



## Full wwPDB EM Validation Report ⓘ

Mar 20, 2024 – 05:25 PM JST

PDB ID : 7E80  
EMDB ID : EMD-31006  
Title : Cryo-EM structure of the flagellar rod with hook and export apparatus from Salmonella  
Authors : Tan, J.X.; Chang, S.H.; Wang, X.F.; Xu, C.H.; Zhou, Y.; Zhang, X.; Zhu, Y.Q.  
Deposited on : 2021-02-28  
Resolution : 3.67 Å(reported)

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

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

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

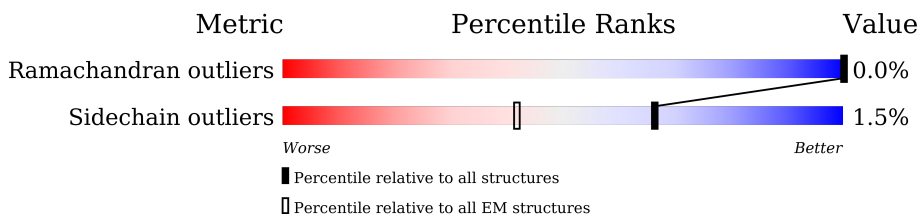
EMDB validation analysis : 0.0.1.dev70  
MolProbity : 4.02b-467  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
MapQ : 1.9.13  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP)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

# 1 Overall quality at a glance

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

The reported resolution of this entry is 3.67 Å.

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



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

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

Mol	Chain	Length	Quality of chain
1	A	260	98%
1	B	260	99%
1	C	260	99%
1	D	260	100%
1	E	260	98%
1	F	260	100%
1	G	260	98%
1	H	260	99%
1	I	260	98%

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Mol	Chain	Length	Quality of chain
1	J	260	99%
1	K	260	98%
1	L	260	98%
1	M	260	99%
1	N	260	96%
1	O	260	97%
1	P	260	95% . 5%
1	Q	260	93% . 5%
1	R	260	94% . .
1	S	260	94% . 5%
1	T	260	97%
1	U	260	99%
1	V	260	98%
1	W	260	98%
1	X	260	6% 98%
2	a	251	97%
2	b	251	9% 97%
2	c	251	97%
2	d	251	98%
2	e	251	6% 98%
3	0	15	13% 100%
3	1	15	20% 100%
3	2	15	100%
3	3	15	100%
3	4	15	20% 100%

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Mol	Chain	Length	Quality of chain
4	5	21	10% 100%
4	6	21	38% 100%
4	7	21	24% 100%
4	8	21	19% 100%
4	9	21	19% 100%
5	f	134	93% • •
5	g	134	9% 95% • •
5	h	134	95% • •
5	i	134	5% 95% • •
5	j	134	5% 94% • 5%
5	p	134	10% 96% •
6	k	138	75% • 22%
6	l	138	7% 75% • 23%
6	m	138	5% 76% • 23%
6	n	138	75% • 22%
6	o	138	9% 76% • 21%
7	q	104	7% 67% • 32%
7	r	104	9% 67% 33%
7	s	104	8% 69% 31%
7	t	104	6% 69% 31%
7	u	104	9% 69% 31%
7	v	104	12% 36% • 63%
8	DA	403	46% 98% •
8	DB	403	33% 99% •
8	DC	403	36% 99%

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Mol	Chain	Length	Quality of chain
8	DD	403	43% 99%
8	DE	403	43% 99%
8	DF	403	45% 99%
8	DG	403	47% 99%
8	DH	403	48% 98%
8	DI	403	50% 98%
8	DJ	403	59% 98%
8	DK	403	60% 99%
9	CE	264	71% 97%
10	CA	89	73% 99%
10	CB	89	81% 98%
10	CC	89	88% 98%
10	CD	89	96% 100%
11	CF	245	39% 85% 14%
11	w	245	26% 84% 14%
11	x	245	43% 84% 14%
11	y	245	26% 82% 14%
11	z	245	28% 85% 14%

## 2 Entry composition [i](#)

There are 11 unique types of molecules in this entry. The entry contains 114468 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 Flagellar basal-body rod protein FlgG.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	260	1949	1202	341	400	6	0	0
1	B	260	1949	1202	341	400	6	0	0
1	C	260	1949	1202	341	400	6	0	0
1	D	260	1949	1202	341	400	6	0	0
1	E	260	1949	1202	341	400	6	0	0
1	F	260	1949	1202	341	400	6	0	0
1	G	260	1949	1202	341	400	6	0	0
1	H	260	1949	1202	341	400	6	0	0
1	I	260	1949	1202	341	400	6	0	0
1	J	260	1949	1202	341	400	6	0	0
1	K	260	1949	1202	341	400	6	0	0
1	L	259	1941	1197	340	399	5	0	0
1	M	260	1949	1202	341	400	6	0	0
1	N	251	1887	1167	330	384	6	0	0
1	O	252	1894	1172	331	385	6	0	0
1	P	248	1862	1151	327	379	5	0	0
1	Q	247	1858	1149	326	378	5	0	0

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Mol	Chain	Residues	Atoms					AltConf	Trace
1	R	250	Total	C	N	O	S	0	0
			1875	1159	329	382	5		
1	S	247	Total	C	N	O	S	0	0
			1858	1149	326	378	5		
1	T	253	Total	C	N	O	S	0	0
			1902	1176	333	388	5		
1	U	259	Total	C	N	O	S	0	0
			1941	1197	340	399	5		
1	V	259	Total	C	N	O	S	0	0
			1941	1197	340	399	5		
1	W	259	Total	C	N	O	S	0	0
			1941	1197	340	399	5		
1	X	259	Total	C	N	O	S	0	0
			1941	1197	340	399	5		

- Molecule 2 is a protein called Flagellar basal-body rod protein FlgF.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	a	249	Total	C	N	O	S	0	0
			1812	1111	325	368	8		
2	b	248	Total	C	N	O	S	0	0
			1804	1106	324	367	7		
2	c	249	Total	C	N	O	S	0	0
			1812	1111	325	368	8		
2	d	249	Total	C	N	O	S	0	0
			1812	1111	325	368	8		
2	e	249	Total	C	N	O	S	0	0
			1812	1111	325	368	8		

- Molecule 3 is a protein called Flagellar MS ring L2.

Mol	Chain	Residues	Atoms				AltConf	Trace
3	0	15	Total	C	N	O	0	0
			75	45	15	15		
3	1	15	Total	C	N	O	0	0
			75	45	15	15		
3	2	15	Total	C	N	O	0	0
			75	45	15	15		
3	3	15	Total	C	N	O	0	0
			75	45	15	15		
3	4	15	Total	C	N	O	0	0
			75	45	15	15		

- Molecule 4 is a protein called Flagellar MS ring L1.

Mol	Chain	Residues	Atoms				AltConf	Trace
4	5	21	Total	C	N	O	0	0
			140	88	24	28		
4	6	21	Total	C	N	O	0	0
			140	88	24	28		
4	7	21	Total	C	N	O	0	0
			140	88	24	28		
4	9	21	Total	C	N	O	0	0
			140	88	24	28		
4	8	21	Total	C	N	O	0	0
			140	88	24	28		

- Molecule 5 is a protein called Flagellar basal-body rod protein FlgC.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	f	128	Total	C	N	O	S	0	0
			936	585	160	186	5		
5	g	130	Total	C	N	O	S	0	0
			949	592	163	189	5		
5	h	128	Total	C	N	O	S	0	0
			935	584	160	186	5		
5	j	127	Total	C	N	O	S	0	0
			931	582	159	185	5		
5	p	129	Total	C	N	O	S	0	0
			940	587	161	187	5		
5	i	129	Total	C	N	O	S	0	0
			944	589	162	188	5		

- Molecule 6 is a protein called Flagellar basal body rod protein FlgB.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	l	106	Total	C	N	O	S	0	0
			833	515	150	163	5		
6	m	106	Total	C	N	O	S	0	0
			833	515	150	163	5		
6	o	109	Total	C	N	O	S	0	0
			856	529	156	166	5		
6	k	108	Total	C	N	O	S	0	0
			852	527	155	165	5		
6	n	107	Total	C	N	O	S	0	0
			844	521	154	164	5		

- Molecule 7 is a protein called Flagellar hook-basal body complex protein FliE.



Mol	Chain	Residues	Atoms					AltConf	Trace
7	q	71	Total	C	N	O	S	0	0
			536	330	98	102	6		
7	r	70	Total	C	N	O	S	0	0
			526	323	97	100	6		
7	s	72	Total	C	N	O	S	0	0
			543	335	99	103	6		
7	t	72	Total	C	N	O	S	0	0
			543	335	99	103	6		
7	u	72	Total	C	N	O	S	0	0
			543	335	99	103	6		
7	v	38	Total	C	N	O	S	0	0
			289	178	50	55	6		

- Molecule 8 is a protein called Flagellar hook protein FlgE.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	DA	401	Total	C	N	O	S	0	0
			2947	1814	507	618	8		
8	DB	401	Total	C	N	O	S	0	0
			2947	1814	507	618	8		
8	DC	401	Total	C	N	O	S	0	0
			2947	1814	507	618	8		
8	DD	401	Total	C	N	O	S	0	0
			2947	1814	507	618	8		
8	DE	401	Total	C	N	O	S	0	0
			2947	1814	507	618	8		
8	DF	401	Total	C	N	O	S	0	0
			2947	1814	507	618	8		
8	DG	401	Total	C	N	O	S	0	0
			2947	1814	507	618	8		
8	DH	401	Total	C	N	O	S	0	0
			2947	1814	507	618	8		
8	DI	401	Total	C	N	O	S	0	0
			2947	1814	507	618	8		
8	DJ	401	Total	C	N	O	S	0	0
			2947	1814	507	618	8		
8	DK	401	Total	C	N	O	S	0	0
			2947	1814	507	618	8		

- Molecule 9 is a protein called Flagellar biosynthetic protein FliR.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	CE	260	Total	C	N	O	S	0	0
			2002	1339	317	330	16		

- Molecule 10 is a protein called Flagellar biosynthetic protein FliQ.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	CA	89	Total	C	N	O	S	0	0
			670	449	100	114	7		
10	CB	89	Total	C	N	O	S	0	0
			670	449	100	114	7		
10	CC	89	Total	C	N	O	S	0	0
			670	449	100	114	7		
10	CD	89	Total	C	N	O	S	0	0
			670	449	100	114	7		

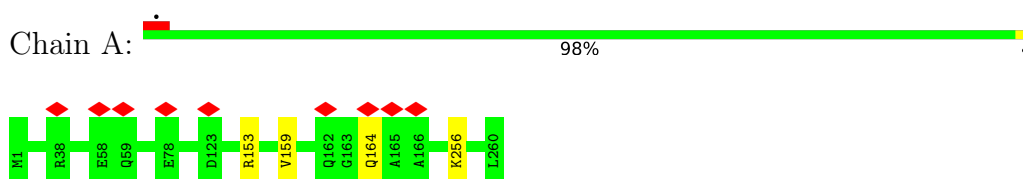
- Molecule 11 is a protein called Flagellar biosynthetic protein FliP.

Mol	Chain	Residues	Atoms					AltConf	Trace
11	w	211	Total	C	N	O	S	0	0
			1636	1091	254	279	12		
11	x	211	Total	C	N	O	S	0	0
			1636	1091	254	279	12		
11	y	211	Total	C	N	O	S	0	0
			1636	1091	254	279	12		
11	z	211	Total	C	N	O	S	0	0
			1636	1091	254	279	12		
11	CF	211	Total	C	N	O	S	0	0
			1636	1091	254	279	12		

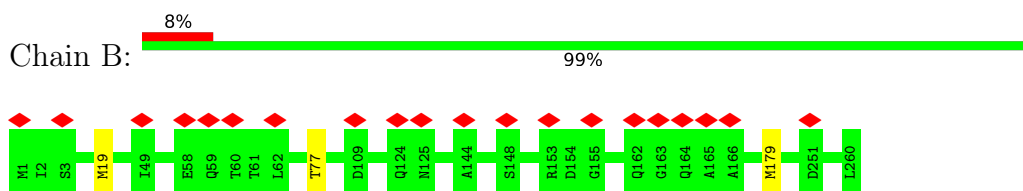
### 3 Residue-property plots [i](#)

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

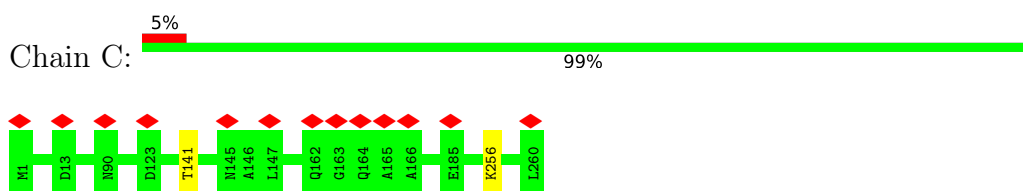
- Molecule 1: Flagellar basal-body rod protein FlgG



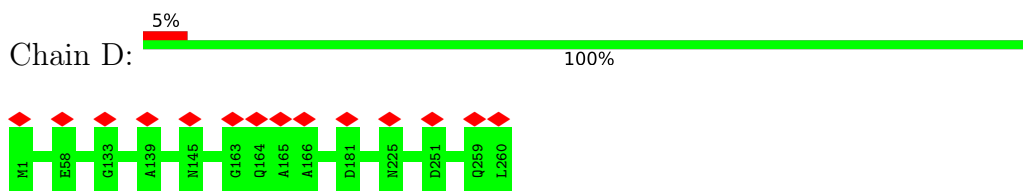
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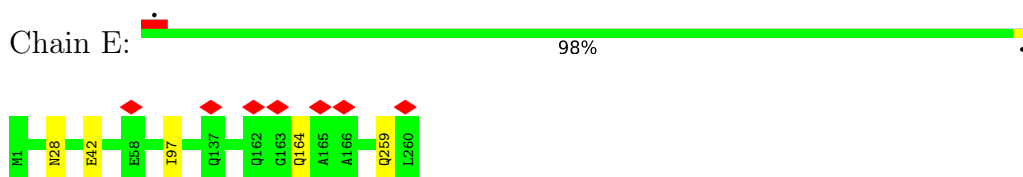
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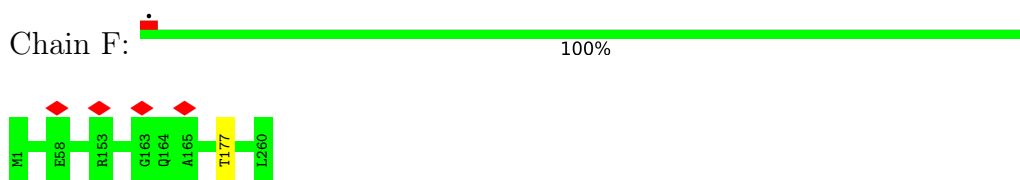
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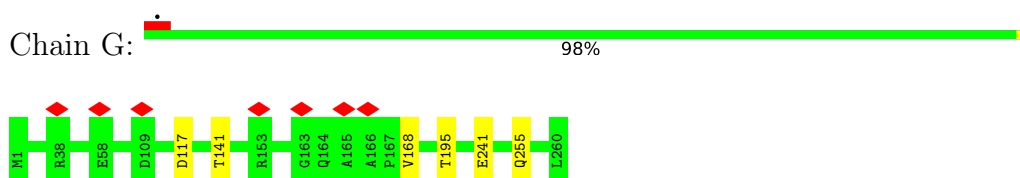
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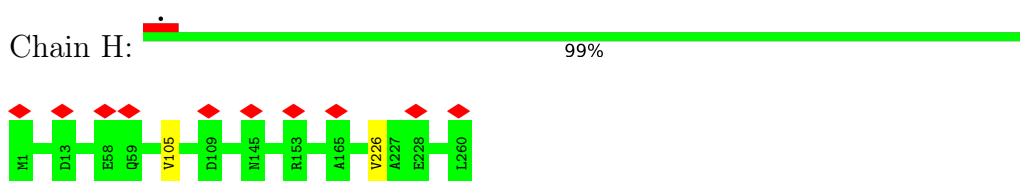
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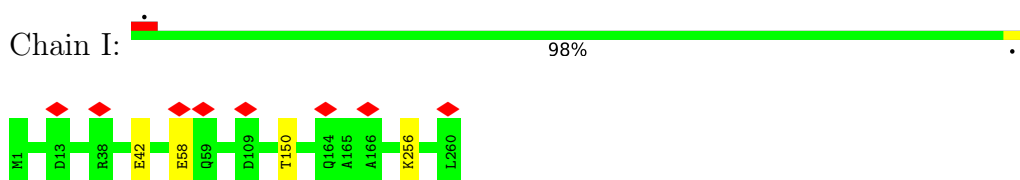
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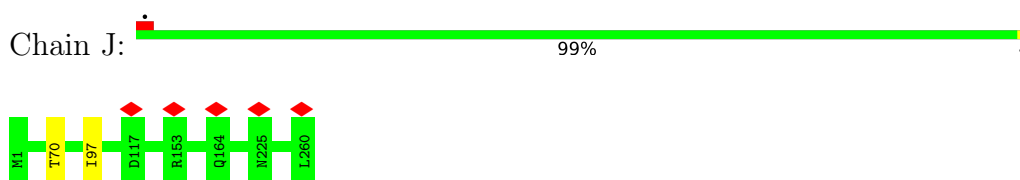
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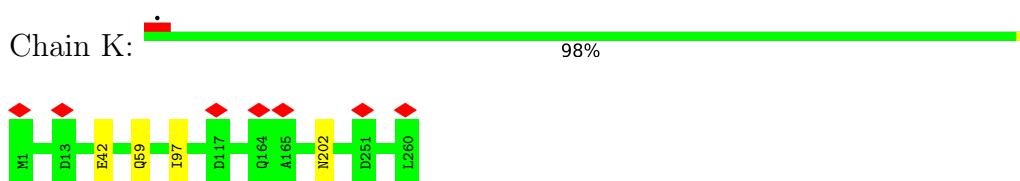
- Molecule 1: Flagellar basal-body rod protein FlgG



