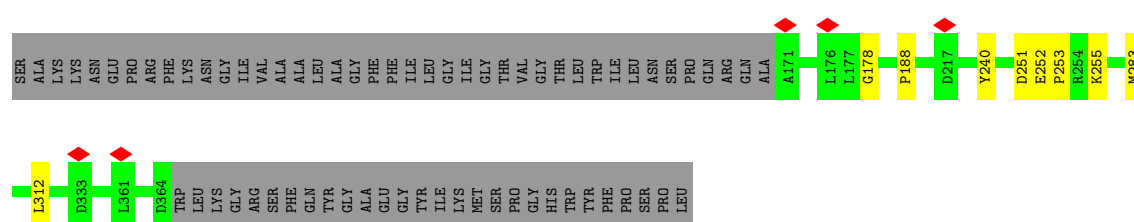
• Molecule 2: Lipoprotein PrgK



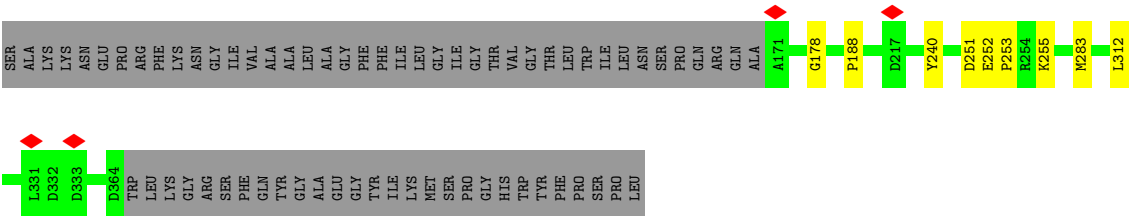
• Molecule 2: Lipoprotein PrgK



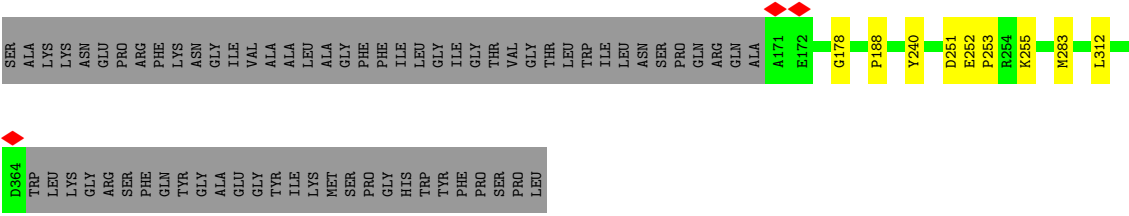
• Molecule 3: Protein PrgH



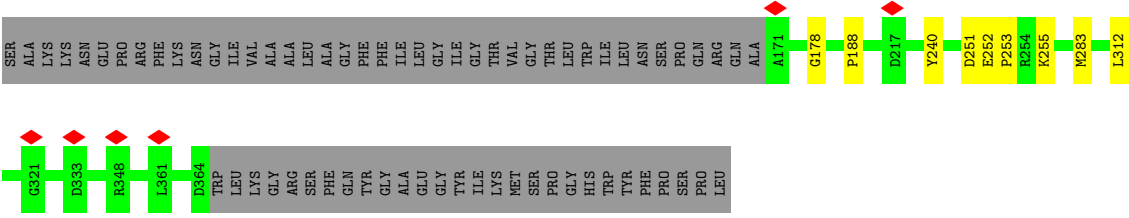
• Molecule 3: Protein PrgH



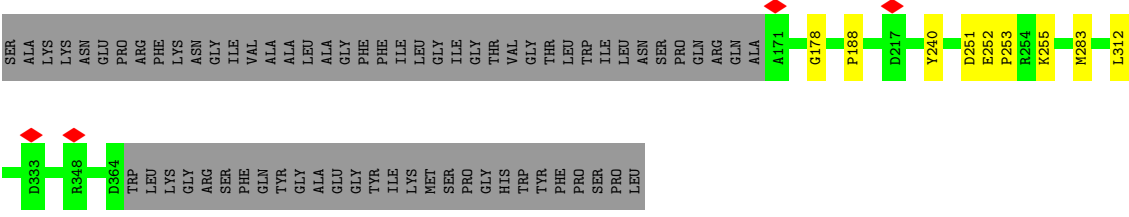
• Molecule 3: Protein PrgH



• Molecule 3: Protein PrgH

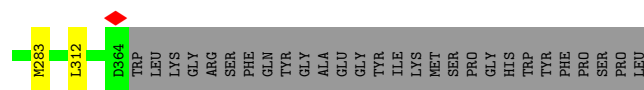


• Molecule 3: Protein PrgH

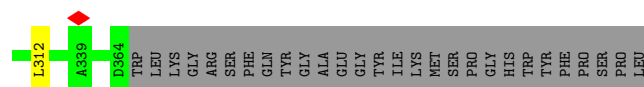
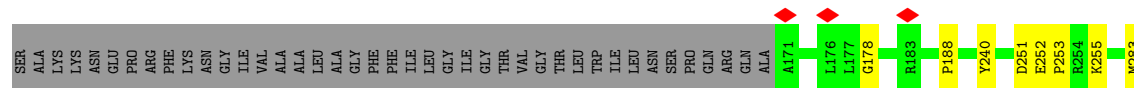


• Molecule 3: Protein PrgH

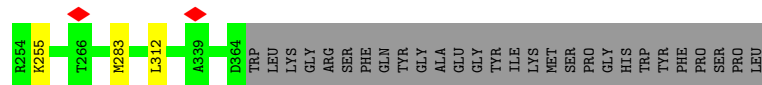
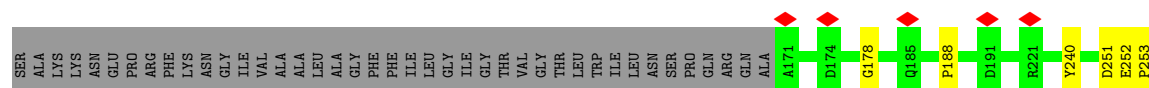




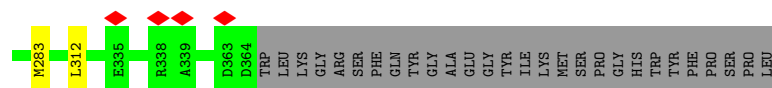
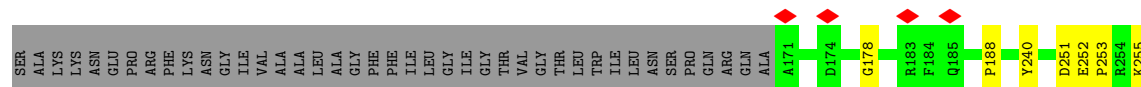
• Molecule 3: Protein PrgH



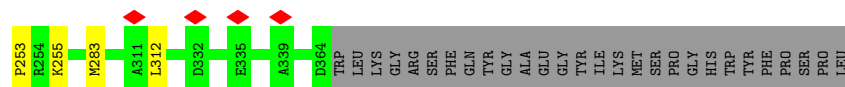
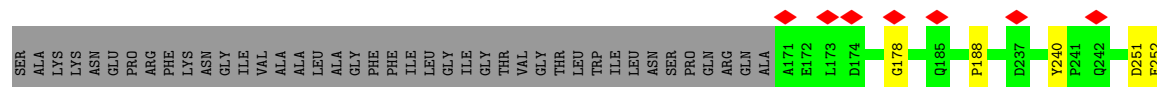
• Molecule 3: Protein PrgH



• Molecule 3: Protein PrgH

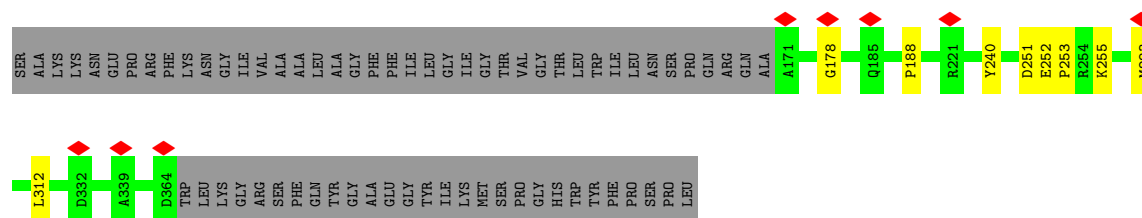


• Molecule 3: Protein PrgH



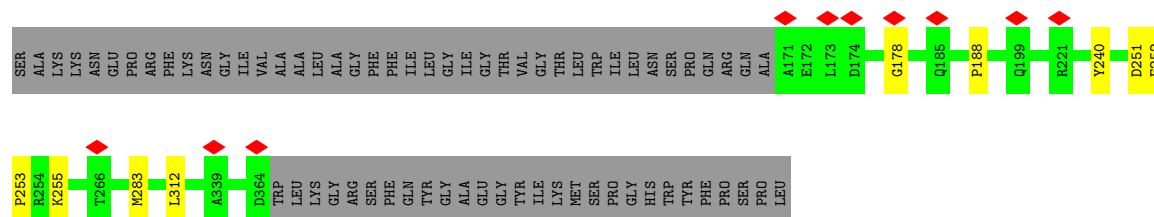
• Molecule 3: Protein PrgH

Chain j: 



