



wwPDB EM Validation Summary Report ⓘ

Mar 8, 2026 – 03:59 PM UTC

PDB ID : 7MUV / pdb_00007muv
EMDB ID : EMD-24023
Title : Reconstruction of the Legionella pneumophila Dot/Icm T4SS 3DVA Map 3
Authors : Sheedlo, M.J.; Durie, C.L.; Swanson, M.; Lacy, D.B.; Ohi, M.D.
Deposited on : 2021-05-14
Resolution : 4.60 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

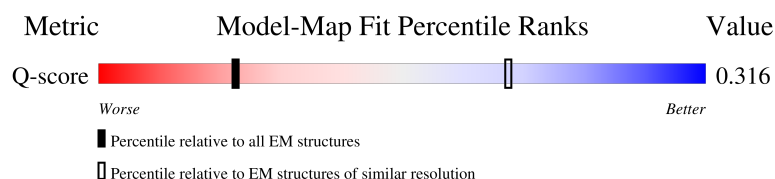
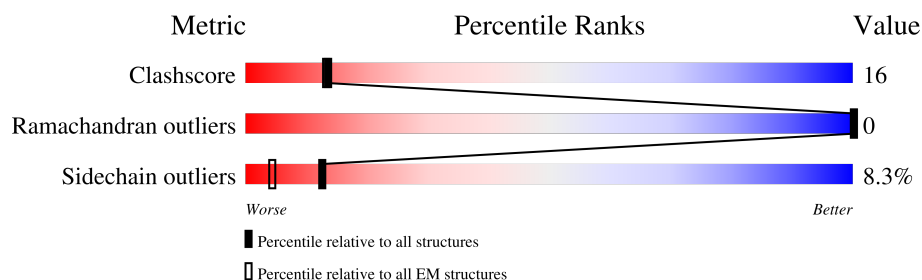
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 4.60 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.



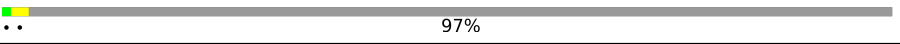
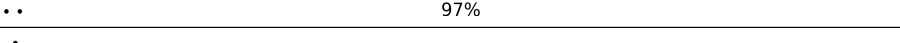
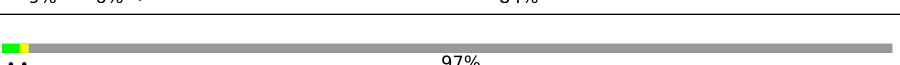
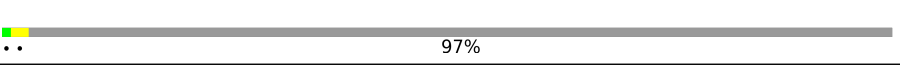
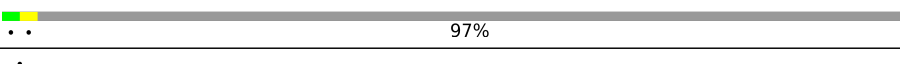
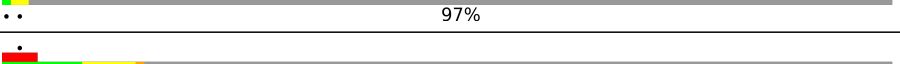
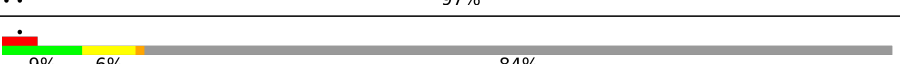
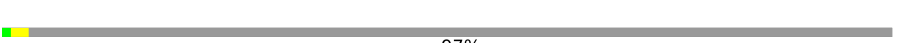
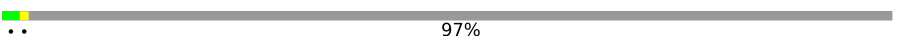
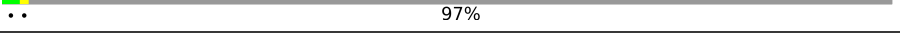
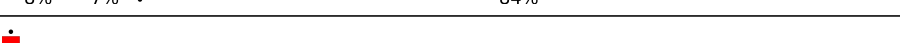
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	2407 (4.10 - 5.10)

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	AG	1048	9% 6% . 84%
1	Ag	1048	. . 97%
1	BG	1048	10% 6% . 84%
1	Bg	1048	. . 97%