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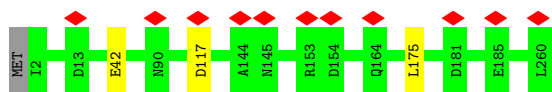


- Molecule 1: Flagellar basal-body rod protein FlgG

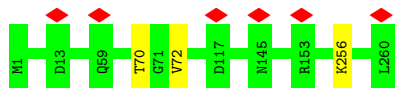


- Molecule 1: Flagellar basal-body rod protein FlgG

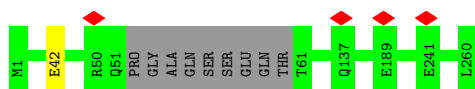




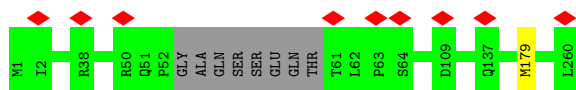
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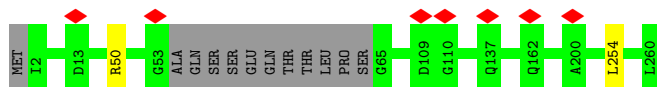
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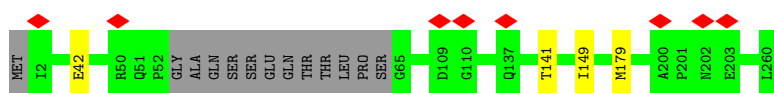
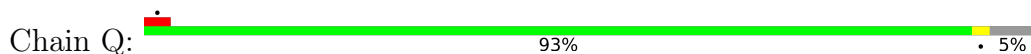
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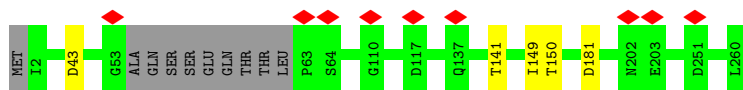
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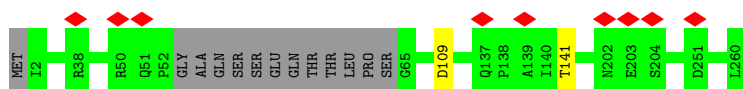


- Molecule 1: Flagellar basal-body rod protein FlgG



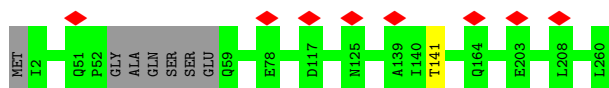
- Molecule 1: Flagellar basal-body rod protein FlgG

Chain S:  94% 5%



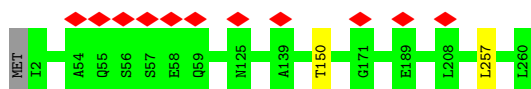
- Molecule 1: Flagellar basal-body rod protein FlgG

Chain T:  97%



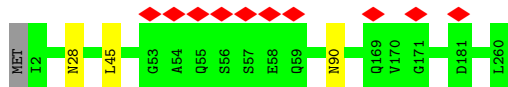
- Molecule 1: Flagellar basal-body rod protein FlgG

Chain U:  99%



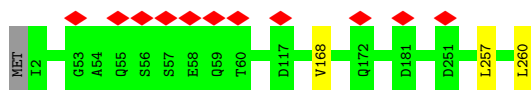
- Molecule 1: Flagellar basal-body rod protein FlgG

Chain V:  98%



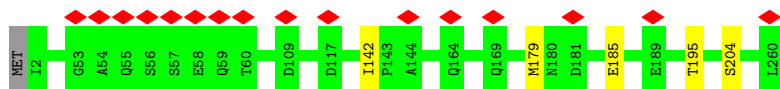
- Molecule 1: Flagellar basal-body rod protein FlgG

Chain W:  98%



- Molecule 1: Flagellar basal-body rod protein FlgG

Chain X:  6% 98%

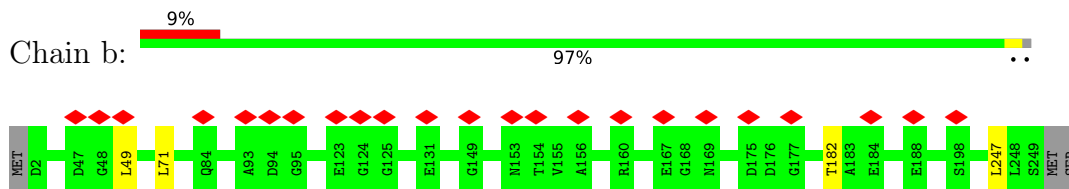


- Molecule 2: Flagellar basal-body rod protein FlgF

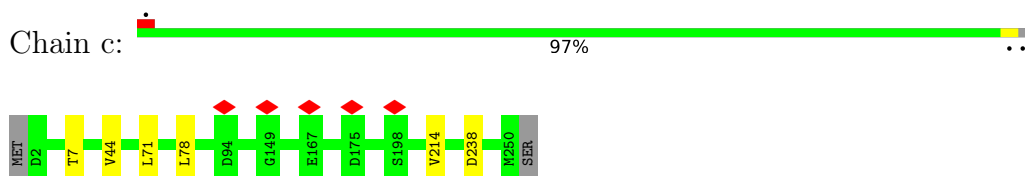
Chain a:  97%



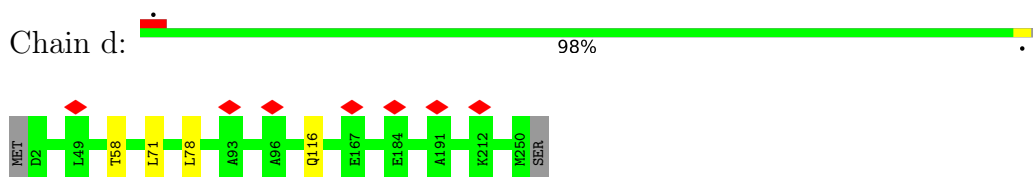
- Molecule 2: Flagellar basal-body rod protein FlgF



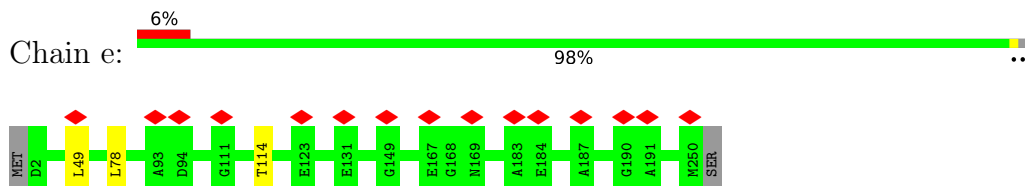
- Molecule 2: Flagellar basal-body rod protein FlgF



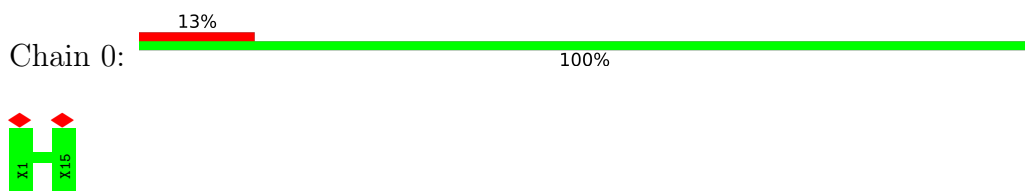
- Molecule 2: Flagellar basal-body rod protein FlgF



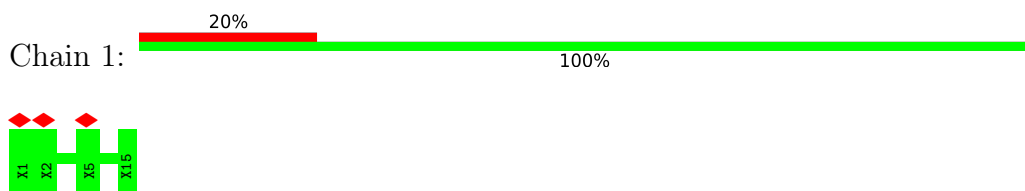
- Molecule 2: Flagellar basal-body rod protein FlgF



- Molecule 3: Flagellar MS ring L2



- Molecule 3: Flagellar MS ring L2



- Molecule 3: Flagellar MS ring L2



There are no outlier residues recorded for this chain.

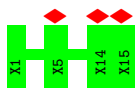
- Molecule 3: Flagellar MS ring L2

Chain 3:  100%

There are no outlier residues recorded for this chain.

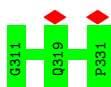
- Molecule 3: Flagellar MS ring L2

Chain 4:  20%  
100%



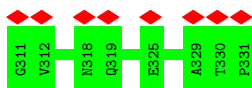
- Molecule 4: Flagellar MS ring L1

Chain 5:  10%  
100%



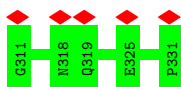
- Molecule 4: Flagellar MS ring L1

Chain 6:  38%  
100%



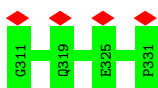
- Molecule 4: Flagellar MS ring L1

Chain 7:  24%  
100%



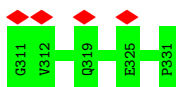
- Molecule 4: Flagellar MS ring L1

Chain 9:  19%  
100%



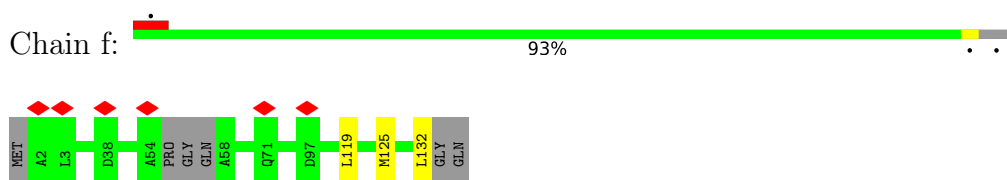
- Molecule 4: Flagellar MS ring L1

Chain 8:  19%  
100%

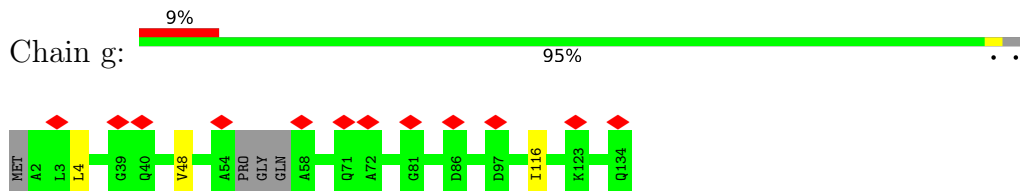




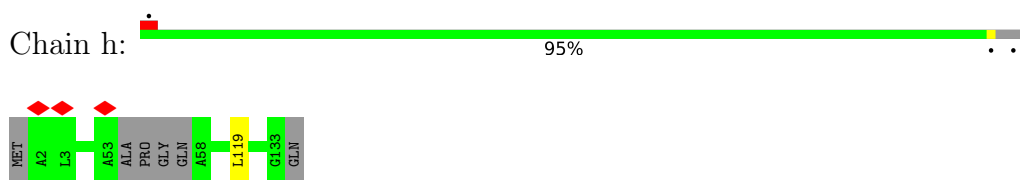
- Molecule 5: Flagellar basal-body rod protein FlgC



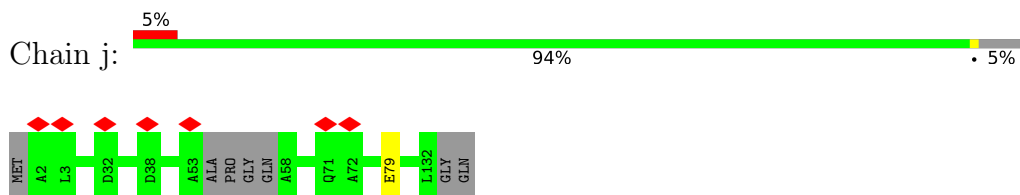
- Molecule 5: Flagellar basal-body rod protein FlgC



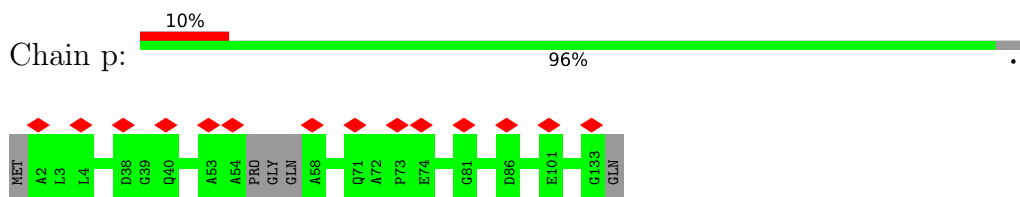
- Molecule 5: Flagellar basal-body rod protein FlgC



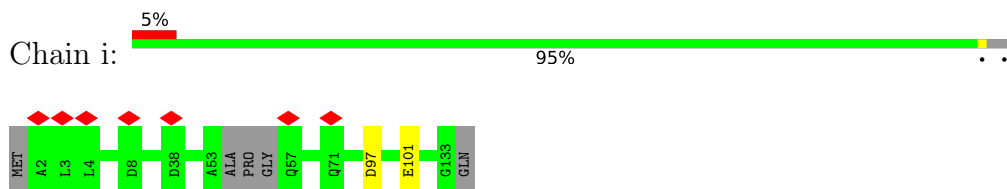
- Molecule 5: Flagellar basal-body rod protein FlgC



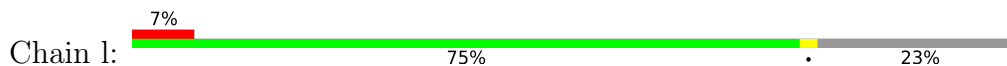
- Molecule 5: Flagellar basal-body rod protein FlgC

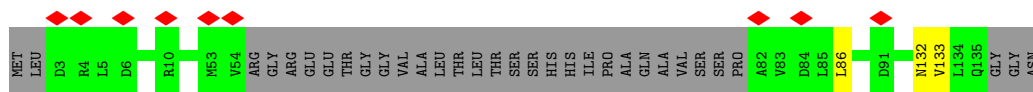


- Molecule 5: Flagellar basal-body rod protein FlgC

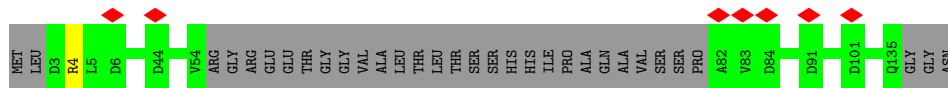
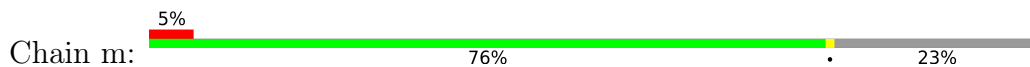


- Molecule 6: Flagellar basal body rod protein FlgB

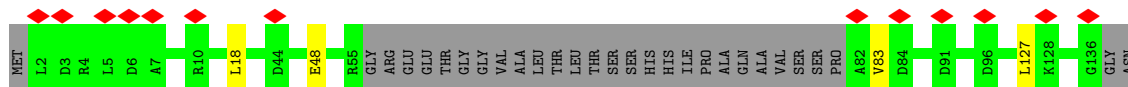
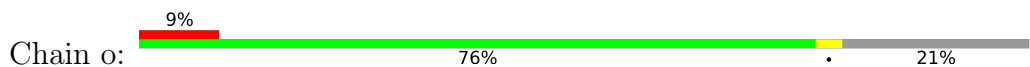




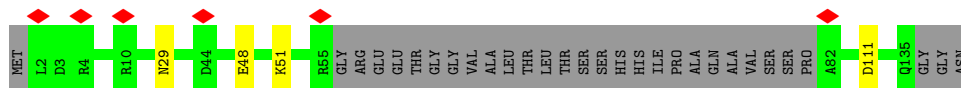
• Molecule 6: Flagellar basal body rod protein FlgB



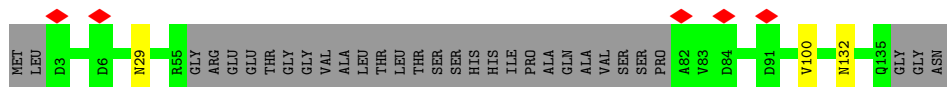
• Molecule 6: Flagellar basal body rod protein FlgB