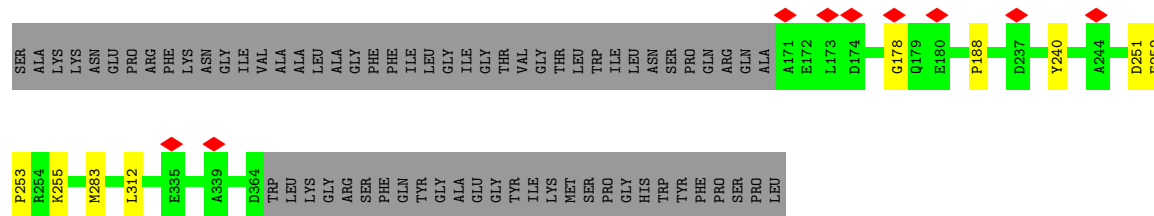
• Molecule 3: Protein PrgH

Chain l: 



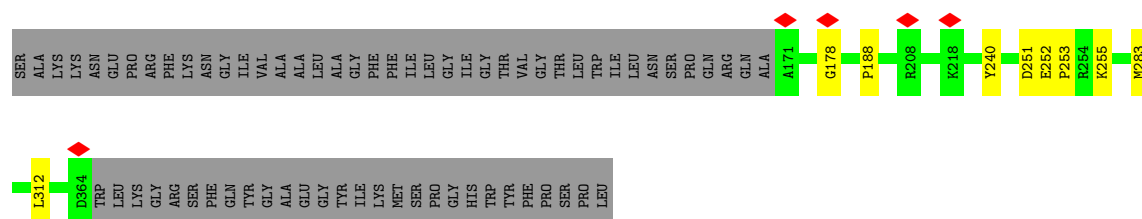
• Molecule 3: Protein PrgH

Chain n: 



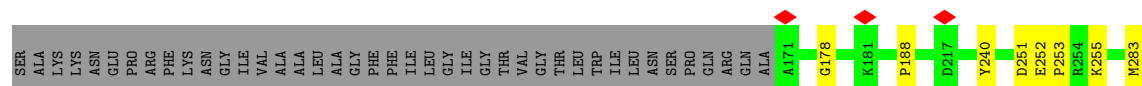
• Molecule 3: Protein PrgH

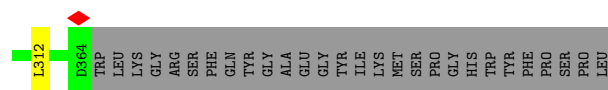
Chain p: 



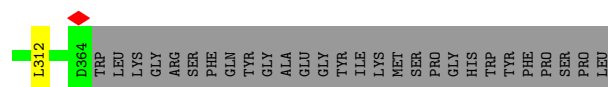
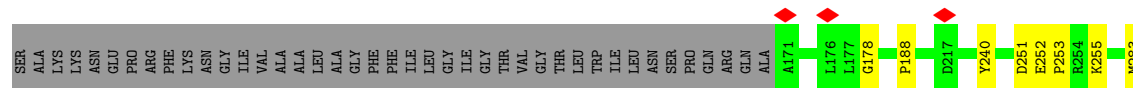
• Molecule 3: Protein PrgH

Chain r: 

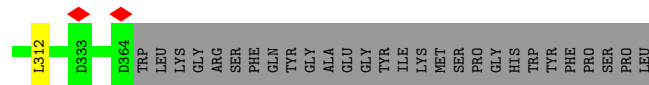
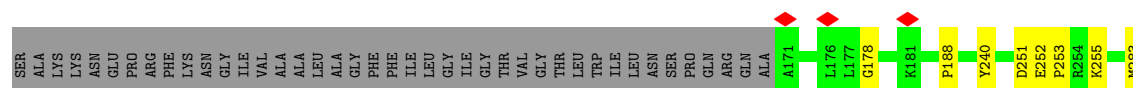




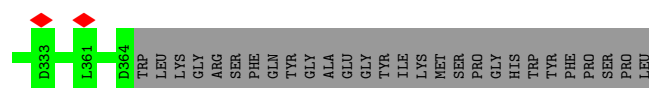
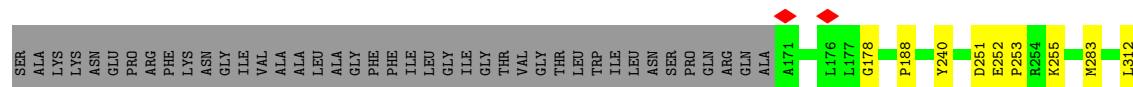
• Molecule 3: Protein PrgH



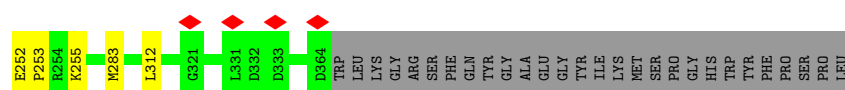
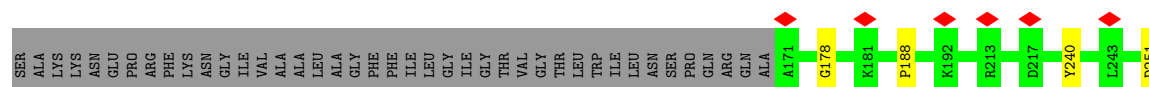
• Molecule 3: Protein PrgH



• Molecule 3: Protein PrgH

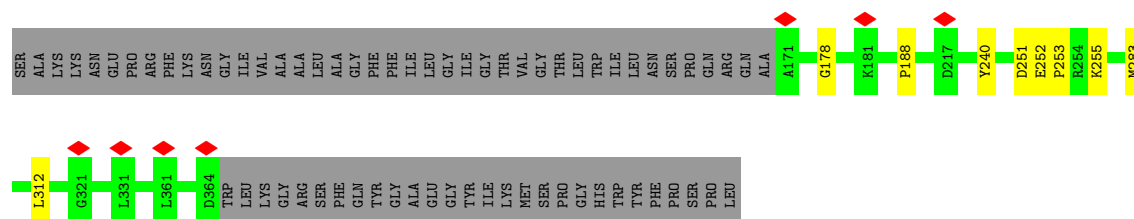


• Molecule 3: Protein PrgH



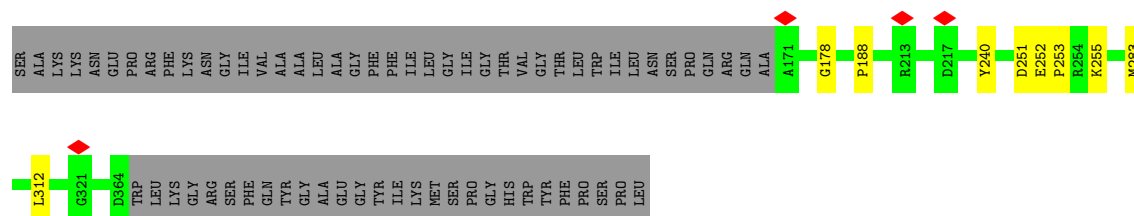
• Molecule 3: Protein PrgH

Chain 2:  70% 26%



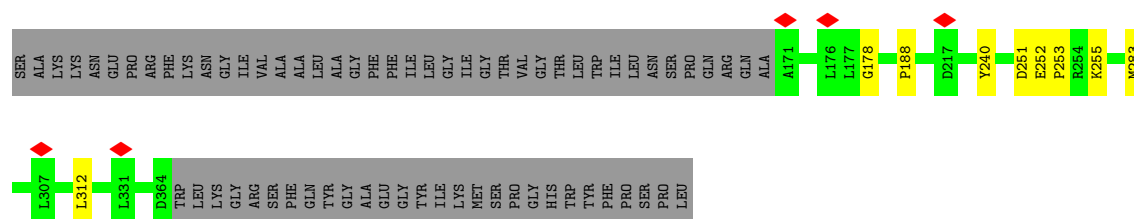
• Molecule 3: Protein PrgH

Chain 4:  70% 26%



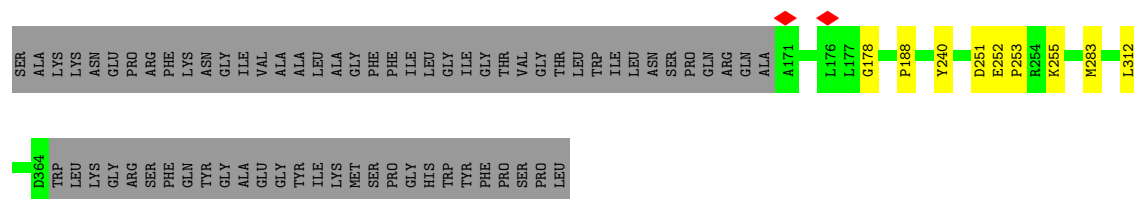
• Molecule 3: Protein PrgH

Chain 6:  70% 26%



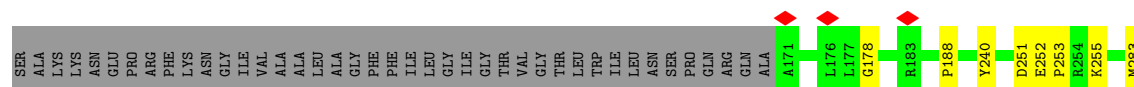
• Molecule 3: Protein PrgH

Chain 8:  70% 26%



• Molecule 3: Protein PrgH

Chain 10:  70% 26%



L312	L331	D364	TRP	LEU	LYS	GLY	ARG	SER	PHE	GLN	TYR	GLY	ALA	GLU	GLY	TYR	ILE	LYS	MET	SER	PRO	GLY	HIS	TRP	TYR	PHE	PRO	SER	PRO	LEU
------	------	------	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	67800	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 ($e^-/\text{\AA}^2$)	1.3	Depositor
Minimum defocus (nm)	1300	Depositor
Maximum defocus (nm)	3200	Depositor
Magnification	29240	Depositor
Image detector	GATAN K2 QUANTUM (4k x 4k)	Depositor
Maximum map value	0.187	Depositor
Minimum map value	-0.072	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.010	Depositor
Recommended contour level	0.05	Depositor
Map size (Å)	513.0, 513.0, 513.0	wwPDB
Map dimensions	300, 300, 300	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.71, 1.71, 1.71	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	0.86	1/3780 (0.0%)	1.11	30/5115 (0.6%)
1	B	0.86	1/3780 (0.0%)	1.11	30/5115 (0.6%)
1	C	0.86	1/3780 (0.0%)	1.11	30/5115 (0.6%)
1	D	0.86	1/3780 (0.0%)	1.11	30/5115 (0.6%)
1	E	0.86	1/3780 (0.0%)	1.11	30/5115 (0.6%)
1	F	0.86	1/3780 (0.0%)	1.11	30/5115 (0.6%)
1	G	0.86	1/3780 (0.0%)	1.11	30/5115 (0.6%)
1	H	0.86	1/3780 (0.0%)	1.11	30/5115 (0.6%)
1	I	0.86	1/3780 (0.0%)	1.11	30/5115 (0.6%)
1	J	0.85	1/3780 (0.0%)	1.11	30/5115 (0.6%)
1	K	0.86	1/3780 (0.0%)	1.11	30/5115 (0.6%)
1	L	0.86	1/3780 (0.0%)	1.11	30/5115 (0.6%)
1	M	0.85	1/3780 (0.0%)	1.11	30/5115 (0.6%)
1	N	0.86	1/3780 (0.0%)	1.11	30/5115 (0.6%)
1	O	0.85	1/3780 (0.0%)	1.11	30/5115 (0.6%)
2	1	0.77	0/1465	1.19	20/1989 (1.0%)
2	3	0.77	0/1465	1.19	20/1989 (1.0%)
2	5	0.77	0/1465	1.20	20/1989 (1.0%)
2	7	0.77	0/1465	1.19	20/1989 (1.0%)
2	9	0.77	0/1465	1.19	20/1989 (1.0%)
2	P	0.77	0/1465	1.19	20/1989 (1.0%)
2	R	0.77	0/1465	1.20	20/1989 (1.0%)
2	T	0.77	0/1465	1.19	20/1989 (1.0%)
2	V	0.77	0/1465	1.19	20/1989 (1.0%)
2	X	0.77	0/1465	1.19	20/1989 (1.0%)
2	a	0.77	0/1465	1.20	20/1989 (1.0%)
2	c	0.77	0/1465	1.19	20/1989 (1.0%)
2	e	0.77	0/1465	1.19	20/1989 (1.0%)
2	g	0.77	0/1465	1.19	20/1989 (1.0%)
2	i	0.77	0/1465	1.19	20/1989 (1.0%)
2	k	0.77	0/1465	1.19	20/1989 (1.0%)
2	m	0.77	0/1465	1.19	20/1989 (1.0%)
2	o	0.77	0/1465	1.19	20/1989 (1.0%)
2	q	0.77	0/1465	1.19	20/1989 (1.0%)

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
2	s	0.77	0/1465	1.19	20/1989 (1.0%)
2	u	0.77	0/1465	1.19	20/1989 (1.0%)
2	w	0.77	0/1465	1.19	20/1989 (1.0%)
2	y	0.77	0/1465	1.19	20/1989 (1.0%)
2	z	0.77	0/1465	1.20	20/1989 (1.0%)
3	0	0.76	0/1632	0.99	10/2204 (0.5%)
3	10	0.76	0/1632	0.99	10/2204 (0.5%)
3	2	0.76	0/1632	0.99	10/2204 (0.5%)
3	4	0.76	0/1632	0.99	10/2204 (0.5%)
3	6	0.76	0/1632	0.99	10/2204 (0.5%)
3	8	0.76	0/1632	0.99	10/2204 (0.5%)
3	Q	0.76	0/1632	0.99	10/2204 (0.5%)
3	S	0.76	0/1632	0.99	10/2204 (0.5%)
3	U	0.76	0/1632	0.99	10/2204 (0.5%)
3	W	0.76	0/1632	0.99	10/2204 (0.5%)
3	Y	0.76	0/1632	0.99	10/2204 (0.5%)
3	Z	0.76	0/1632	0.99	10/2204 (0.5%)
3	b	0.76	0/1632	0.99	10/2204 (0.5%)
3	d	0.76	0/1632	0.99	10/2204 (0.5%)
3	f	0.76	0/1632	0.99	10/2204 (0.5%)
3	h	0.76	0/1632	0.99	10/2204 (0.5%)
3	j	0.76	0/1632	0.99	10/2204 (0.5%)
3	l	0.76	0/1632	0.99	10/2204 (0.5%)
3	n	0.76	0/1632	0.99	10/2204 (0.5%)
3	p	0.76	0/1632	0.99	10/2204 (0.5%)
3	r	0.76	0/1632	0.99	10/2204 (0.5%)
3	t	0.76	0/1632	0.99	10/2204 (0.5%)
3	v	0.76	0/1632	0.99	10/2204 (0.5%)
3	x	0.76	0/1632	0.99	10/2204 (0.5%)
All	All	0.80	15/131028 (0.0%)	1.10	1170/177357 (0.7%)

The worst 5 of 15 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	79	ALA	CA-CB	-5.11	1.44	1.53
1	K	79	ALA	CA-CB	-5.11	1.44	1.53
1	M	79	ALA	CA-CB	-5.10	1.44	1.53
1	G	79	ALA	CA-CB	-5.09	1.44	1.53
1	B	79	ALA	CA-CB	-5.09	1.44	1.53

The worst 5 of 1170 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	P	141	ARG	CA-C-N	10.30	127.08	119.66
2	P	141	ARG	C-N-CA	10.30	127.08	119.66
2	e	141	ARG	CA-C-N	10.30	127.08	119.66
2	e	141	ARG	C-N-CA	10.30	127.08	119.66
2	R	141	ARG	CA-C-N	10.29	127.07	119.66