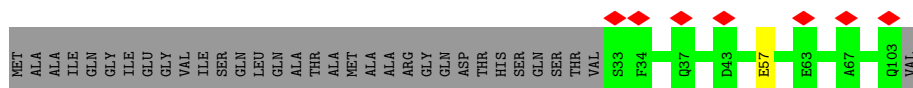
• Molecule 6: Flagellar basal body rod protein FlgB



• Molecule 6: Flagellar basal body rod protein FlgB



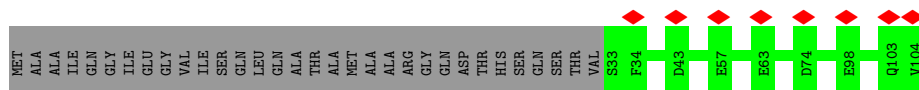
• Molecule 7: Flagellar hook-basal body complex protein FliE



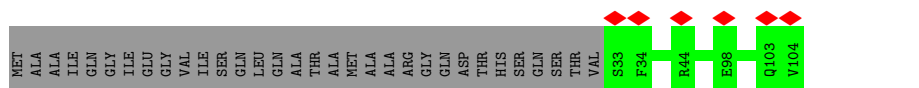
• Molecule 7: Flagellar hook-basal body complex protein FliE



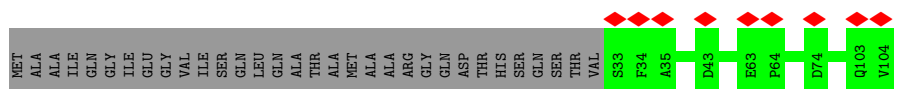
• Molecule 7: Flagellar hook-basal body complex protein FliE



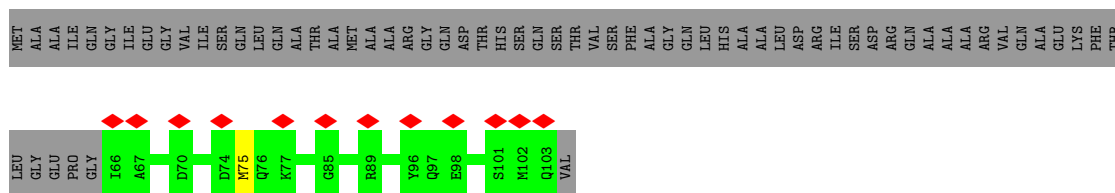
• Molecule 7: Flagellar hook-basal body complex protein FliE



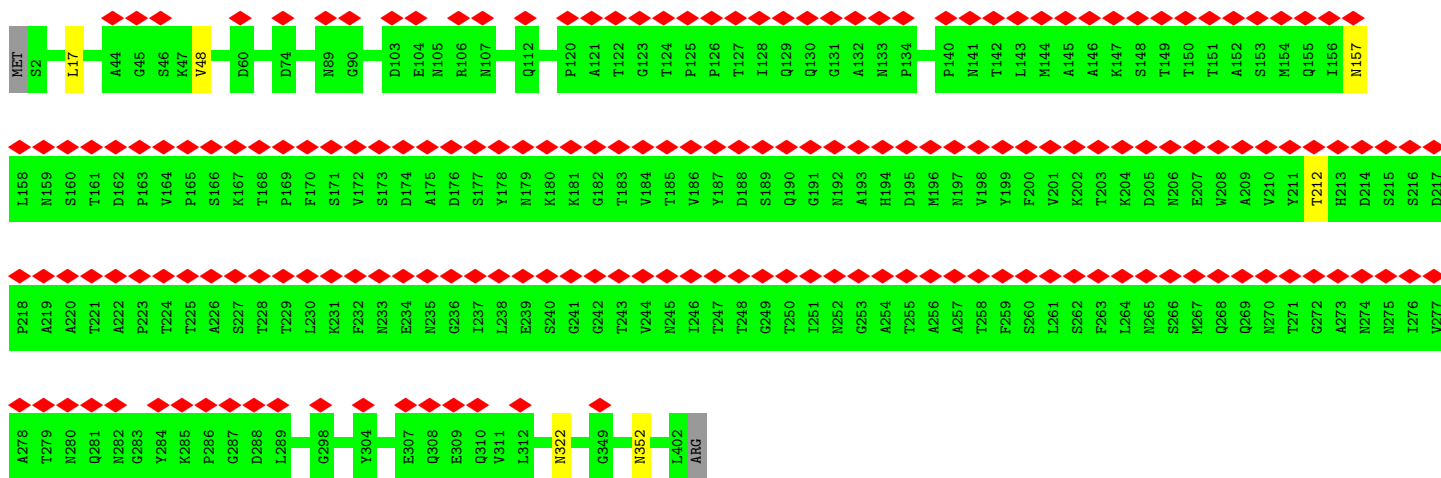
• Molecule 7: Flagellar hook-basal body complex protein FliE



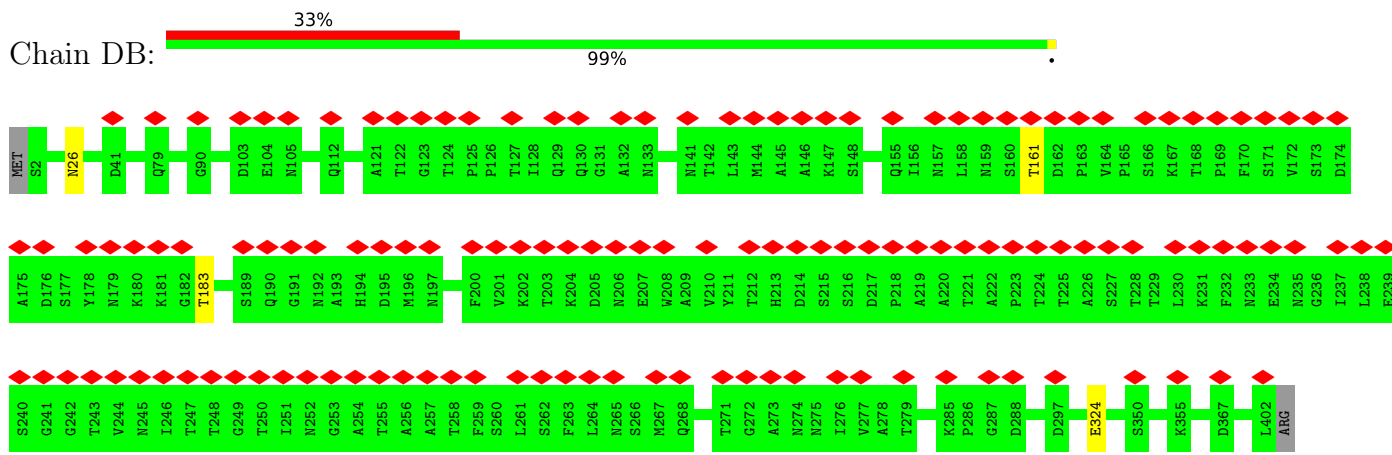
• Molecule 7: Flagellar hook-basal body complex protein FliE



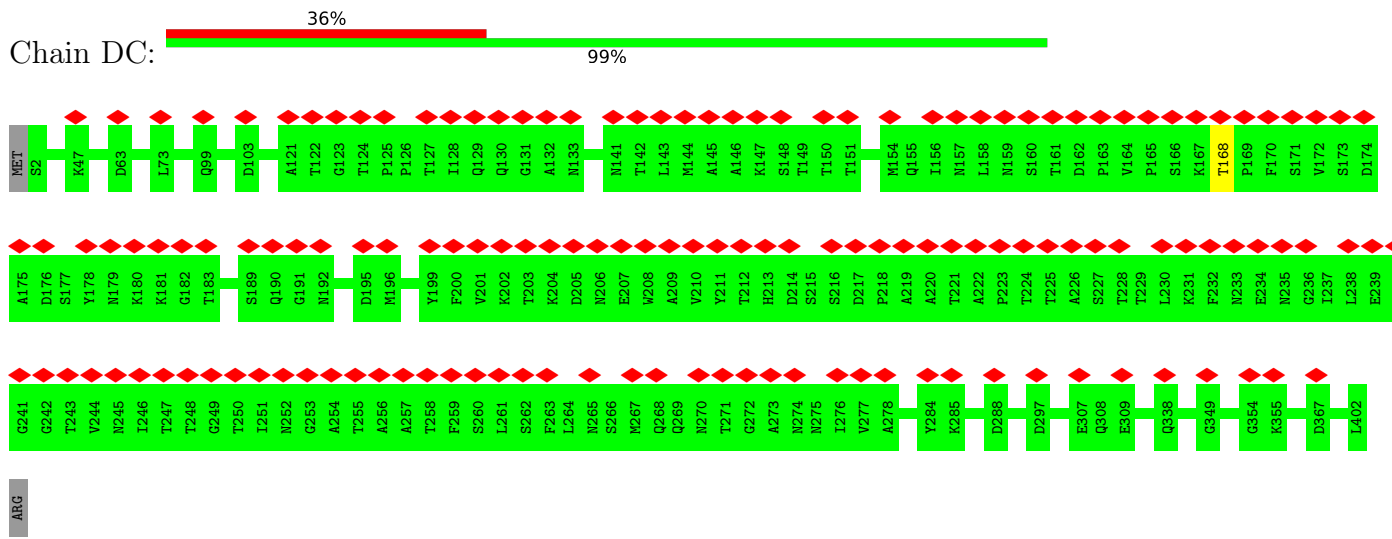
• Molecule 8: Flagellar hook protein FlgE



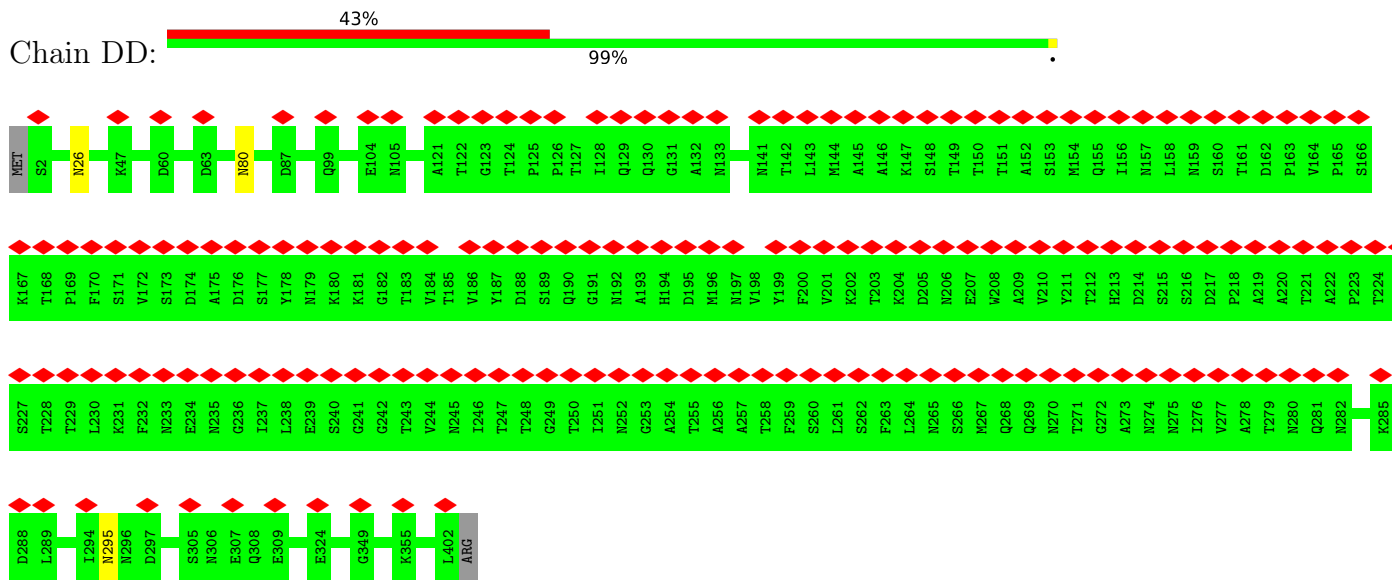
• Molecule 8: Flagellar hook protein FlgE



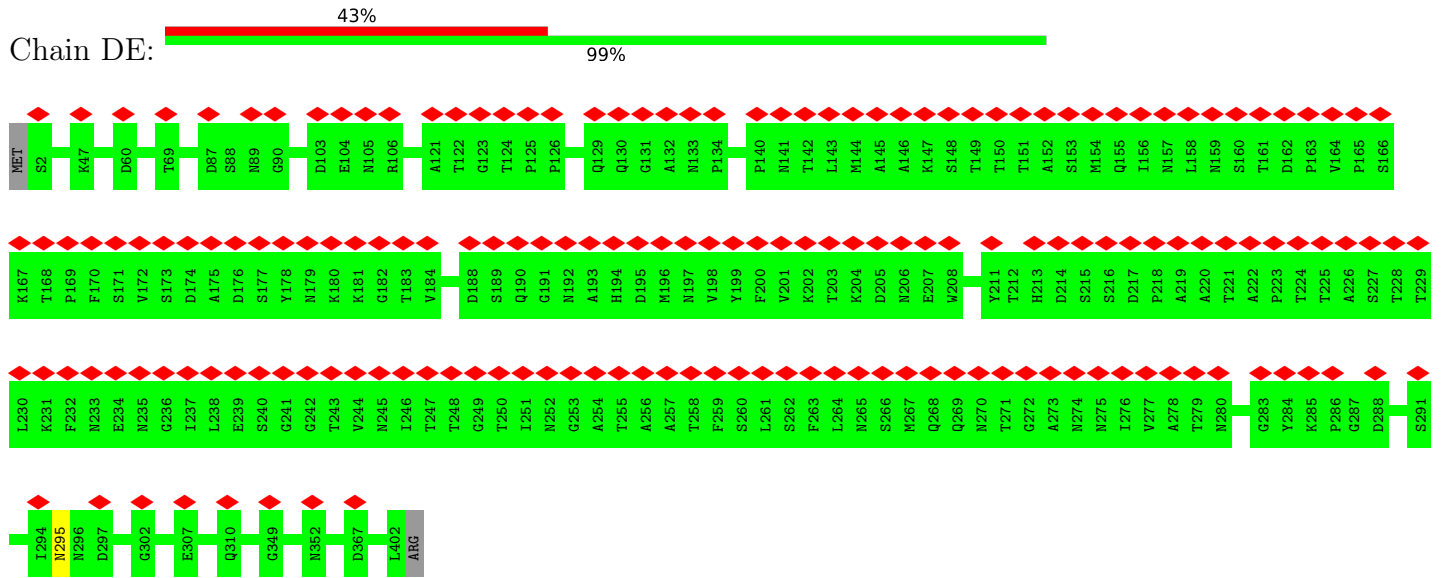
• Molecule 8: Flagellar hook protein FlgE



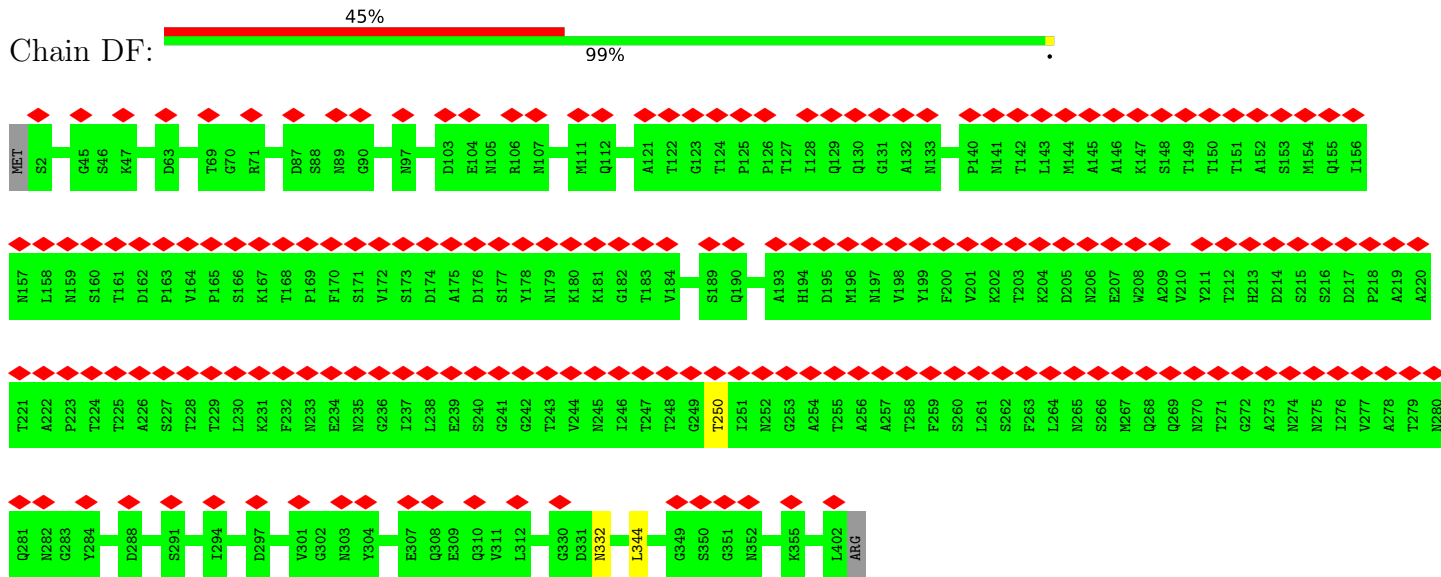
• Molecule 8: Flagellar hook protein FlgE