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	3718	0	3777	10	0
1	B	3718	0	3777	10	0
1	C	3718	0	3777	10	0
1	D	3718	0	3777	9	0
1	E	3718	0	3777	9	0
1	F	3718	0	3777	10	0
1	G	3718	0	3777	10	0
1	H	3718	0	3777	10	0
1	I	3718	0	3777	10	0
1	J	3718	0	3777	10	0
1	K	3718	0	3777	9	0
1	L	3718	0	3777	10	0
1	M	3718	0	3777	10	0
1	N	3718	0	3777	11	0
1	O	3718	0	3777	10	0
2	1	1437	0	1434	4	0
2	3	1437	0	1434	4	0
2	5	1437	0	1434	4	0
2	7	1437	0	1434	5	0
2	9	1437	0	1434	5	0
2	P	1437	0	1434	5	0
2	R	1437	0	1434	4	0
2	T	1437	0	1434	5	0
2	V	1437	0	1434	5	0
2	X	1437	0	1434	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	a	1437	0	1434	5	0
2	c	1437	0	1434	5	0
2	e	1437	0	1434	5	0
2	g	1437	0	1434	5	0
2	i	1437	0	1434	4	0
2	k	1437	0	1434	4	0
2	m	1437	0	1434	4	0
2	o	1437	0	1434	4	0
2	q	1437	0	1434	5	0
2	s	1437	0	1434	4	0
2	u	1437	0	1434	5	0
2	w	1437	0	1434	4	0
2	y	1437	0	1434	4	0
2	z	1437	0	1434	5	0
3	0	1600	0	1580	0	0
3	10	1600	0	1580	0	0
3	2	1600	0	1580	0	0
3	4	1600	0	1580	0	0
3	6	1600	0	1580	0	0
3	8	1600	0	1580	0	0
3	Q	1600	0	1580	0	0
3	S	1600	0	1580	0	0
3	U	1600	0	1580	0	0
3	W	1600	0	1580	0	0
3	Y	1600	0	1580	0	0
3	Z	1600	0	1580	0	0
3	b	1600	0	1580	0	0
3	d	1600	0	1580	0	0
3	f	1600	0	1580	0	0
3	h	1600	0	1580	0	0
3	j	1600	0	1580	0	0
3	l	1600	0	1580	0	0
3	n	1600	0	1580	0	0
3	p	1600	0	1580	0	0
3	r	1600	0	1580	0	0
3	t	1600	0	1580	0	0
3	v	1600	0	1580	0	0
3	x	1600	0	1580	0	0
All	All	128658	0	128991	256	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 1.

The worst 5 of 256 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:E:173:ILE:HG22	1:E:173:ILE:O	1.92	0.70
1:H:173:ILE:O	1:H:173:ILE:HG22	1.92	0.70
1:I:173:ILE:HG22	1:I:173:ILE:O	1.92	0.70
1:D:173:ILE:O	1:D:173:ILE:HG22	1.92	0.70
1:G:173:ILE:HG22	1:G:173:ILE:O	1.91	0.69

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	468/562 (83%)	444 (95%)	15 (3%)	9 (2%)	6	32
1	B	468/562 (83%)	444 (95%)	15 (3%)	9 (2%)	6	32
1	C	468/562 (83%)	444 (95%)	15 (3%)	9 (2%)	6	32
1	D	468/562 (83%)	444 (95%)	15 (3%)	9 (2%)	6	32
1	E	468/562 (83%)	444 (95%)	15 (3%)	9 (2%)	6	32
1	F	468/562 (83%)	444 (95%)	15 (3%)	9 (2%)	6	32
1	G	468/562 (83%)	444 (95%)	15 (3%)	9 (2%)	6	32
1	H	468/562 (83%)	444 (95%)	15 (3%)	9 (2%)	6	32
1	I	468/562 (83%)	444 (95%)	15 (3%)	9 (2%)	6	32
1	J	468/562 (83%)	444 (95%)	15 (3%)	9 (2%)	6	32
1	K	468/562 (83%)	444 (95%)	15 (3%)	9 (2%)	6	32
1	L	468/562 (83%)	444 (95%)	15 (3%)	9 (2%)	6	32
1	M	468/562 (83%)	444 (95%)	15 (3%)	9 (2%)	6	32
1	N	468/562 (83%)	444 (95%)	15 (3%)	9 (2%)	6	32

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	O	468/562 (83%)	444 (95%)	15 (3%)	9 (2%)	6	32
2	1	182/235 (77%)	175 (96%)	7 (4%)	0	100	100
2	3	182/235 (77%)	175 (96%)	7 (4%)	0	100	100
2	5	182/235 (77%)	175 (96%)	7 (4%)	0	100	100
2	7	182/235 (77%)	175 (96%)	7 (4%)	0	100	100
2	9	182/235 (77%)	175 (96%)	7 (4%)	0	100	100
2	P	182/235 (77%)	175 (96%)	7 (4%)	0	100	100
2	R	182/235 (77%)	175 (96%)	7 (4%)	0	100	100
2	T	182/235 (77%)	175 (96%)	7 (4%)	0	100	100
2	V	182/235 (77%)	175 (96%)	7 (4%)	0	100	100
2	X	182/235 (77%)	175 (96%)	7 (4%)	0	100	100
2	a	182/235 (77%)	175 (96%)	7 (4%)	0	100	100
2	c	182/235 (77%)	175 (96%)	7 (4%)	0	100	100
2	e	182/235 (77%)	175 (96%)	7 (4%)	0	100	100
2	g	182/235 (77%)	175 (96%)	7 (4%)	0	100	100
2	i	182/235 (77%)	175 (96%)	7 (4%)	0	100	100
2	k	182/235 (77%)	175 (96%)	7 (4%)	0	100	100
2	m	182/235 (77%)	175 (96%)	7 (4%)	0	100	100
2	o	182/235 (77%)	175 (96%)	7 (4%)	0	100	100
2	q	182/235 (77%)	175 (96%)	7 (4%)	0	100	100
2	s	182/235 (77%)	175 (96%)	7 (4%)	0	100	100
2	u	182/235 (77%)	175 (96%)	7 (4%)	0	100	100
2	w	182/235 (77%)	175 (96%)	7 (4%)	0	100	100
2	y	182/235 (77%)	175 (96%)	7 (4%)	0	100	100
2	z	182/235 (77%)	175 (96%)	7 (4%)	0	100	100
3	0	192/263 (73%)	182 (95%)	6 (3%)	4 (2%)	5	29
3	10	192/263 (73%)	182 (95%)	6 (3%)	4 (2%)	5	29
3	2	192/263 (73%)	182 (95%)	6 (3%)	4 (2%)	5	29
3	4	192/263 (73%)	182 (95%)	6 (3%)	4 (2%)	5	29
3	6	192/263 (73%)	182 (95%)	6 (3%)	4 (2%)	5	29
3	8	192/263 (73%)	182 (95%)	6 (3%)	4 (2%)	5	29

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	Q	192/263 (73%)	182 (95%)	6 (3%)	4 (2%)	5	29
3	S	192/263 (73%)	182 (95%)	6 (3%)	4 (2%)	5	29
3	U	192/263 (73%)	182 (95%)	6 (3%)	4 (2%)	5	29
3	W	192/263 (73%)	182 (95%)	6 (3%)	4 (2%)	5	29
3	Y	192/263 (73%)	182 (95%)	6 (3%)	4 (2%)	5	29
3	Z	192/263 (73%)	182 (95%)	6 (3%)	4 (2%)	5	29
3	b	192/263 (73%)	182 (95%)	6 (3%)	4 (2%)	5	29
3	d	192/263 (73%)	182 (95%)	6 (3%)	4 (2%)	5	29
3	f	192/263 (73%)	182 (95%)	6 (3%)	4 (2%)	5	29
3	h	192/263 (73%)	182 (95%)	6 (3%)	4 (2%)	5	29
3	j	192/263 (73%)	182 (95%)	6 (3%)	4 (2%)	5	29
3	l	192/263 (73%)	182 (95%)	6 (3%)	4 (2%)	5	29
3	n	192/263 (73%)	182 (95%)	6 (3%)	4 (2%)	5	29
3	p	192/263 (73%)	182 (95%)	6 (3%)	4 (2%)	5	29
3	r	192/263 (73%)	182 (95%)	6 (3%)	4 (2%)	5	29
3	t	192/263 (73%)	182 (95%)	6 (3%)	4 (2%)	5	29
3	v	192/263 (73%)	182 (95%)	6 (3%)	4 (2%)	5	29
3	x	192/263 (73%)	182 (95%)	6 (3%)	4 (2%)	5	29
All	All	15996/20382 (78%)	15228 (95%)	537 (3%)	231 (1%)	11	40

5 of 231 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	446	ALA
1	B	446	ALA
1	B	515	PRO
1	C	446	ALA
1	C	515	PRO

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	413/477 (87%)	413 (100%)	0	100	100
1	B	413/477 (87%)	413 (100%)	0	100	100
1	C	413/477 (87%)	413 (100%)	0	100	100
1	D	413/477 (87%)	413 (100%)	0	100	100
1	E	413/477 (87%)	413 (100%)	0	100	100
1	F	413/477 (87%)	413 (100%)	0	100	100
1	G	413/477 (87%)	413 (100%)	0	100	100
1	H	413/477 (87%)	413 (100%)	0	100	100
1	I	413/477 (87%)	413 (100%)	0	100	100
1	J	413/477 (87%)	413 (100%)	0	100	100
1	K	413/477 (87%)	413 (100%)	0	100	100
1	L	413/477 (87%)	413 (100%)	0	100	100
1	M	413/477 (87%)	413 (100%)	0	100	100
1	N	413/477 (87%)	413 (100%)	0	100	100
1	O	413/477 (87%)	413 (100%)	0	100	100
2	1	157/200 (78%)	157 (100%)	0	100	100
2	3	157/200 (78%)	157 (100%)	0	100	100
2	5	157/200 (78%)	157 (100%)	0	100	100
2	7	157/200 (78%)	157 (100%)	0	100	100
2	9	157/200 (78%)	157 (100%)	0	100	100
2	P	157/200 (78%)	157 (100%)	0	100	100
2	R	157/200 (78%)	157 (100%)	0	100	100
2	T	157/200 (78%)	157 (100%)	0	100	100
2	V	157/200 (78%)	157 (100%)	0	100	100
2	X	157/200 (78%)	157 (100%)	0	100	100
2	a	157/200 (78%)	157 (100%)	0	100	100
2	c	157/200 (78%)	157 (100%)	0	100	100
2	e	157/200 (78%)	157 (100%)	0	100	100
2	g	157/200 (78%)	157 (100%)	0	100	100
2	i	157/200 (78%)	157 (100%)	0	100	100
2	k	157/200 (78%)	157 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	m	157/200 (78%)	157 (100%)	0	100	100
2	o	157/200 (78%)	157 (100%)	0	100	100
2	q	157/200 (78%)	157 (100%)	0	100	100
2	s	157/200 (78%)	157 (100%)	0	100	100
2	u	157/200 (78%)	157 (100%)	0	100	100
2	w	157/200 (78%)	157 (100%)	0	100	100
2	y	157/200 (78%)	157 (100%)	0	100	100
2	z	157/200 (78%)	157 (100%)	0	100	100
3	0	167/221 (76%)	167 (100%)	0	100	100
3	10	167/221 (76%)	167 (100%)	0	100	100
3	2	167/221 (76%)	167 (100%)	0	100	100
3	4	167/221 (76%)	167 (100%)	0	100	100
3	6	167/221 (76%)	167 (100%)	0	100	100
3	8	167/221 (76%)	167 (100%)	0	100	100
3	Q	167/221 (76%)	167 (100%)	0	100	100
3	S	167/221 (76%)	167 (100%)	0	100	100
3	U	167/221 (76%)	167 (100%)	0	100	100
3	W	167/221 (76%)	167 (100%)	0	100	100
3	Y	167/221 (76%)	167 (100%)	0	100	100
3	Z	167/221 (76%)	167 (100%)	0	100	100
3	b	167/221 (76%)	167 (100%)	0	100	100
3	d	167/221 (76%)	167 (100%)	0	100	100
3	f	167/221 (76%)	167 (100%)	0	100	100
3	h	167/221 (76%)	167 (100%)	0	100	100
3	j	167/221 (76%)	167 (100%)	0	100	100
3	l	167/221 (76%)	167 (100%)	0	100	100
3	n	167/221 (76%)	167 (100%)	0	100	100
3	p	167/221 (76%)	167 (100%)	0	100	100
3	r	167/221 (76%)	167 (100%)	0	100	100
3	t	167/221 (76%)	167 (100%)	0	100	100
3	v	167/221 (76%)	167 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	x	167/221 (76%)	167 (100%)	0	100	100
All	All	13971/17259 (81%)	13971 (100%)	0	100	100

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

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 37 such sidechains are listed below:

Mol	Chain	Res	Type
2	q	47	ASN
2	7	47	ASN
2	s	47	ASN
2	z	47	ASN
1	H	357	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	K	1
1	L	1
1	M	1
1	J	1
1	N	1
1	O	1
1	I	1
1	A	1
1	H	1
1	B	1
1	G	1
1	C	1
1	F	1
1	D	1
1	E	1

The worst 5 of 15 chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	K	171:ASP	C	172:GLY	N	24.97
1	L	171:ASP	C	172:GLY	N	24.97
1	M	171:ASP	C	172:GLY	N	24.97
1	J	171:ASP	C	172:GLY	N	24.95
1	N	171:ASP	C	172:GLY	N	24.95

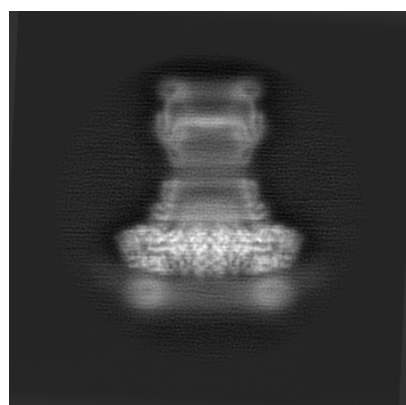
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-8400. These allow visual inspection of the internal detail of the map and identification of artifacts.

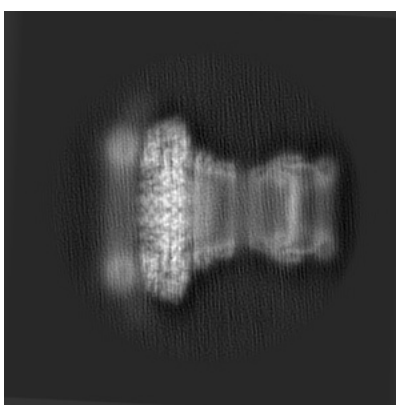
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

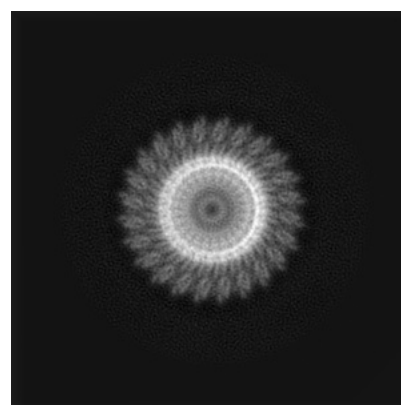
6.1.1 Primary map



X



Y

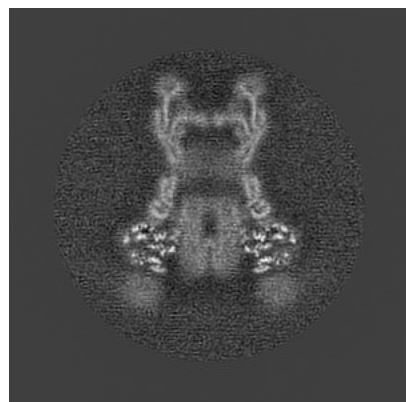


Z

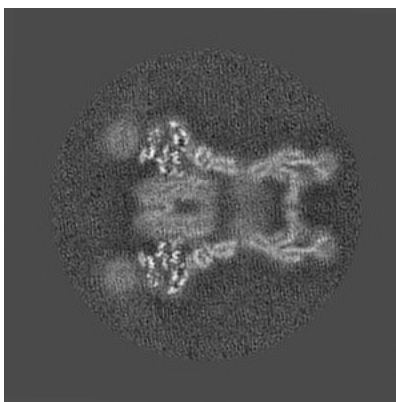
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

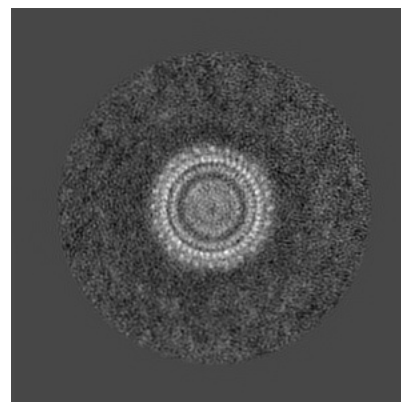
6.2.1 Primary map



X Index: 150



Y Index: 150

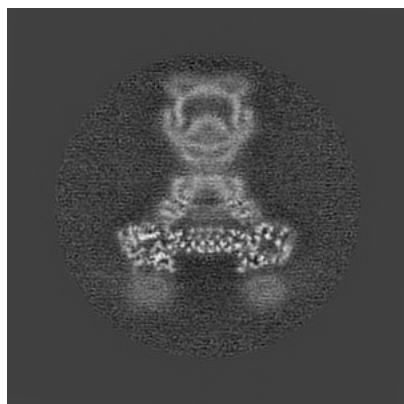


Z Index: 150

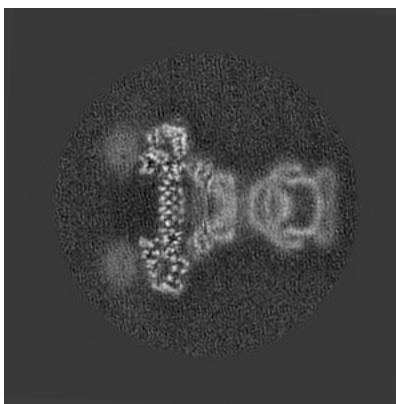
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