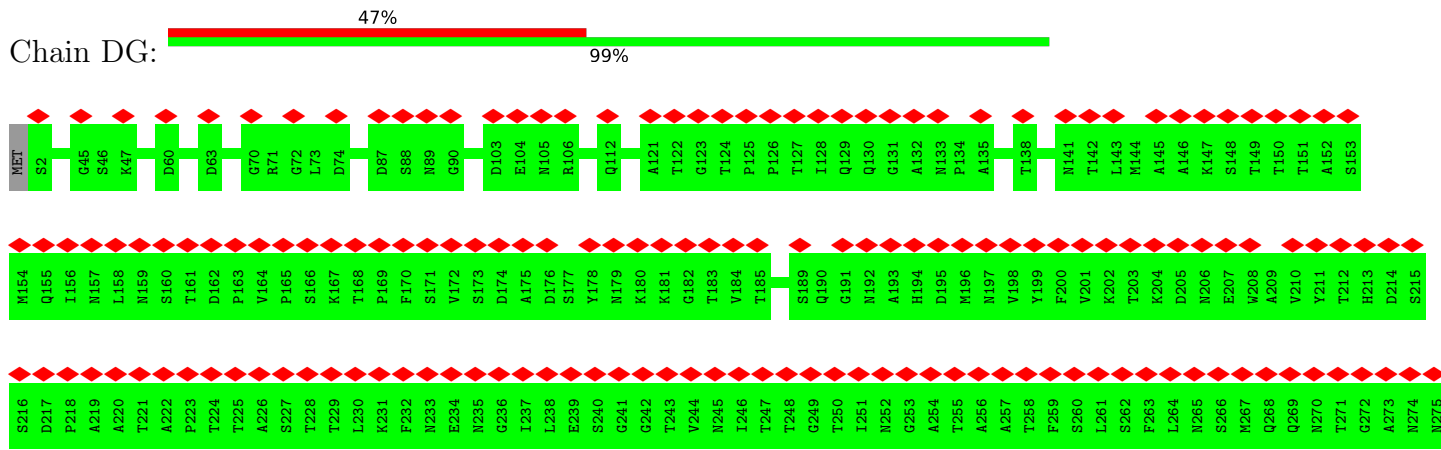
• Molecule 8: Flagellar hook protein FlgE

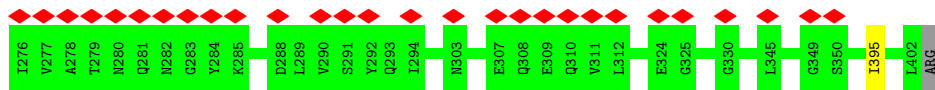


• Molecule 8: Flagellar hook protein FlgE

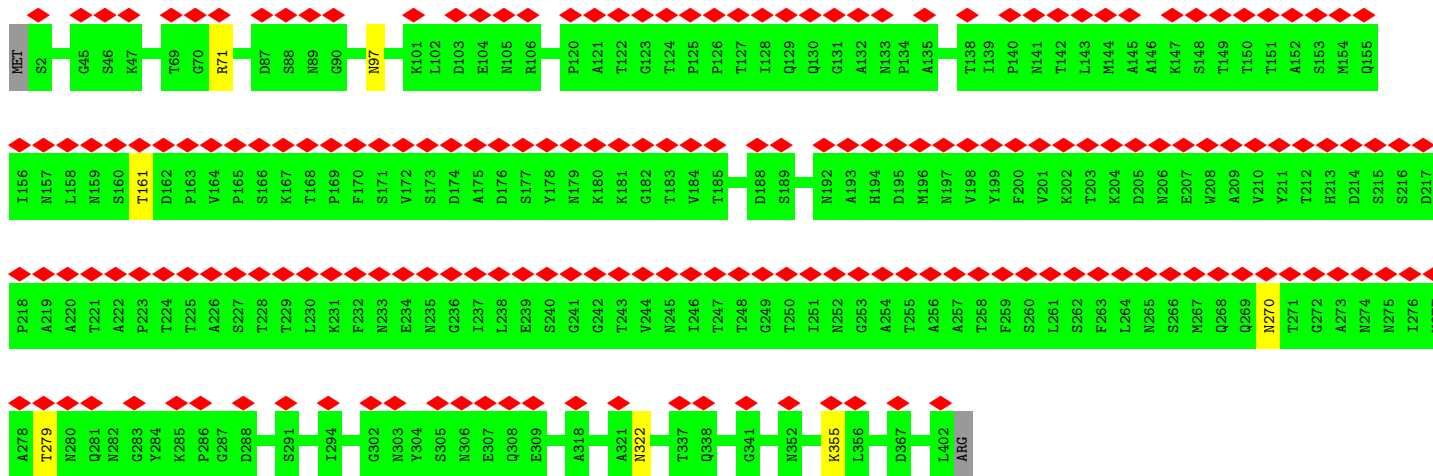


• Molecule 8: Flagellar hook protein FlgE

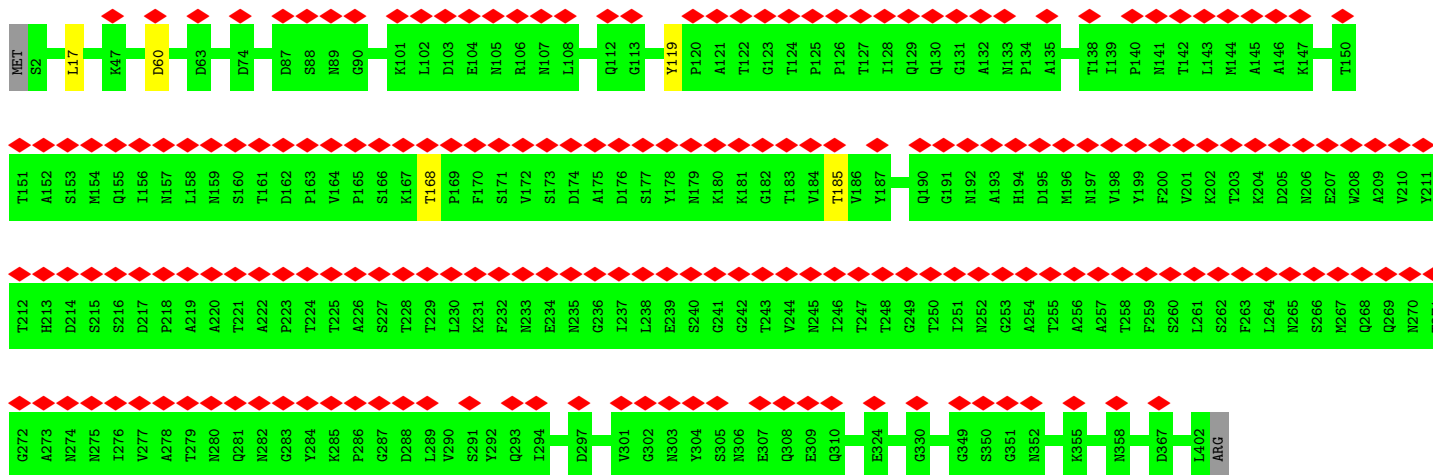




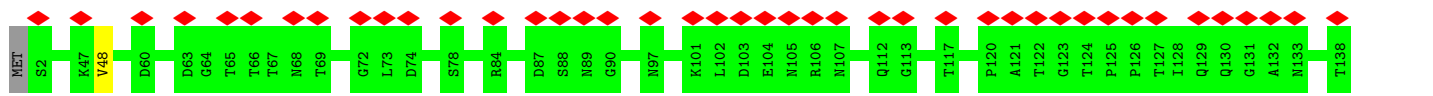
• Molecule 8: Flagellar hook protein FlgE

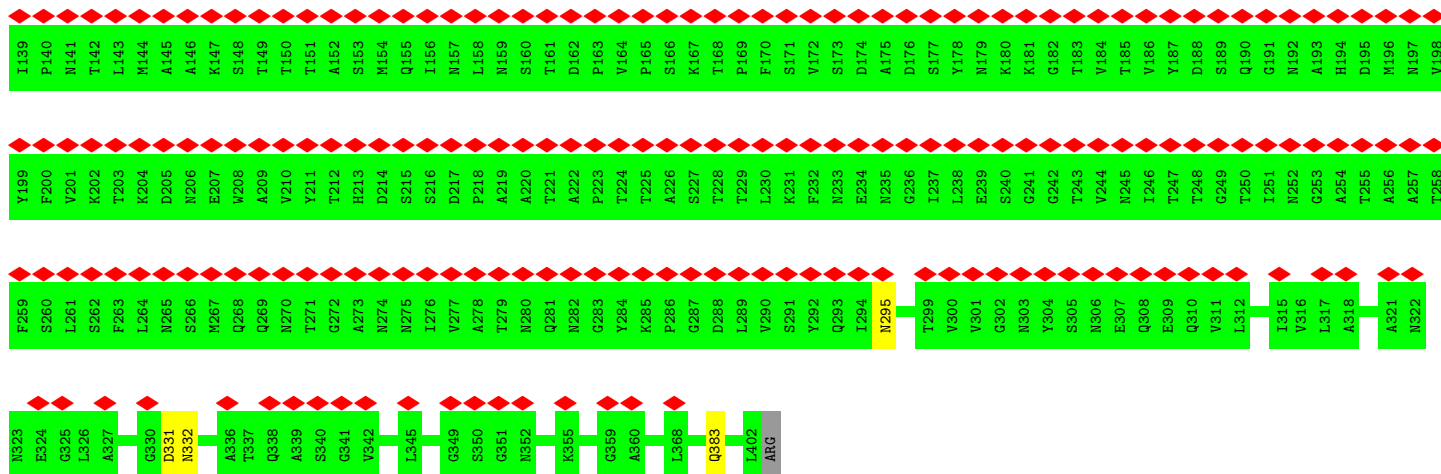


• Molecule 8: Flagellar hook protein FlgE

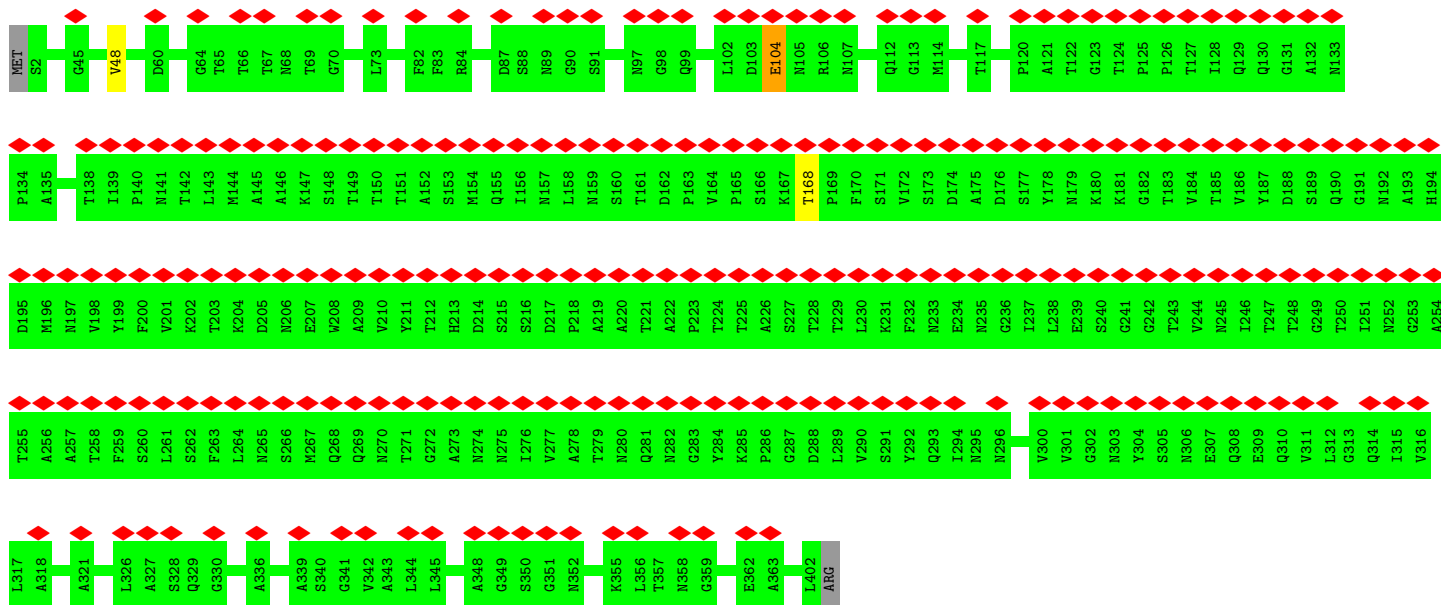


• Molecule 8: Flagellar hook protein FlgE

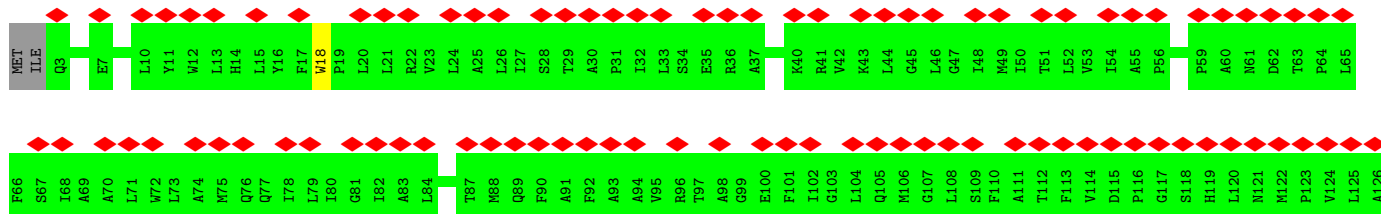


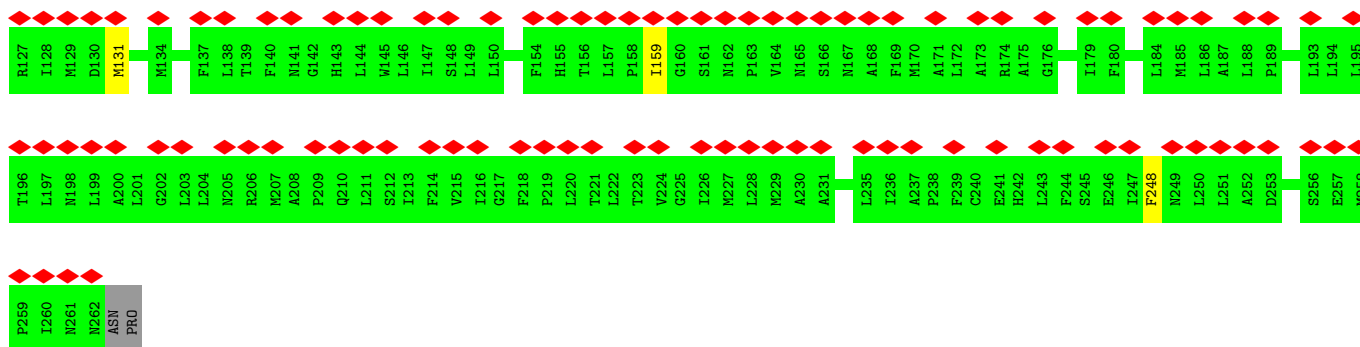


• Molecule 8: Flagellar hook protein FlgE

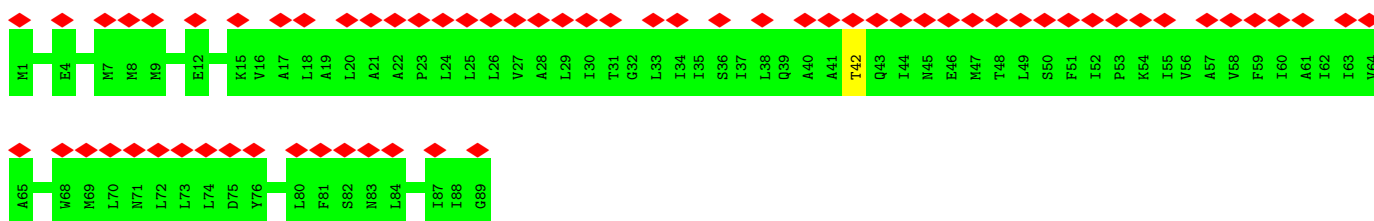
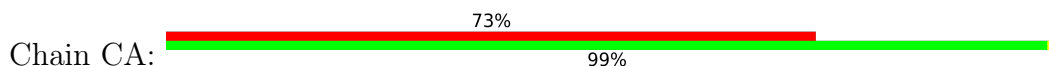