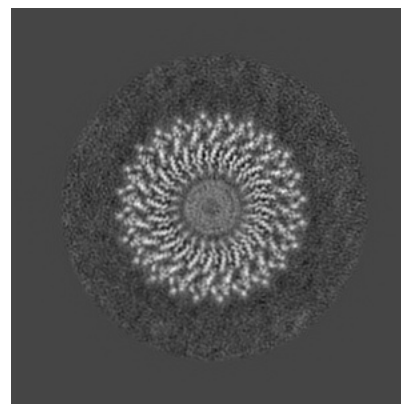
6.3.1 Primary map



X Index: 176



Y Index: 177

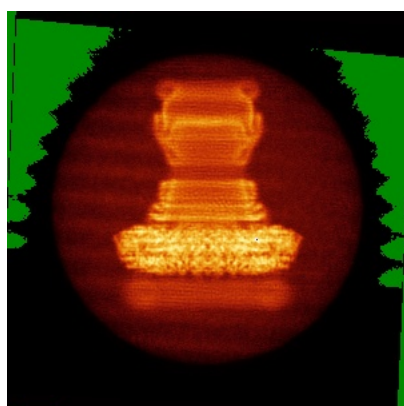


Z Index: 127

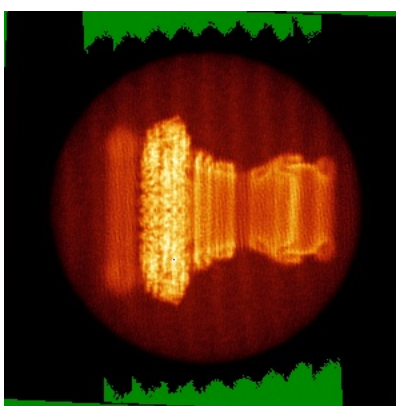
The images above show the largest variance slices of the map in three orthogonal directions.

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

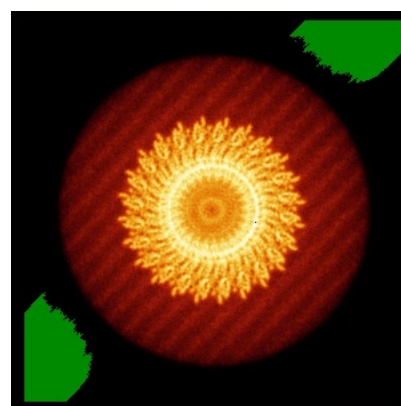
6.4.1 Primary map



X



Y



Z

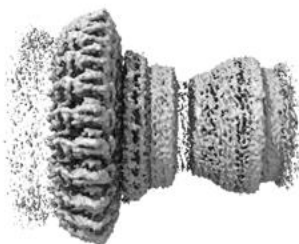
The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



X



Y



Z

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

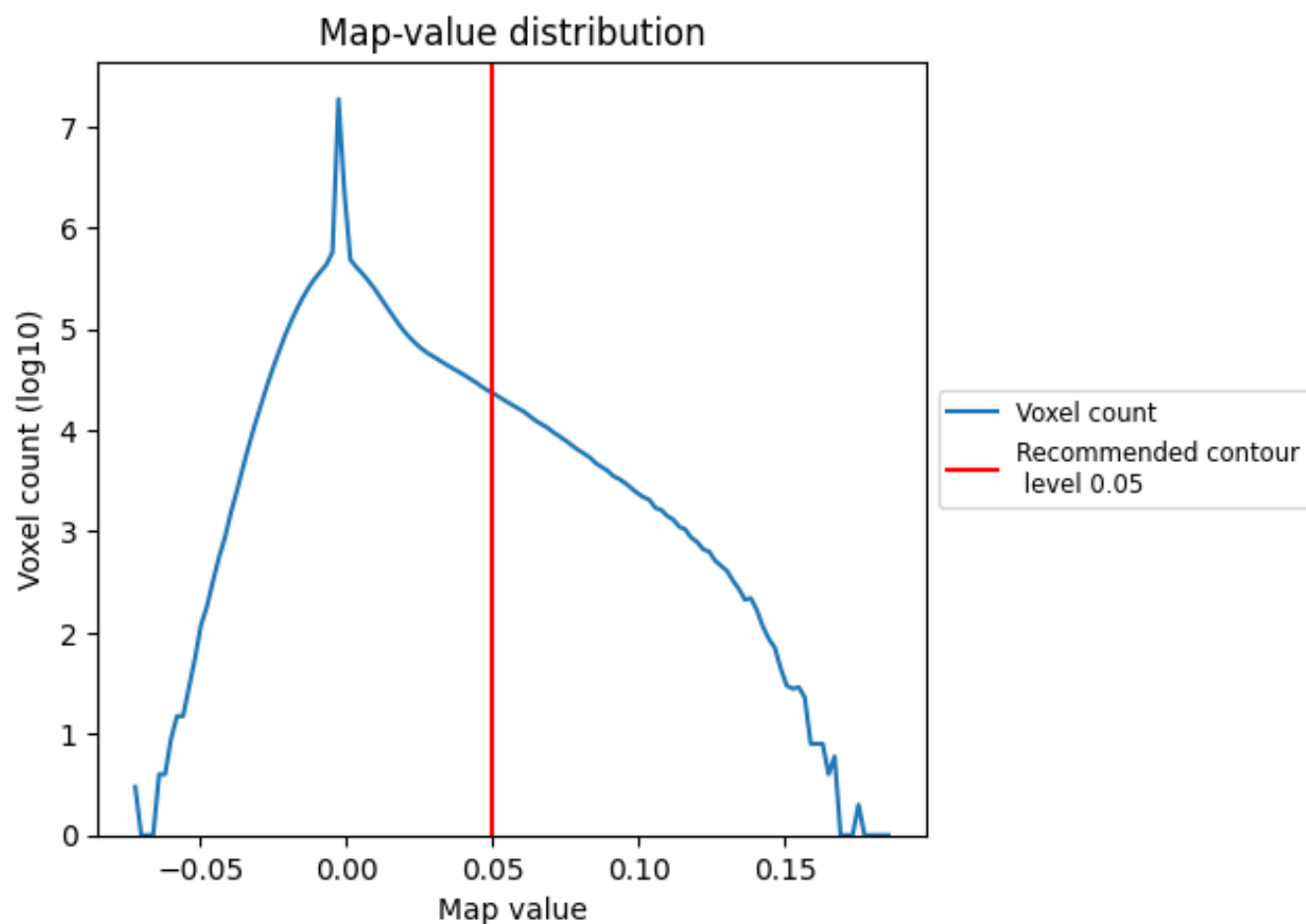
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

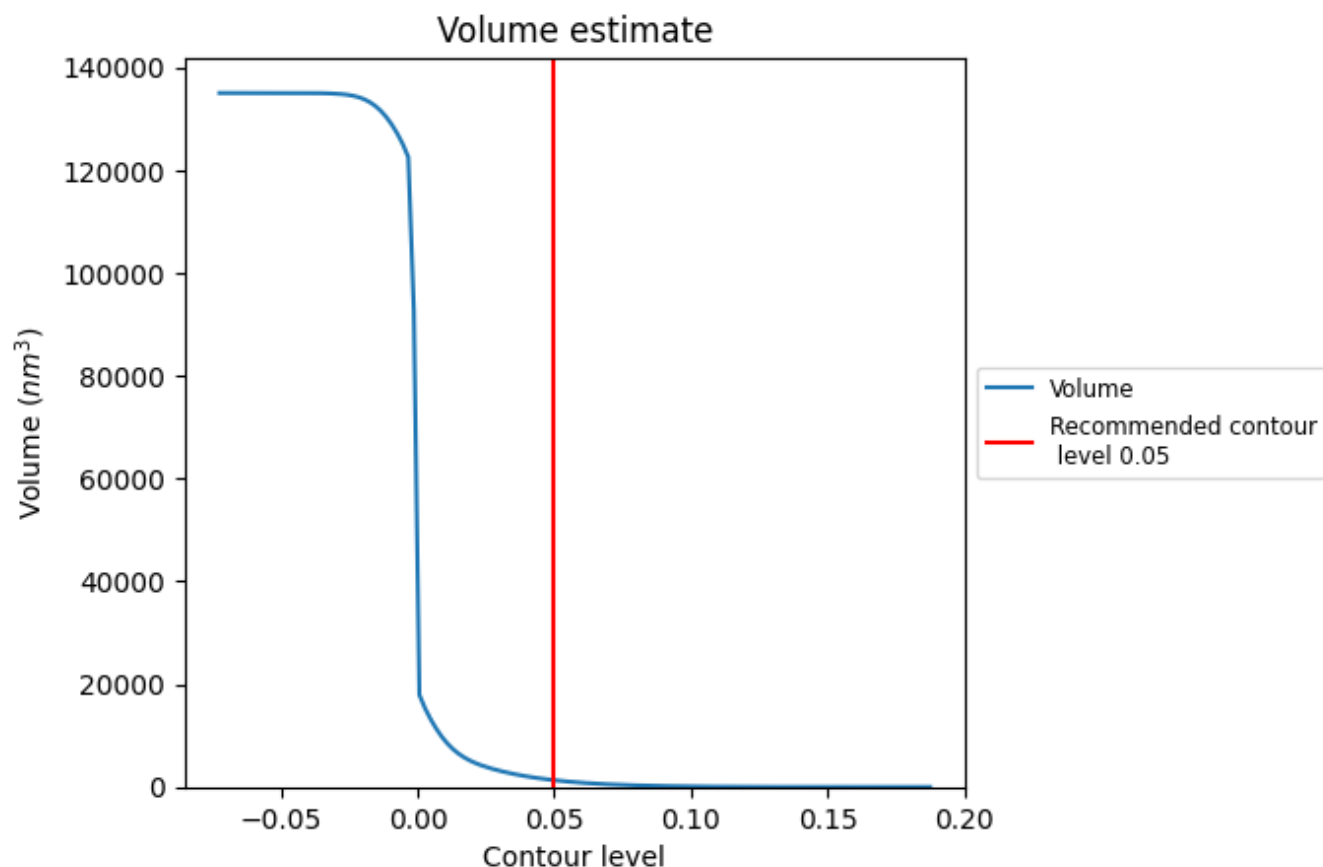
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

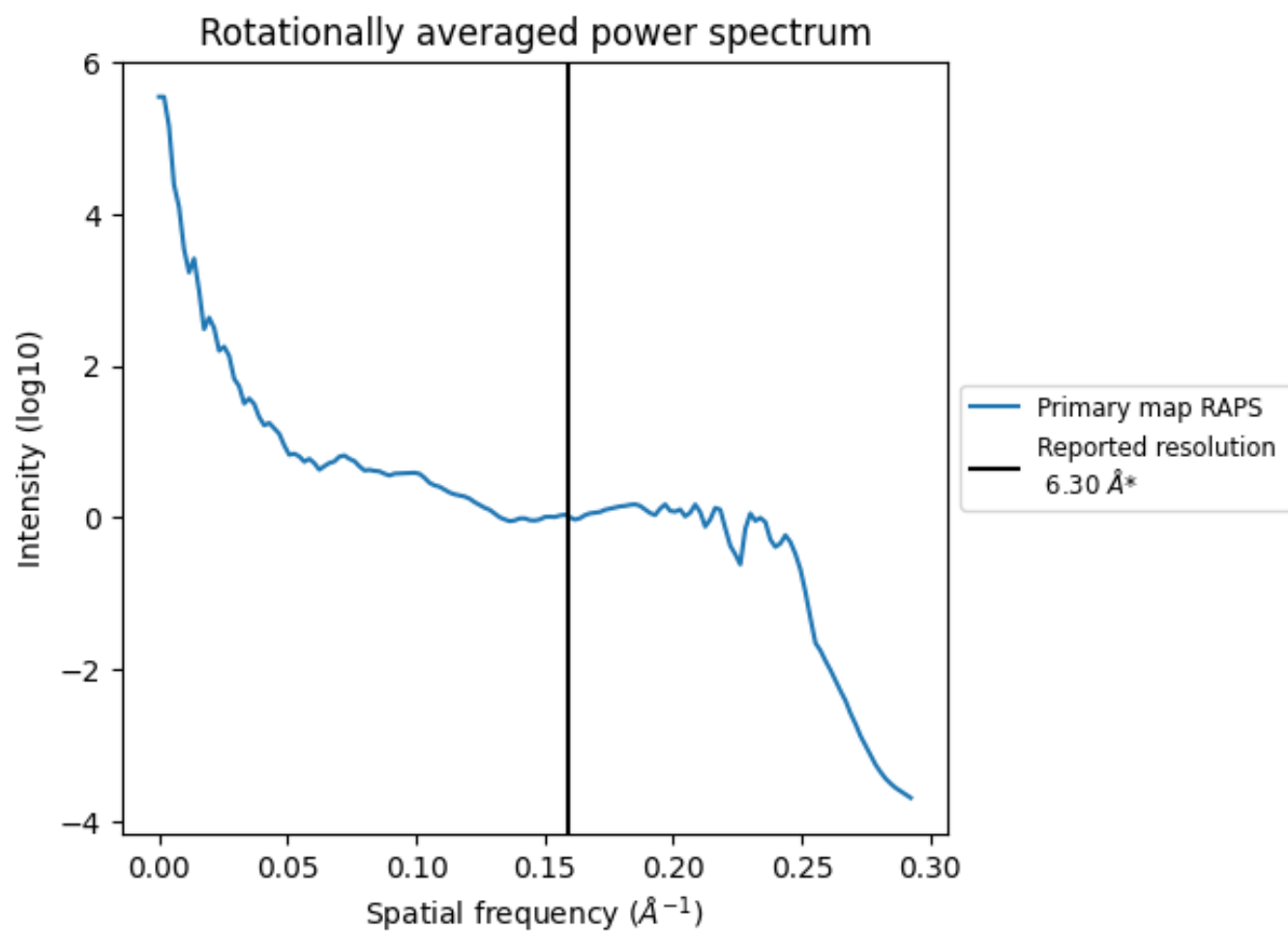
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 1316 nm^3 ; this corresponds to an approximate mass of 1189 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 0.159 Å⁻¹

8 Fourier-Shell correlation

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

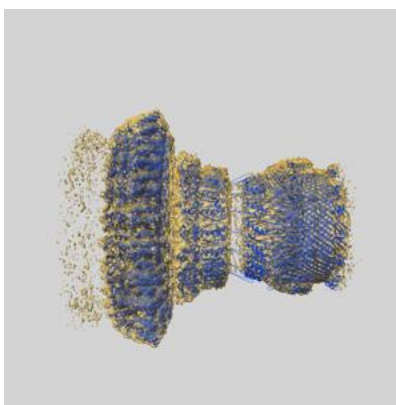
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-8400 and PDB model 5TCR. Per-residue inclusion information can be found in section [3](#) on page [10](#).

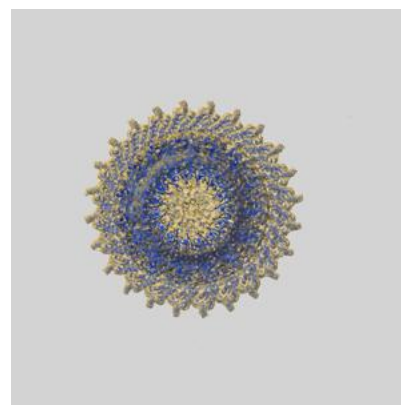
9.1 Map-model overlay [i](#)