• Molecule 9: Flagellar biosynthetic protein FliR

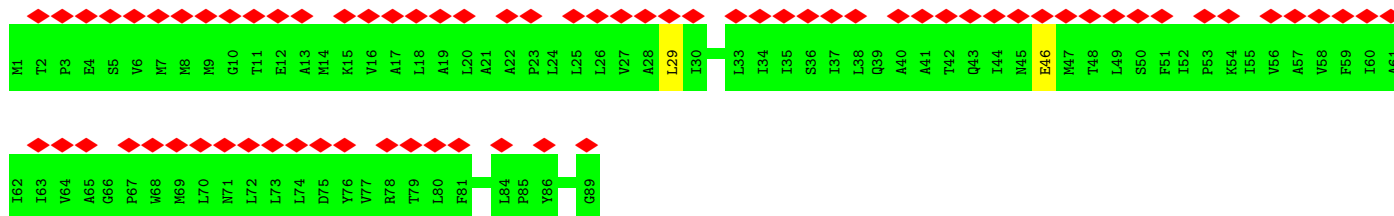
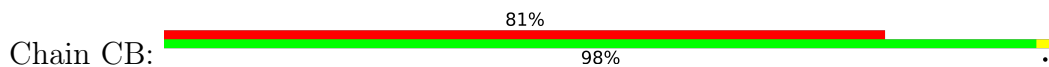




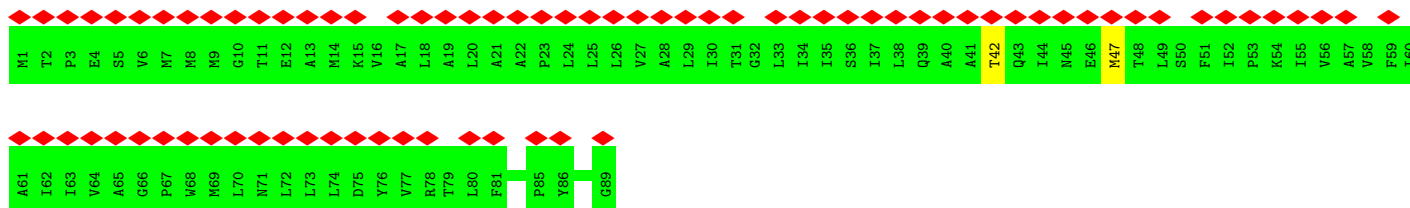
• Molecule 10: Flagellar biosynthetic protein FliQ



• Molecule 10: Flagellar biosynthetic protein FliQ



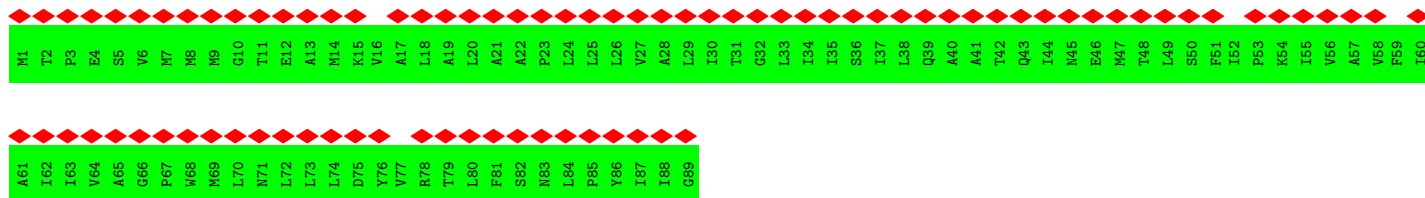
• Molecule 10: Flagellar biosynthetic protein FliQ



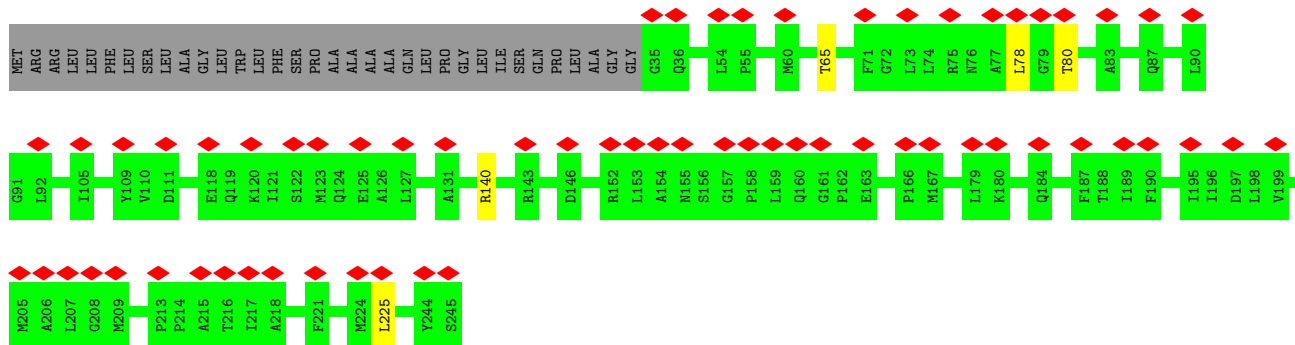
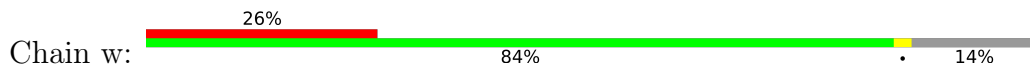
• Molecule 10: Flagellar biosynthetic protein FliQ



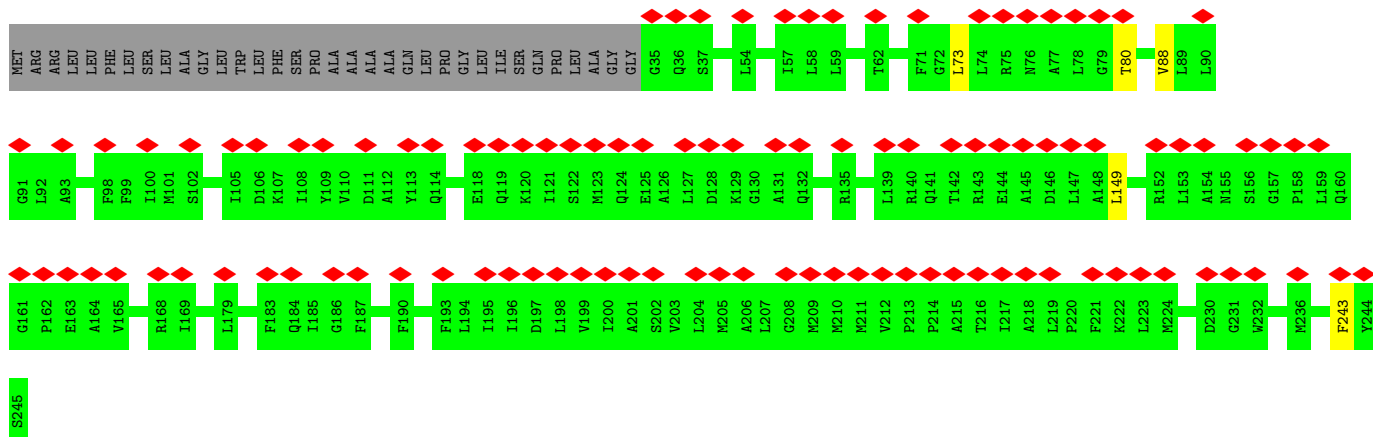
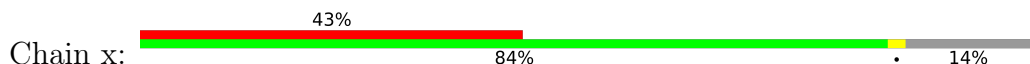




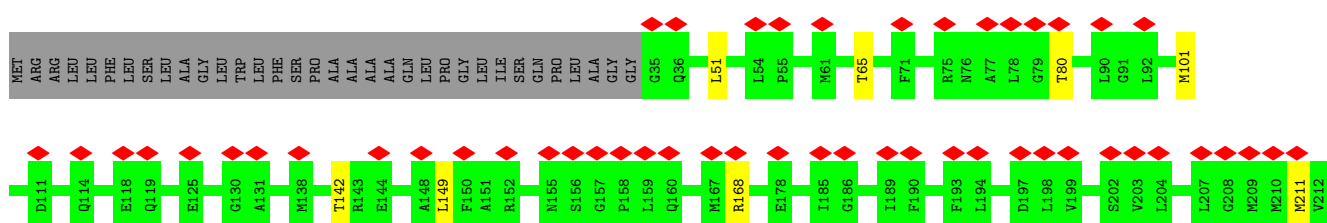
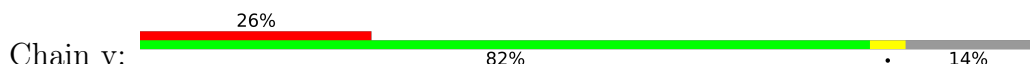
• Molecule 11: Flagellar biosynthetic protein FliP

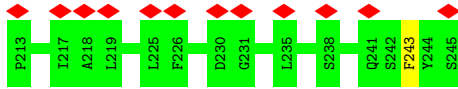


• Molecule 11: Flagellar biosynthetic protein FliP

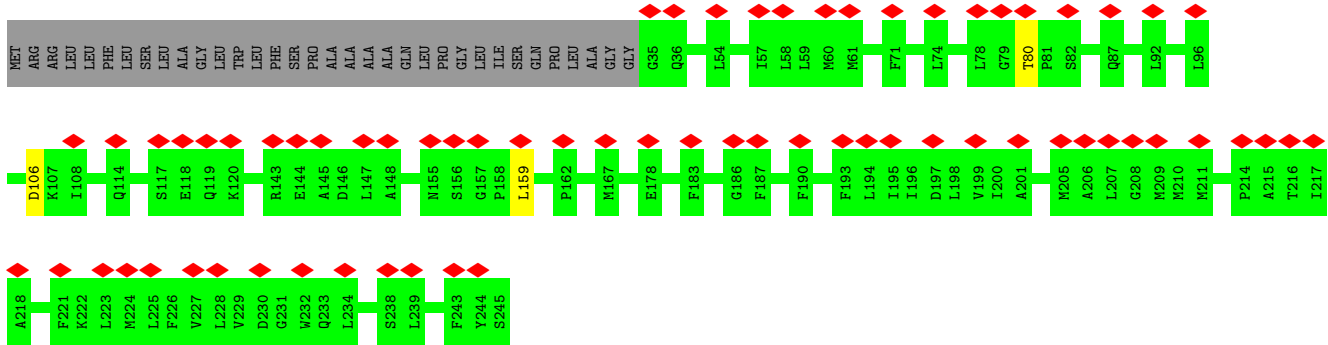
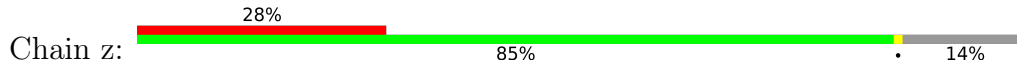


• Molecule 11: Flagellar biosynthetic protein FliP

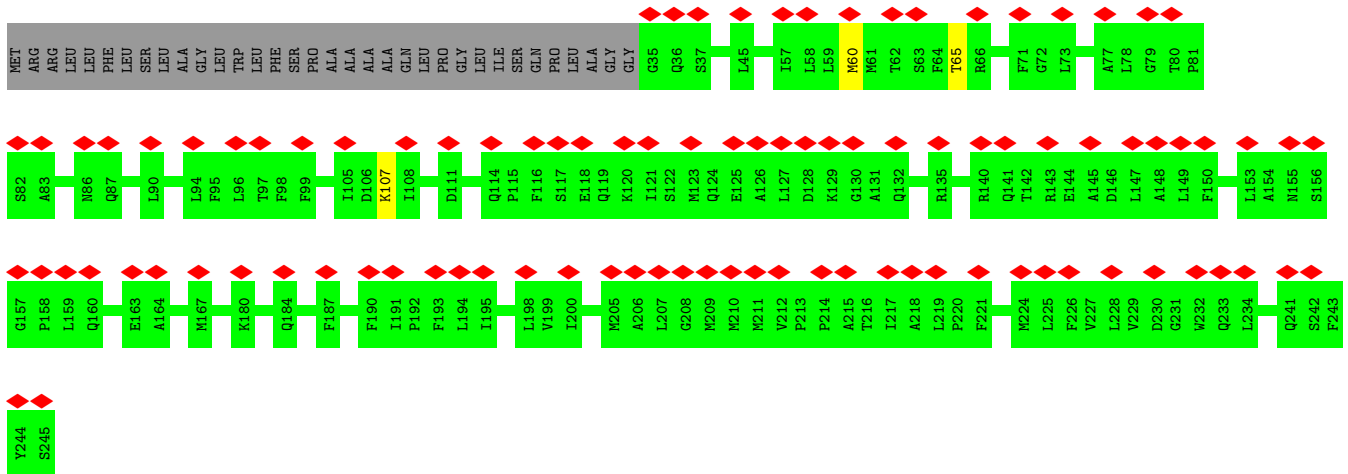
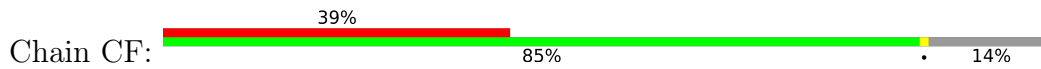




- Molecule 11: Flagellar biosynthetic protein FliP



- Molecule 11: Flagellar biosynthetic protein FliP



## 4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	52714	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 ( $e^-/\text{\AA}^2$ )	47	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	2.440	Depositor
Minimum map value	-1.067	Depositor
Average map value	-0.001	Depositor
Map value standard deviation	0.092	Depositor
Recommended contour level	0.6	Depositor
Map size (Å)	669.184, 669.184, 669.184	wwPDB
Map dimensions	512, 512, 512	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.307, 1.307, 1.307	Depositor

## 5 Model quality [i](#)

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

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
1	A	0.27	0/1973	0.48	0/2682
1	B	0.27	0/1973	0.48	0/2682
1	C	0.27	0/1973	0.48	0/2682
1	D	0.28	0/1973	0.48	0/2682
1	E	0.27	0/1973	0.49	0/2682
1	F	0.28	0/1973	0.48	0/2682
1	G	0.28	0/1973	0.49	0/2682
1	H	0.27	0/1973	0.48	0/2682
1	I	0.28	0/1973	0.49	0/2682
1	J	0.28	0/1973	0.48	0/2682
1	K	0.28	0/1973	0.50	0/2682
1	L	0.28	0/1965	0.47	0/2672
1	M	0.28	0/1973	0.48	0/2682
1	N	0.28	0/1909	0.47	0/2593
1	O	0.28	0/1917	0.48	0/2605
1	P	0.28	0/1884	0.48	0/2559
1	Q	0.28	0/1880	0.48	0/2554
1	R	0.28	0/1898	0.48	0/2578
1	S	0.29	0/1880	0.50	0/2554
1	T	0.28	0/1925	0.47	0/2617
1	U	0.28	0/1965	0.48	0/2672
1	V	0.28	0/1965	0.47	0/2672
1	W	0.29	0/1965	0.49	0/2672
1	X	0.27	0/1965	0.49	1/2672 (0.0%)
2	a	0.27	0/1836	0.50	0/2502
2	b	0.26	0/1828	0.49	0/2492
2	c	0.27	0/1836	0.49	0/2502
2	d	0.27	0/1836	0.50	0/2502
2	e	0.26	0/1836	0.50	0/2502
4	5	0.28	0/145	0.42	0/203
4	6	0.26	0/145	0.43	0/203
4	7	0.40	0/145	0.55	0/203
4	8	0.27	0/145	0.41	0/203
4	9	0.26	0/145	0.44	0/203

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
5	f	0.27	0/946	0.44	0/1285
5	g	0.27	0/959	0.45	0/1302
5	h	0.28	0/945	0.45	0/1283
5	i	0.28	0/954	0.46	0/1295
5	j	0.28	0/941	0.46	0/1278
5	p	0.27	0/950	0.43	0/1290
6	k	0.26	0/859	0.48	0/1156
6	l	0.26	0/840	0.48	0/1131
6	m	0.26	0/840	0.47	0/1131
6	n	0.26	0/851	0.49	0/1145
6	o	0.25	0/863	0.50	0/1161
7	q	0.26	0/540	0.44	0/723
7	r	0.26	0/529	0.46	0/709
7	s	0.26	0/547	0.44	0/733
7	t	0.25	0/547	0.44	0/733
7	u	0.25	0/547	0.46	0/733
7	v	0.24	0/289	0.39	0/385
8	DA	0.26	0/2991	0.45	0/4076
8	DB	0.26	0/2991	0.45	0/4076
8	DC	0.26	0/2991	0.45	0/4076
8	DD	0.26	0/2991	0.45	0/4076
8	DE	0.26	0/2991	0.45	0/4076
8	DF	0.25	0/2991	0.45	0/4076
8	DG	0.25	0/2991	0.45	0/4076
8	DH	0.26	0/2991	0.45	0/4076
8	DI	0.26	0/2991	0.46	0/4076
8	DJ	0.28	0/2991	0.48	1/4076 (0.0%)
8	DK	0.26	0/2991	0.47	1/4076 (0.0%)
9	CE	0.25	0/2052	0.44	0/2803
10	CA	0.24	0/681	0.45	1/930 (0.1%)
10	CB	0.25	0/681	0.44	0/930
10	CC	0.24	0/681	0.46	1/930 (0.1%)
10	CD	0.24	0/681	0.46	0/930
11	CF	0.25	0/1675	0.44	0/2280
11	w	0.27	0/1675	0.44	0/2280
11	x	0.26	0/1675	0.44	0/2280
11	y	0.26	0/1675	0.46	0/2280
11	z	0.26	0/1675	0.44	0/2280
All	All	0.27	0/115690	0.47	5/157351 (0.0%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	204	SER	C-N-CA	5.85	136.33	121.70
8	DJ	383	GLN	CB-CA-C	-5.64	99.12	110.40
8	DK	104	GLU	CB-CA-C	5.37	121.14	110.40
10	CC	42	THR	C-N-CA	5.04	134.30	121.70
10	CA	42	THR	C-N-CA	5.03	134.28	121.70

There are no chirality outliers.