X



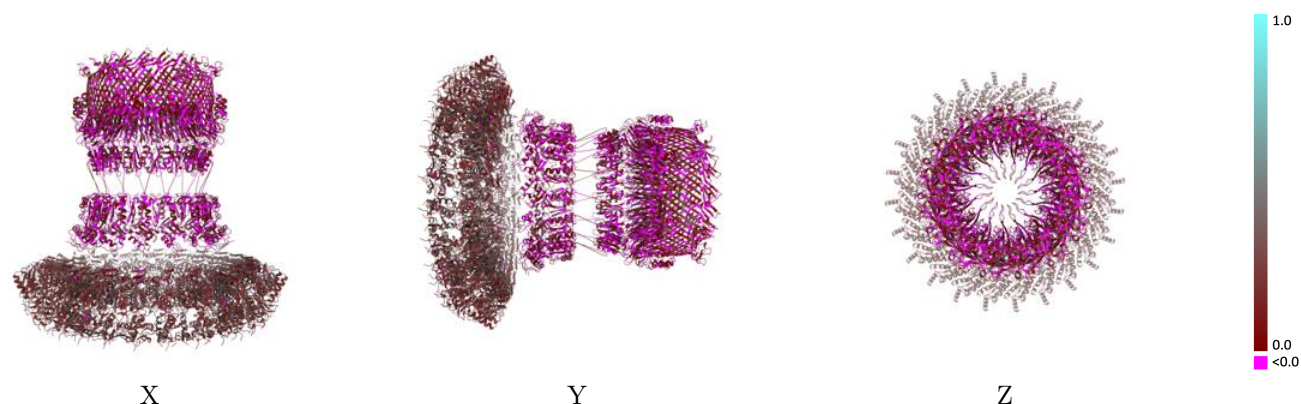
Y



Z

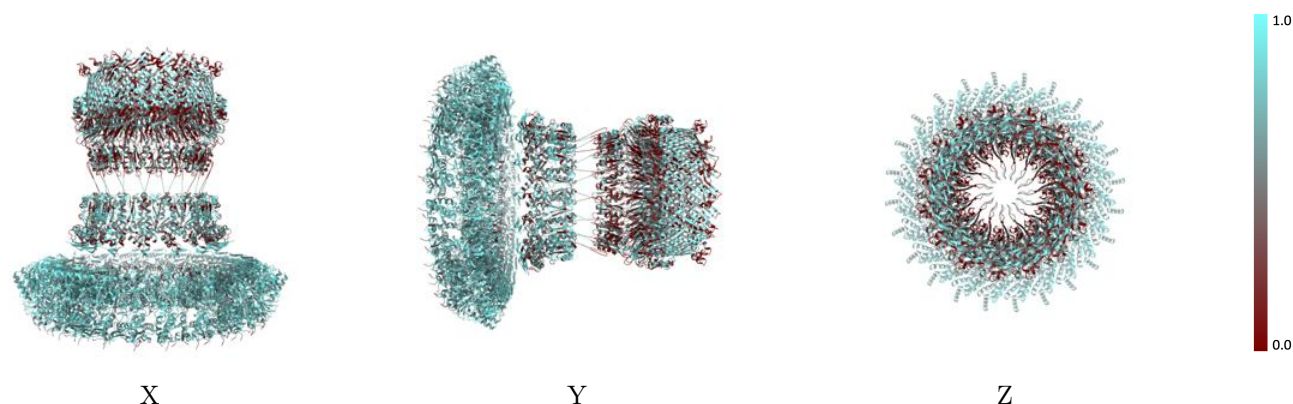
The images above show the 3D surface view of the map at the recommended contour level 0.05 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



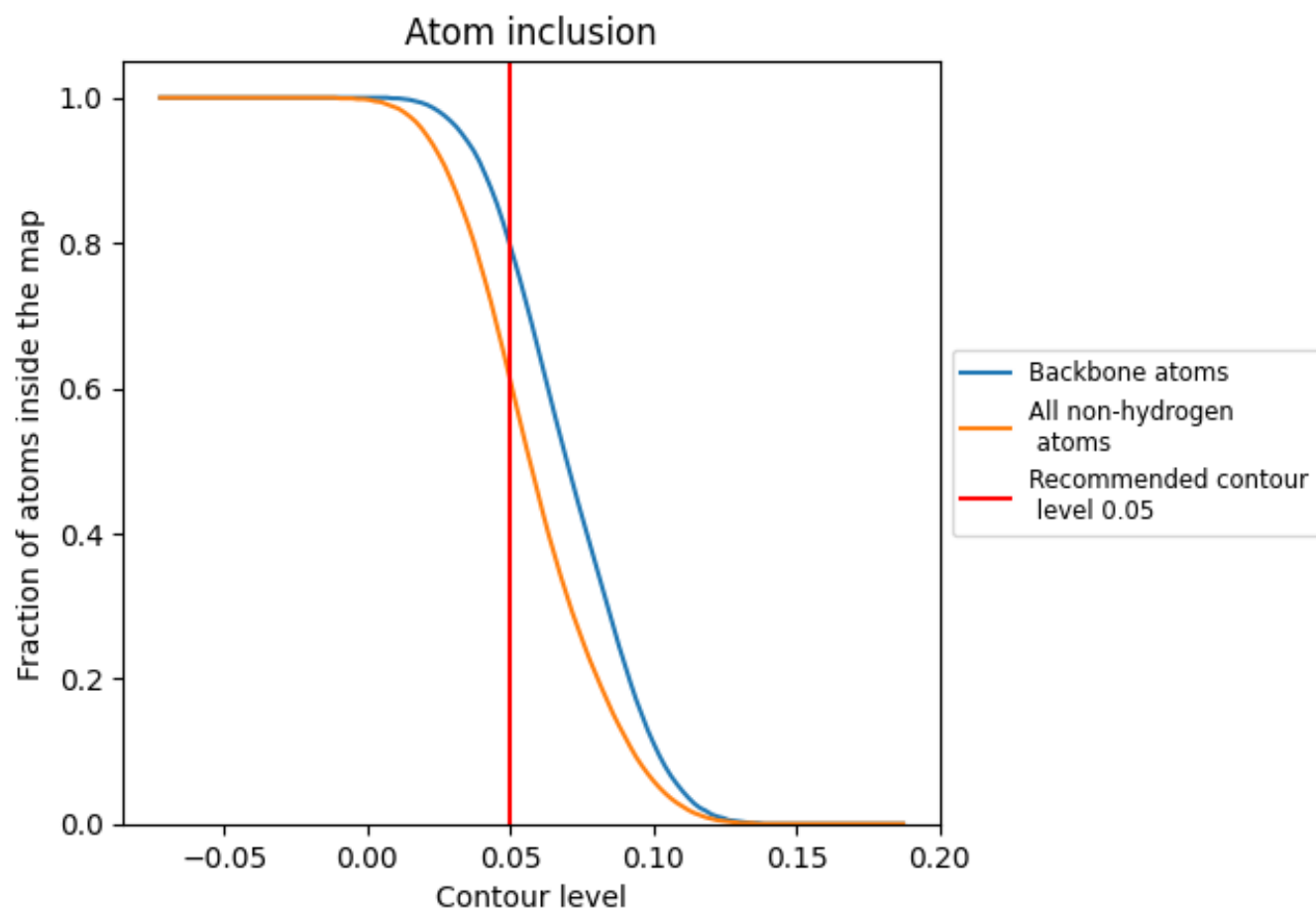
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions at which atoms can be resolved. Low Q-score values may therefore be expected for many entries.

9.3 Atom inclusion mapped to coordinate model [i](#)



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.05).

9.4 Atom inclusion ⓘ



At the recommended contour level, 80% of all backbone atoms, 61% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary ⓘ

The table lists the average atom inclusion at the recommended contour level (0.05) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.6110	 0.1790
0	 0.7240	 0.2520
1	 0.7030	 0.2960
10	 0.7410	 0.2850
2	 0.7410	 0.2650
3	 0.7030	 0.3050
4	 0.7460	 0.2830
5	 0.7170	 0.3110
6	 0.7420	 0.2740
7	 0.7080	 0.3070
8	 0.7510	 0.2980
9	 0.7070	 0.3060
A	 0.4940	 0.0620
B	 0.4970	 0.0670
C	 0.4650	 0.0480
D	 0.4580	 0.0340
E	 0.4620	 0.0460
F	 0.4720	 0.0480
G	 0.4760	 0.0680
H	 0.4750	 0.0570
I	 0.4720	 0.0450
J	 0.4770	 0.0480
K	 0.4680	 0.0420
L	 0.4580	 0.0340
M	 0.4600	 0.0430
N	 0.4730	 0.0410
O	 0.4760	 0.0480
P	 0.7010	 0.2960
Q	 0.7550	 0.2670
R	 0.7160	 0.3040
S	 0.7370	 0.2600
T	 0.7180	 0.3010
U	 0.7260	 0.2480
V	 0.7070	 0.2960
W	 0.7400	 0.2620



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Chain	Atom inclusion	Q-score
X	 0.7130	 0.2990
Y	 0.7270	 0.2470
Z	 0.7350	 0.2750
a	 0.6980	 0.2930
b	 0.7280	 0.2670
c	 0.6880	 0.2960
d	 0.7230	 0.2530
e	 0.6810	 0.2740
f	 0.7130	 0.2510
g	 0.6980	 0.2800
h	 0.7000	 0.2370
i	 0.6810	 0.2720
j	 0.7090	 0.2340
k	 0.6890	 0.2830
l	 0.7000	 0.2350
m	 0.6910	 0.2820
n	 0.7050	 0.2460
o	 0.6950	 0.2940
p	 0.7240	 0.2530
q	 0.7080	 0.2990
r	 0.7440	 0.2740
s	 0.7130	 0.3050
t	 0.7410	 0.2690
u	 0.7180	 0.3020
v	 0.7510	 0.2800
w	 0.7260	 0.3070
x	 0.7550	 0.2790
y	 0.7160	 0.3100
z	 0.7090	 0.2970