There are no planarity outliers.

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

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

## 5.3 Torsion angles [i](#)

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

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	258/260 (99%)	249 (96%)	9 (4%)	0	100	100
1	B	258/260 (99%)	247 (96%)	11 (4%)	0	100	100
1	C	258/260 (99%)	247 (96%)	11 (4%)	0	100	100
1	D	258/260 (99%)	248 (96%)	10 (4%)	0	100	100
1	E	258/260 (99%)	246 (95%)	12 (5%)	0	100	100
1	F	258/260 (99%)	248 (96%)	10 (4%)	0	100	100
1	G	258/260 (99%)	247 (96%)	11 (4%)	0	100	100
1	H	258/260 (99%)	245 (95%)	13 (5%)	0	100	100
1	I	258/260 (99%)	248 (96%)	10 (4%)	0	100	100
1	J	258/260 (99%)	251 (97%)	7 (3%)	0	100	100
1	K	258/260 (99%)	248 (96%)	10 (4%)	0	100	100
1	L	257/260 (99%)	250 (97%)	7 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	M	258/260 (99%)	247 (96%)	11 (4%)	0	100	100
1	N	247/260 (95%)	240 (97%)	7 (3%)	0	100	100
1	O	248/260 (95%)	240 (97%)	8 (3%)	0	100	100
1	P	244/260 (94%)	236 (97%)	8 (3%)	0	100	100
1	Q	243/260 (94%)	237 (98%)	6 (2%)	0	100	100
1	R	246/260 (95%)	238 (97%)	8 (3%)	0	100	100
1	S	243/260 (94%)	235 (97%)	8 (3%)	0	100	100
1	T	249/260 (96%)	241 (97%)	8 (3%)	0	100	100
1	U	257/260 (99%)	243 (95%)	14 (5%)	0	100	100
1	V	257/260 (99%)	249 (97%)	8 (3%)	0	100	100
1	W	257/260 (99%)	242 (94%)	15 (6%)	0	100	100
1	X	257/260 (99%)	244 (95%)	13 (5%)	0	100	100
2	a	247/251 (98%)	235 (95%)	12 (5%)	0	100	100
2	b	246/251 (98%)	236 (96%)	10 (4%)	0	100	100
2	c	247/251 (98%)	237 (96%)	10 (4%)	0	100	100
2	d	247/251 (98%)	232 (94%)	15 (6%)	0	100	100
2	e	247/251 (98%)	231 (94%)	16 (6%)	0	100	100
4	5	19/21 (90%)	19 (100%)	0	0	100	100
4	6	19/21 (90%)	19 (100%)	0	0	100	100
4	7	19/21 (90%)	19 (100%)	0	0	100	100
4	8	19/21 (90%)	18 (95%)	1 (5%)	0	100	100
4	9	19/21 (90%)	19 (100%)	0	0	100	100
5	f	124/134 (92%)	121 (98%)	3 (2%)	0	100	100
5	g	126/134 (94%)	121 (96%)	5 (4%)	0	100	100
5	h	124/134 (92%)	120 (97%)	4 (3%)	0	100	100
5	i	125/134 (93%)	121 (97%)	4 (3%)	0	100	100
5	j	123/134 (92%)	119 (97%)	4 (3%)	0	100	100
5	p	125/134 (93%)	120 (96%)	5 (4%)	0	100	100
6	k	104/138 (75%)	99 (95%)	5 (5%)	0	100	100
6	l	102/138 (74%)	101 (99%)	1 (1%)	0	100	100
6	m	102/138 (74%)	99 (97%)	3 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
6	n	103/138 (75%)	99 (96%)	4 (4%)	0	100	100
6	o	105/138 (76%)	103 (98%)	2 (2%)	0	100	100
7	q	69/104 (66%)	68 (99%)	1 (1%)	0	100	100
7	r	68/104 (65%)	64 (94%)	4 (6%)	0	100	100
7	s	70/104 (67%)	70 (100%)	0	0	100	100
7	t	70/104 (67%)	69 (99%)	1 (1%)	0	100	100
7	u	70/104 (67%)	70 (100%)	0	0	100	100
7	v	36/104 (35%)	36 (100%)	0	0	100	100
8	DA	399/403 (99%)	377 (94%)	22 (6%)	0	100	100
8	DB	399/403 (99%)	377 (94%)	22 (6%)	0	100	100
8	DC	399/403 (99%)	375 (94%)	24 (6%)	0	100	100
8	DD	399/403 (99%)	378 (95%)	21 (5%)	0	100	100
8	DE	399/403 (99%)	380 (95%)	19 (5%)	0	100	100
8	DF	399/403 (99%)	376 (94%)	23 (6%)	0	100	100
8	DG	399/403 (99%)	382 (96%)	17 (4%)	0	100	100
8	DH	399/403 (99%)	378 (95%)	19 (5%)	2 (0%)	29	66
8	DI	399/403 (99%)	381 (96%)	18 (4%)	0	100	100
8	DJ	399/403 (99%)	375 (94%)	23 (6%)	1 (0%)	41	74
8	DK	399/403 (99%)	386 (97%)	13 (3%)	0	100	100
9	CE	258/264 (98%)	239 (93%)	18 (7%)	1 (0%)	34	69
10	CA	87/89 (98%)	77 (88%)	10 (12%)	0	100	100
10	CB	87/89 (98%)	78 (90%)	9 (10%)	0	100	100
10	CC	87/89 (98%)	77 (88%)	10 (12%)	0	100	100
10	CD	87/89 (98%)	77 (88%)	10 (12%)	0	100	100
11	CF	209/245 (85%)	203 (97%)	6 (3%)	0	100	100
11	w	209/245 (85%)	203 (97%)	6 (3%)	0	100	100
11	x	209/245 (85%)	197 (94%)	12 (6%)	0	100	100
11	y	209/245 (85%)	201 (96%)	8 (4%)	0	100	100
11	z	209/245 (85%)	204 (98%)	5 (2%)	0	100	100
All	All	15116/15996 (94%)	14452 (96%)	660 (4%)	4 (0%)	100	100

All (4) Ramachandran outliers are listed below:



Mol	Chain	Res	Type
8	DJ	332	ASN
8	DH	270	ASN
9	CE	159	ILE
8	DH	71	ARG

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

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	215/215 (100%)	211 (98%)	4 (2%)	57	76
1	B	215/215 (100%)	212 (99%)	3 (1%)	67	82
1	C	215/215 (100%)	213 (99%)	2 (1%)	78	88
1	D	215/215 (100%)	215 (100%)	0	100	100
1	E	215/215 (100%)	210 (98%)	5 (2%)	50	71
1	F	215/215 (100%)	214 (100%)	1 (0%)	88	94
1	G	215/215 (100%)	209 (97%)	6 (3%)	43	67
1	H	215/215 (100%)	213 (99%)	2 (1%)	78	88
1	I	215/215 (100%)	211 (98%)	4 (2%)	57	76
1	J	215/215 (100%)	213 (99%)	2 (1%)	78	88
1	K	215/215 (100%)	211 (98%)	4 (2%)	57	76
1	L	214/215 (100%)	211 (99%)	3 (1%)	67	82
1	M	215/215 (100%)	212 (99%)	3 (1%)	67	82
1	N	208/215 (97%)	207 (100%)	1 (0%)	88	94
1	O	209/215 (97%)	208 (100%)	1 (0%)	88	94
1	P	204/215 (95%)	202 (99%)	2 (1%)	76	86
1	Q	204/215 (95%)	200 (98%)	4 (2%)	55	74
1	R	206/215 (96%)	201 (98%)	5 (2%)	49	70
1	S	204/215 (95%)	202 (99%)	2 (1%)	76	86
1	T	210/215 (98%)	209 (100%)	1 (0%)	88	94
1	U	214/215 (100%)	212 (99%)	2 (1%)	78	88

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	V	214/215 (100%)	211 (99%)	3 (1%)	67	82
1	W	214/215 (100%)	211 (99%)	3 (1%)	67	82
1	X	214/215 (100%)	210 (98%)	4 (2%)	57	76
2	a	191/193 (99%)	185 (97%)	6 (3%)	40	65
2	b	190/193 (98%)	186 (98%)	4 (2%)	53	73
2	c	191/193 (99%)	185 (97%)	6 (3%)	40	65
2	d	191/193 (99%)	187 (98%)	4 (2%)	53	73
2	e	191/193 (99%)	188 (98%)	3 (2%)	62	79
4	5	15/15 (100%)	15 (100%)	0	100	100
4	6	15/15 (100%)	15 (100%)	0	100	100
4	7	15/15 (100%)	15 (100%)	0	100	100
4	8	15/15 (100%)	15 (100%)	0	100	100
4	9	15/15 (100%)	15 (100%)	0	100	100
5	f	101/105 (96%)	98 (97%)	3 (3%)	41	66
5	g	102/105 (97%)	99 (97%)	3 (3%)	42	66
5	h	101/105 (96%)	100 (99%)	1 (1%)	76	86
5	i	102/105 (97%)	100 (98%)	2 (2%)	55	74
5	j	101/105 (96%)	100 (99%)	1 (1%)	76	86
5	p	101/105 (96%)	101 (100%)	0	100	100
6	k	91/113 (80%)	87 (96%)	4 (4%)	28	58
6	l	89/113 (79%)	86 (97%)	3 (3%)	37	64
6	m	89/113 (79%)	88 (99%)	1 (1%)	73	85
6	n	90/113 (80%)	87 (97%)	3 (3%)	38	64
6	o	91/113 (80%)	87 (96%)	4 (4%)	28	58
7	q	55/79 (70%)	54 (98%)	1 (2%)	59	77
7	r	54/79 (68%)	54 (100%)	0	100	100
7	s	56/79 (71%)	56 (100%)	0	100	100
7	t	56/79 (71%)	56 (100%)	0	100	100
7	u	56/79 (71%)	56 (100%)	0	100	100
7	v	32/79 (40%)	31 (97%)	1 (3%)	40	65
8	DA	321/323 (99%)	315 (98%)	6 (2%)	57	76

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
8	DB	321/323 (99%)	317 (99%)	4 (1%)	71	84
8	DC	321/323 (99%)	320 (100%)	1 (0%)	92	96
8	DD	321/323 (99%)	318 (99%)	3 (1%)	78	88
8	DE	321/323 (99%)	320 (100%)	1 (0%)	92	96
8	DF	321/323 (99%)	318 (99%)	3 (1%)	78	88
8	DG	321/323 (99%)	320 (100%)	1 (0%)	92	96
8	DH	321/323 (99%)	316 (98%)	5 (2%)	62	79
8	DI	321/323 (99%)	316 (98%)	5 (2%)	62	79
8	DJ	321/323 (99%)	318 (99%)	3 (1%)	78	88
8	DK	321/323 (99%)	318 (99%)	3 (1%)	78	88
9	CE	217/221 (98%)	214 (99%)	3 (1%)	67	82
10	CA	74/74 (100%)	74 (100%)	0	100	100
10	CB	74/74 (100%)	72 (97%)	2 (3%)	44	68
10	CC	74/74 (100%)	73 (99%)	1 (1%)	67	82
10	CD	74/74 (100%)	74 (100%)	0	100	100
11	CF	180/204 (88%)	177 (98%)	3 (2%)	60	79
11	w	180/204 (88%)	175 (97%)	5 (3%)	43	67
11	x	180/204 (88%)	175 (97%)	5 (3%)	43	67
11	y	180/204 (88%)	171 (95%)	9 (5%)	24	55
11	z	180/204 (88%)	177 (98%)	3 (2%)	60	79
All	All	12435/12959 (96%)	12252 (98%)	183 (2%)	66	81

All (183) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	153	ARG
1	A	159	VAL
1	A	164	GLN
1	A	256	LYS
1	B	19	MET
1	B	77	THR
1	B	179	MET
1	C	141	THR
1	C	256	LYS
1	E	28	ASN

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	E	42	GLU
1	E	97	ILE
1	E	164	GLN
1	E	259	GLN
1	F	177	THR
1	G	117	ASP
1	G	141	THR
1	G	168	VAL
1	G	195	THR
1	G	241	GLU
1	G	255	GLN
1	H	105	VAL
1	H	226	VAL
1	I	42	GLU
1	I	58	GLU
1	I	150	THR
1	I	256	LYS
1	J	70	THR
1	J	97	ILE
1	K	42	GLU
1	K	59	GLN
1	K	97	ILE
1	K	202	ASN
1	L	42	GLU
1	L	117	ASP
1	L	175	LEU
1	M	70	THR
1	M	72	VAL
1	M	256	LYS
1	N	42	GLU
1	O	179	MET
1	P	50	ARG
1	P	254	LEU
1	Q	42	GLU
1	Q	141	THR
1	Q	149	ILE
1	Q	179	MET
1	R	43	ASP
1	R	141	THR
1	R	149	ILE
1	R	150	THR
1	R	181	ASP

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	S	109	ASP
1	S	141	THR
1	T	141	THR
1	U	150	THR
1	U	257	LEU
2	a	51	LEU
2	a	71	LEU
2	a	78	LEU
2	a	150	ASP
2	a	153	ASN
2	a	247	LEU
2	b	49	LEU
2	b	71	LEU
2	b	182	THR
2	b	247	LEU
2	c	7	THR
2	c	44	VAL
2	c	71	LEU
2	c	78	LEU
2	c	214	VAL
2	c	238	ASP
2	d	58	THR
2	d	71	LEU
2	d	78	LEU
2	d	116	GLN
2	e	49	LEU
2	e	78	LEU
2	e	114	THR
5	f	119	LEU
5	f	125	MET
5	f	132	LEU
5	g	4	LEU
5	g	48	VAL
5	g	116	ILE
5	h	119	LEU
5	j	79	GLU
6	l	86	LEU
6	l	132	ASN
6	l	133	VAL
6	m	4	ARG
6	o	18	LEU
6	o	48	GLU

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
6	o	83	VAL
6	o	127	LEU
5	i	97	ASP
5	i	101	GLU
6	k	29	ASN
6	k	48	GLU
6	k	51	LYS
6	k	111	ASP
6	n	29	ASN
6	n	100	VAL
6	n	132	ASN
1	V	28	ASN
1	V	45	LEU
1	V	90	ASN
1	W	168	VAL
1	W	257	LEU
1	W	260	LEU
1	X	142	ILE
1	X	179	MET
1	X	185	GLU
1	X	195	THR
7	q	57	GLU
7	v	75	MET
8	DA	17	LEU
8	DA	48	VAL
8	DA	157	ASN
8	DA	212	THR
8	DA	322	ASN
8	DA	352	ASN
8	DB	26	ASN
8	DB	161	THR
8	DB	183	THR
8	DB	324	GLU
8	DC	168	THR
8	DD	26	ASN
8	DD	80	ASN
8	DD	295	ASN
8	DE	295	ASN
8	DF	250	THR
8	DF	332	ASN
8	DF	344	LEU
8	DG	395	ILE

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
8	DH	97	ASN
8	DH	161	THR
8	DH	279	THR
8	DH	322	ASN
8	DH	355	LYS
8	DI	17	LEU
8	DI	60	ASP
8	DI	119	TYR
8	DI	168	THR
8	DI	185	THR
8	DJ	48	VAL
8	DJ	295	ASN
8	DJ	331	ASP
8	DK	48	VAL
8	DK	104	GLU
8	DK	168	THR
9	CE	18	TRP
9	CE	131	MET
9	CE	248	PHE
10	CB	29	LEU
10	CB	46	GLU
10	CC	47	MET
11	w	65	THR
11	w	78	LEU
11	w	80	THR
11	w	140	ARG
11	w	225	LEU
11	x	73	LEU
11	x	80	THR
11	x	88	VAL
11	x	149	LEU
11	x	243	PHE
11	y	51	LEU
11	y	65	THR
11	y	80	THR
11	y	101	MET
11	y	142	THR
11	y	149	LEU
11	y	168	ARG
11	y	211	MET
11	y	243	PHE
11	z	80	THR

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Mol	Chain	Res	Type
11	z	106	ASP
11	z	159	LEU
11	CF	60	MET
11	CF	65	THR
11	CF	107	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (47) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	255	GLN
1	C	259	GLN
1	E	164	GLN
1	F	145	ASN
1	F	161	GLN
1	F	259	GLN
1	J	255	GLN
1	J	259	GLN
1	K	25	ASN
1	K	59	GLN
1	T	135	GLN
1	T	172	GLN
5	g	82	ASN
5	h	30	ASN
6	m	29	ASN
6	o	29	ASN
5	i	120	ASN
1	W	169	GLN
7	v	76	GLN
8	DA	97	ASN
8	DB	107	ASN
8	DB	129	GLN
8	DD	22	ASN
8	DD	26	ASN
8	DD	107	ASN
8	DE	107	ASN
8	DF	22	ASN
8	DF	26	ASN
8	DF	365	ASN
8	DG	129	GLN
8	DH	80	ASN
8	DH	197	ASN
8	DH	213	HIS

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Mol	Chain	Res	Type
8	DH	310	GLN
8	DH	319	ASN
8	DH	365	ASN
8	DI	129	GLN
8	DJ	310	GLN
8	DJ	352	ASN
8	DK	97	ASN
8	DK	105	ASN
8	DK	332	ASN
8	DK	383	GLN
10	CB	39	GLN
10	CB	45	ASN
11	y	184	GLN
11	z	87	GLN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

### 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

### 5.6 Ligand geometry [i](#)

There are no ligands in this entry.

### 5.7 Other polymers [i](#)

There are no such residues in this entry.

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

There are no chain breaks in this entry.

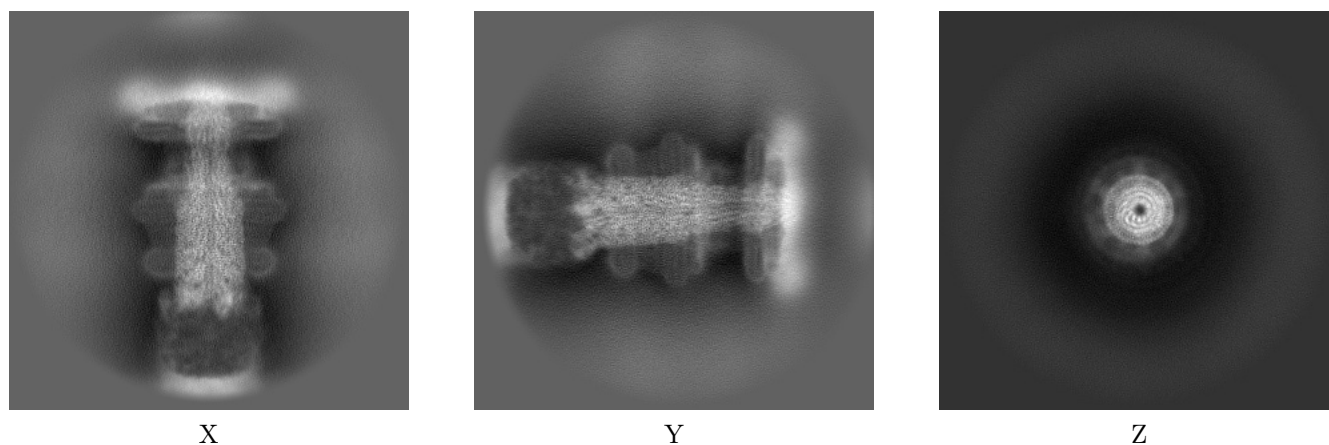
## 6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-31006. 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](#)

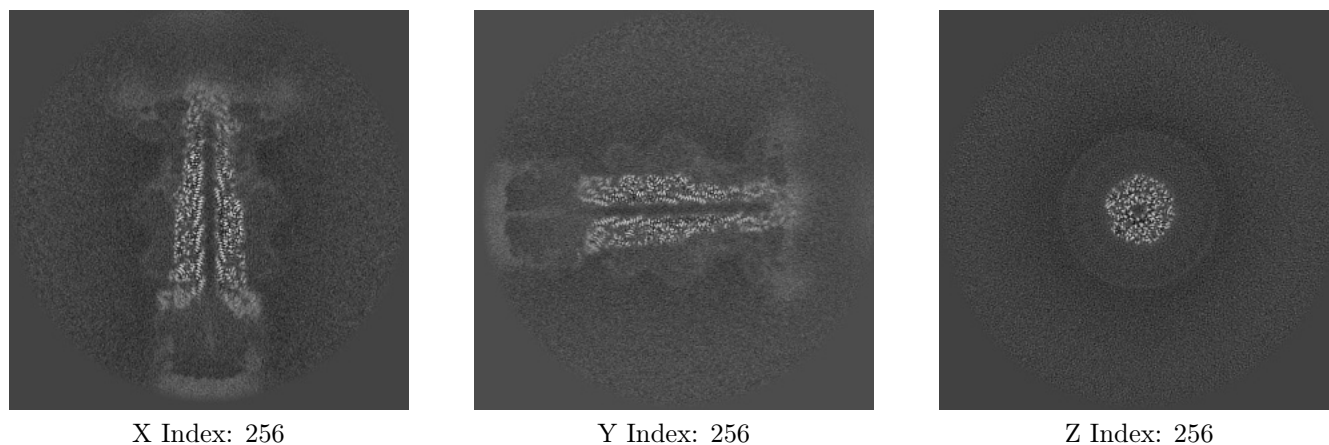
#### 6.1.1 Primary map



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

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

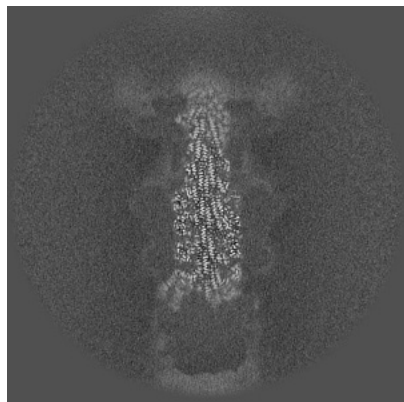
#### 6.2.1 Primary map



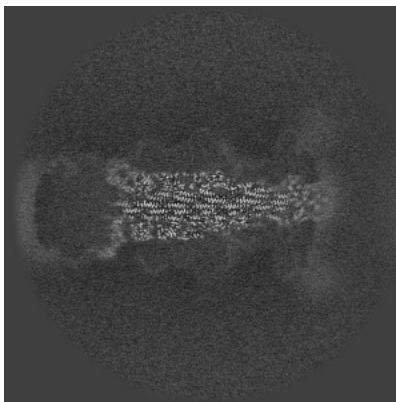
The images above show central slices of the map in three orthogonal directions.

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

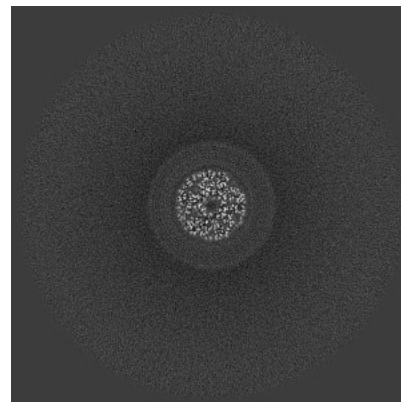
### 6.3.1 Primary map



X Index: 269



Y Index: 243

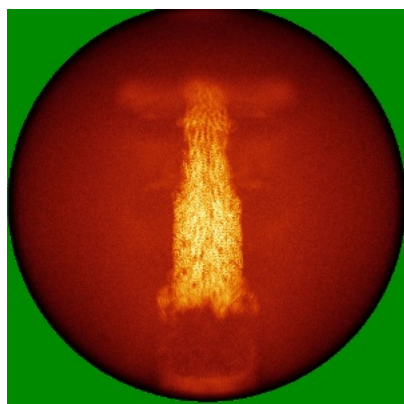


Z Index: 228

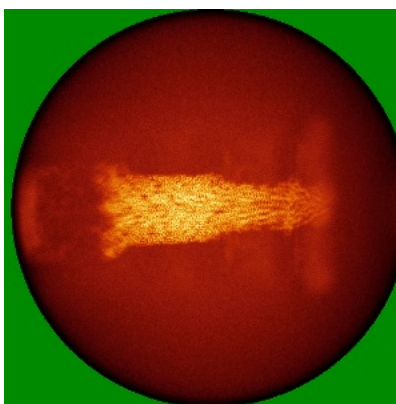
The images above show the largest variance slices of the map in three orthogonal directions.

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

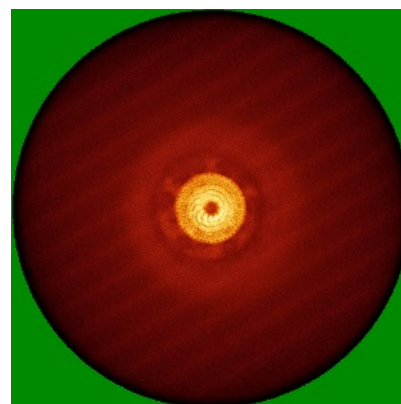
### 6.4.1 Primary map



X



Y

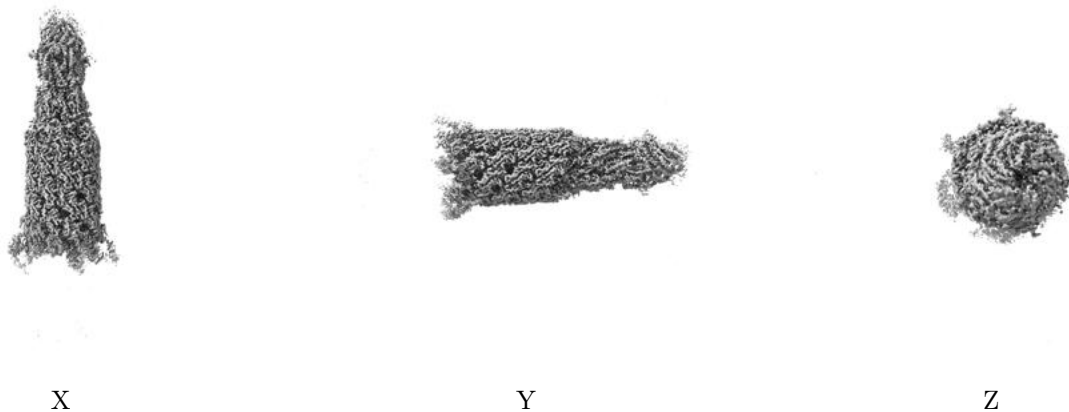


Z

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

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

### 6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.6. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

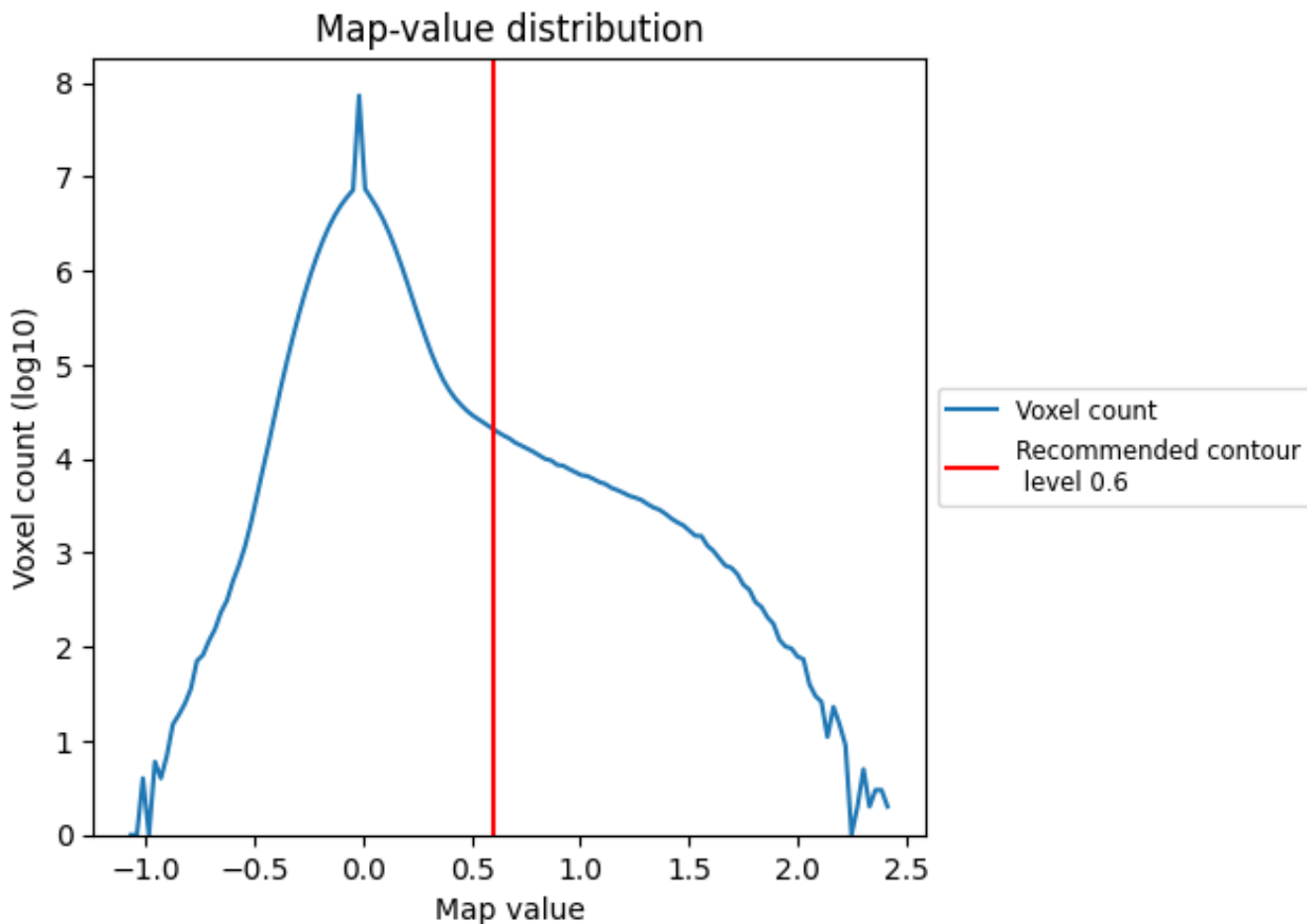
## 6.6 Mask visualisation [i](#)

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

## 7 Map analysis [i](#)

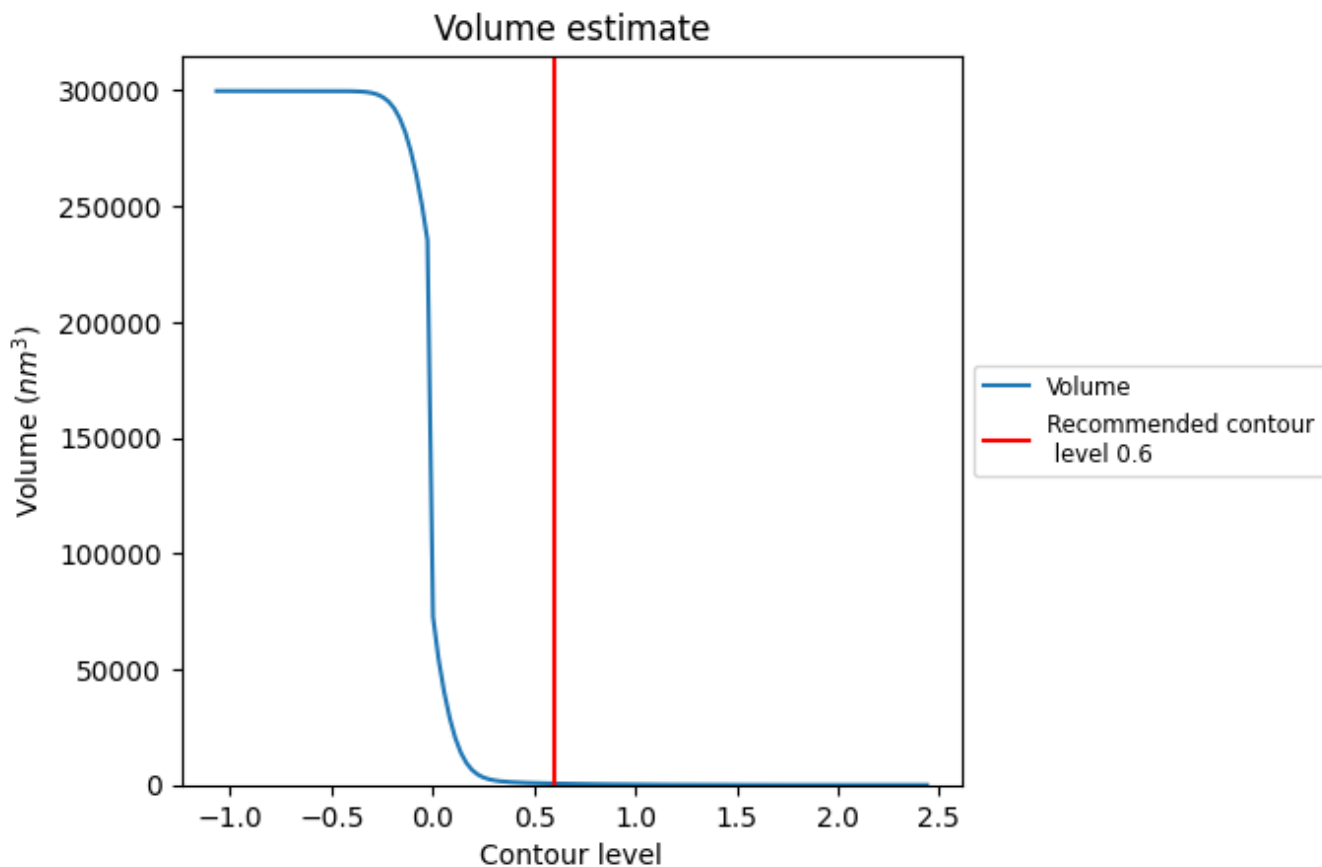
This section contains the results of statistical analysis of the map.

### 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.

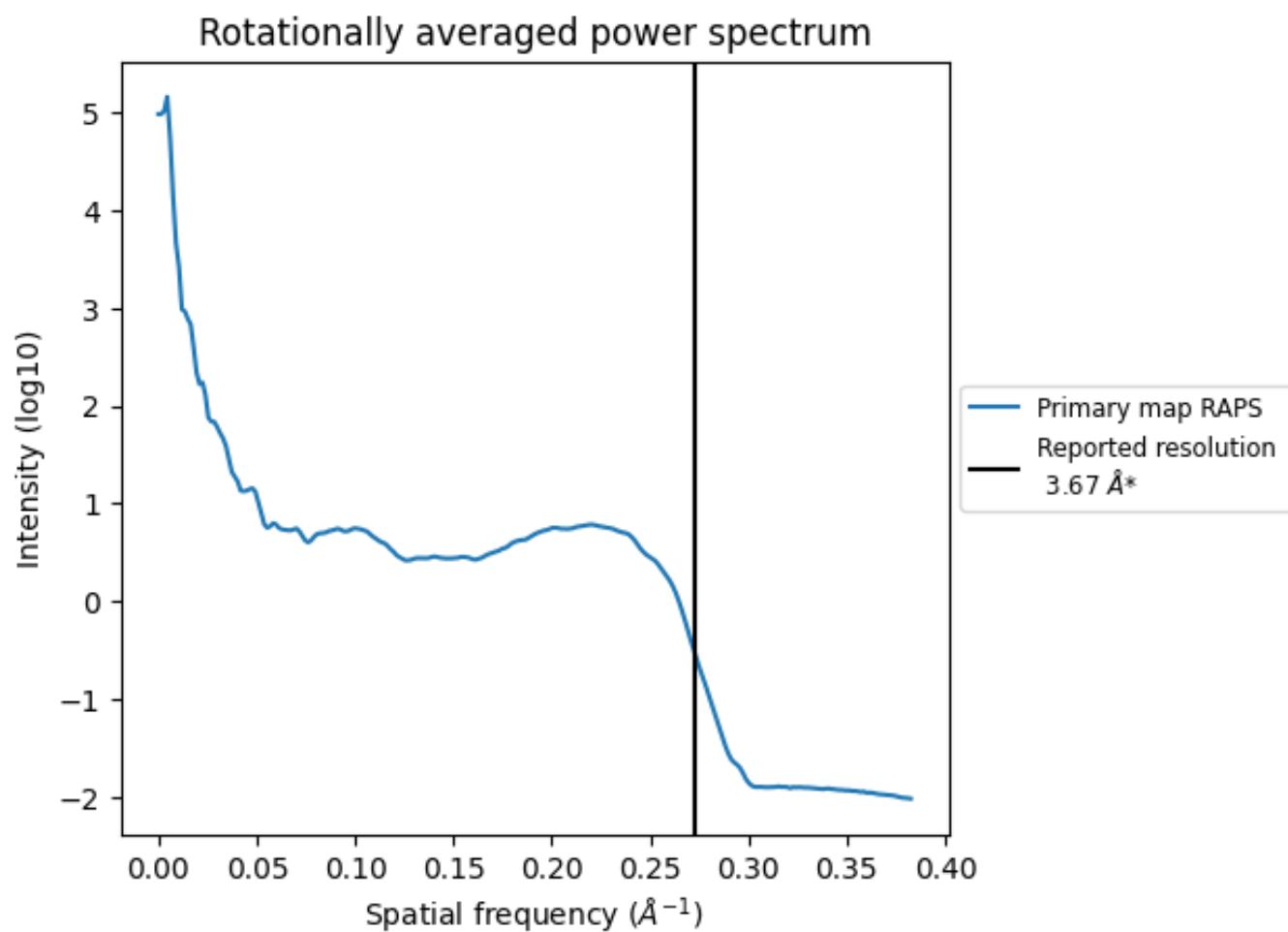
## 7.2 Volume estimate [\(i\)](#)



The volume at the recommended contour level is  $603 \text{ nm}^3$ ; this corresponds to an approximate mass of 545 kDa.

The volume estimate graph shows how the enclosed volume varies with the contour level. The recommended contour level is shown as a vertical line and the intersection between the line and the curve gives the volume of the enclosed surface at the given level.

### 7.3 Rotationally averaged power spectrum [i](#)



\*Reported resolution corresponds to spatial frequency of  $0.272 \text{\AA}^{-1}$

## 8 Fourier-Shell correlation

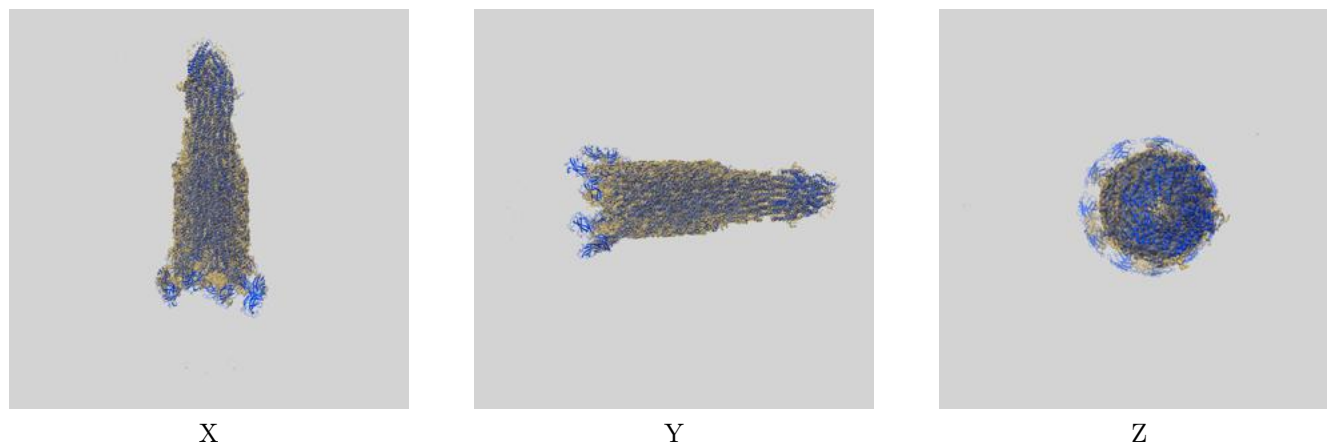
This section was not generated. No FSC curve or half-maps provided.



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

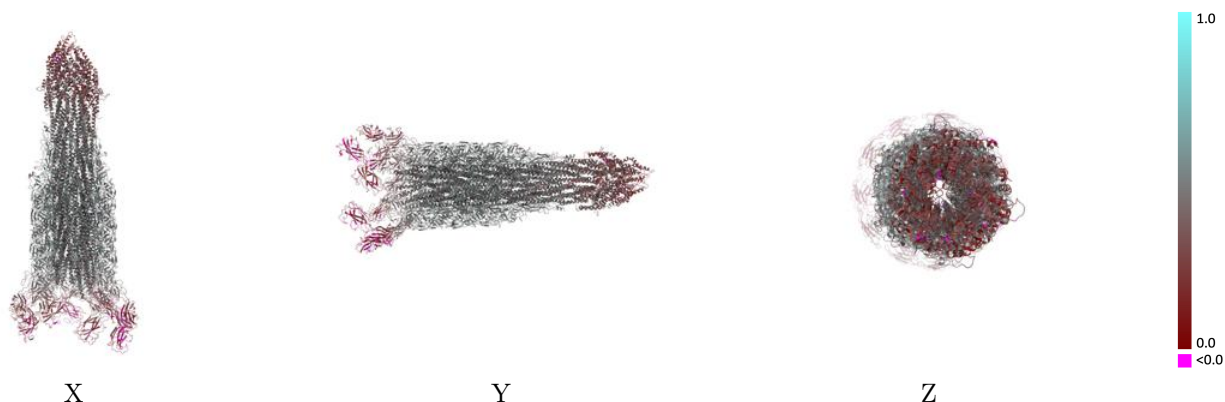
This section contains information regarding the fit between EMDB map EMD-31006 and PDB model 7E80. Per-residue inclusion information can be found in section 3 on page 11.

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



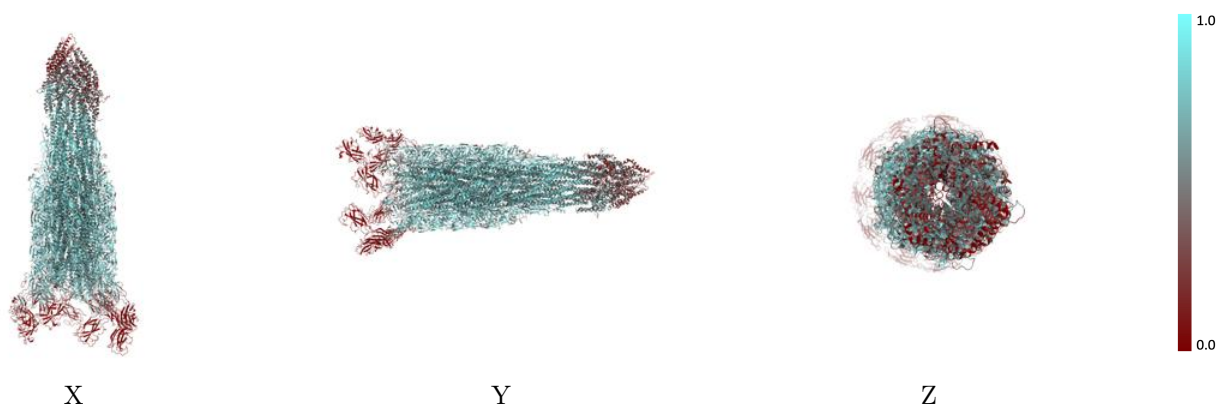
The images above show the 3D surface view of the map at the recommended contour level 0.6 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



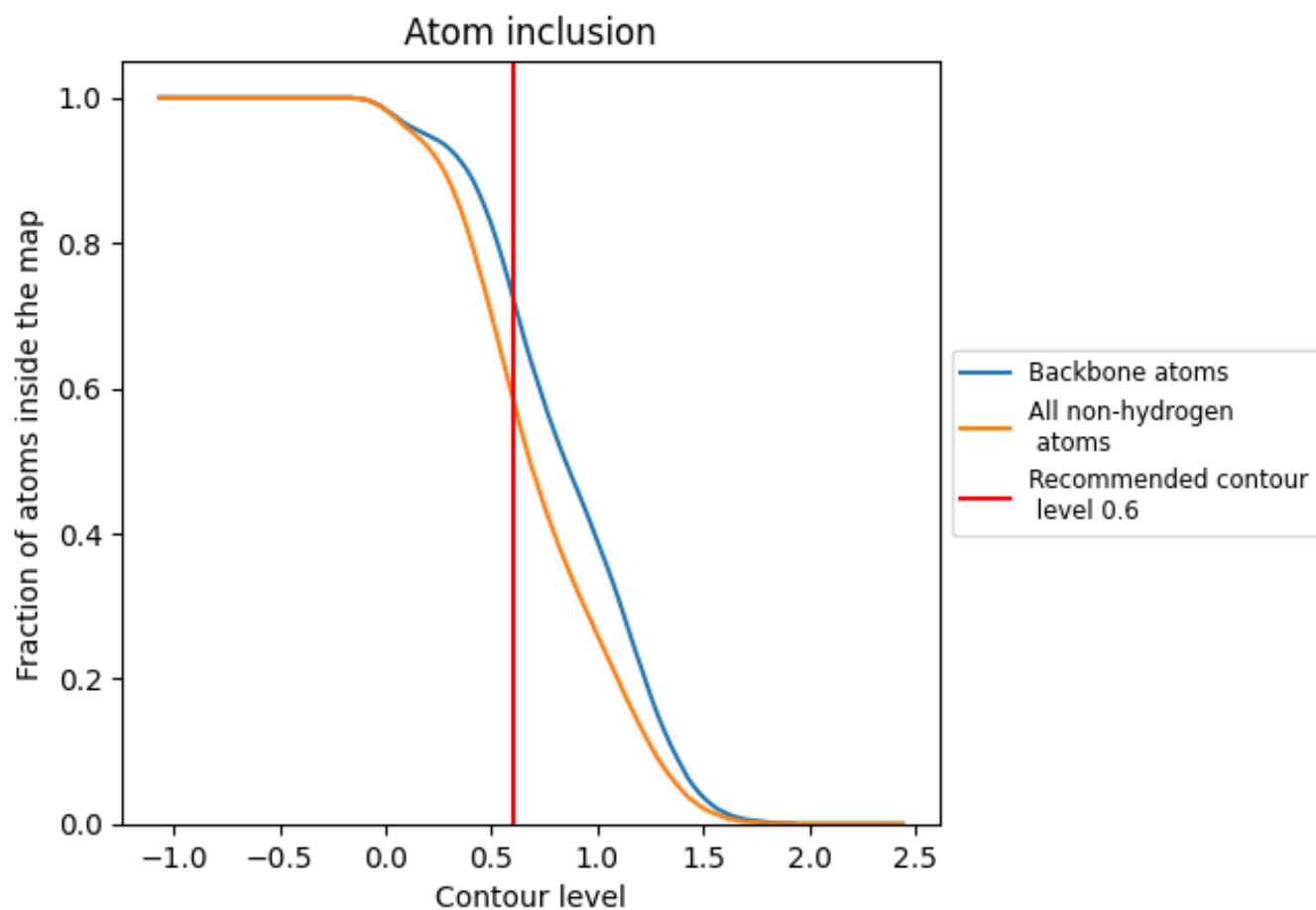
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.6).







































































## 9.4 Atom inclusion [i](#)



At the recommended contour level, 73% of all backbone atoms, 59% 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 (0.6) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.5860	 0.4230
0	 0.7330	 0.4950
1	 0.7330	 0.4700
2	 0.8670	 0.5240
3	 0.8130	 0.4680
4	 0.6930	 0.4640
5	 0.7070	 0.4870
6	 0.4930	 0.4500
7	 0.6640	 0.4700
8	 0.6360	 0.4830
9	 0.5290	 0.4590
A	 0.6970	 0.4820
B	 0.6480	 0.4640
C	 0.7030	 0.4790
CA	 0.2990	 0.2650
CB	 0.2360	 0.2680
CC	 0.2100	 0.2600
CD	 0.1080	 0.2520
CE	 0.2840	 0.2720
CF	 0.4310	 0.3320
D	 0.7140	 0.4810
DA	 0.4060	 0.3830
DB	 0.4970	 0.3960
DC	 0.4970	 0.3830
DD	 0.4360	 0.3500
DE	 0.4270	 0.3560
DF	 0.4120	 0.3580
DG	 0.4100	 0.3500
DH	 0.3960	 0.3480
DI	 0.3790	 0.3100
DJ	 0.3150	 0.3040
DK	 0.2870	 0.2850
E	 0.7260	 0.4810
F	 0.7200	 0.4810
G	 0.7180	 0.4820





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Chain	Atom inclusion	Q-score
H	 0.7160	 0.4900
I	 0.7230	 0.4880
J	 0.7160	 0.4900
K	 0.7190	 0.4890
L	 0.7140	 0.4890
M	 0.7120	 0.4900
N	 0.7300	 0.4960
O	 0.7130	 0.4890
P	 0.7150	 0.4870
Q	 0.7130	 0.4880
R	 0.7120	 0.4830
S	 0.7110	 0.4870
T	 0.7110	 0.4870
U	 0.7030	 0.4900
V	 0.7200	 0.4820
W	 0.7160	 0.4880
X	 0.6880	 0.4780
a	 0.7420	 0.4850
b	 0.6800	 0.4740
c	 0.7360	 0.4860
d	 0.7430	 0.4800
e	 0.6950	 0.4720
f	 0.7280	 0.4770
g	 0.6700	 0.4390
h	 0.7270	 0.4790
i	 0.7200	 0.4790
j	 0.7310	 0.4770
k	 0.7320	 0.4580
l	 0.6740	 0.4500
m	 0.7240	 0.4600
n	 0.7360	 0.4630
o	 0.6830	 0.4460
p	 0.6740	 0.4580
q	 0.6510	 0.4310
r	 0.6170	 0.3770
s	 0.6650	 0.4380
t	 0.6440	 0.4320
u	 0.6200	 0.4200
v	 0.5100	 0.3790
w	 0.5070	 0.3620
x	 0.4070	 0.3360
y	 0.5140	 0.3650

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Chain	Atom inclusion	Q-score
z	 0.4910	 0.3630