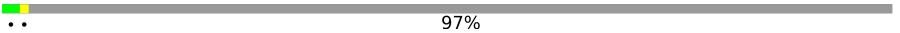
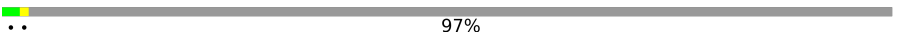
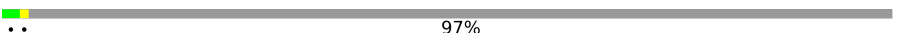
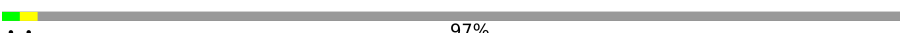
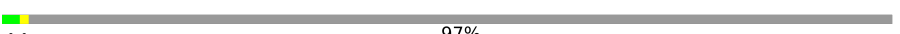
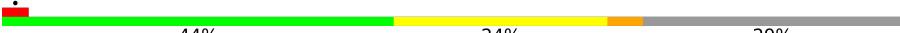
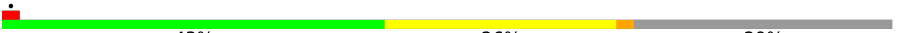
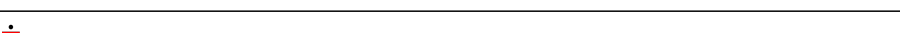
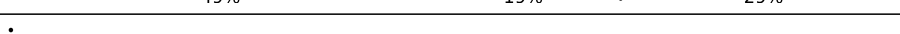
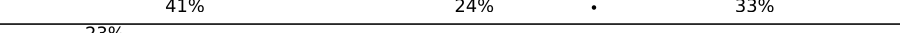
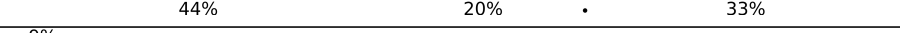
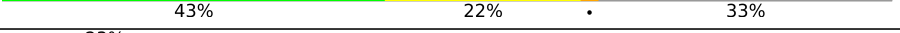
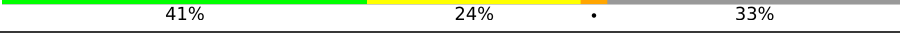
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Mol	Chain	Length	Quality of chain
1	CG	1048	 9% 6% . 84%
1	Cg	1048	 97%
1	DG	1048	 9% 6% . 84%
1	Dg	1048	 97%
1	EG	1048	 9% 6% . 84%
1	Eg	1048	 97%
1	FG	1048	 9% 6% . 84%
1	Fg	1048	 97%
1	GG	1048	 9% 6% . 84%
1	Gg	1048	 97%
1	HG	1048	 8% 7% . 84%
1	Hg	1048	 97%
1	IG	1048	 8% 7% . 84%
1	Ig	1048	 97%
1	JG	1048	 9% 6% . 84%
1	Jg	1048	 97%
1	KG	1048	 9% 6% . 84%
1	Kg	1048	 97%
1	LG	1048	 9% 6% . 84%
1	Lg	1048	 97%
1	MG	1048	 10% 6% . 84%
1	Mg	1048	 97%
1	NG	1048	 8% 7% . 84%
1	OG	1048	 8% 7% . 84%
1	PG	1048	 9% 6% . 84%

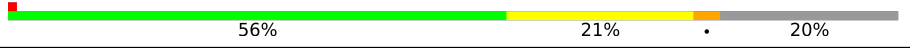
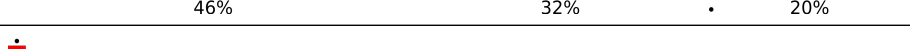
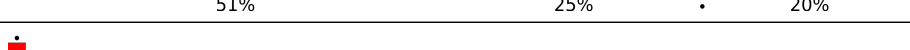
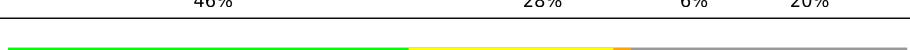
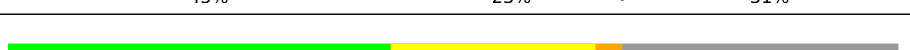
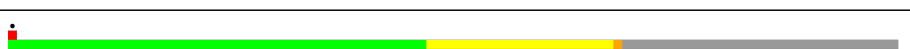
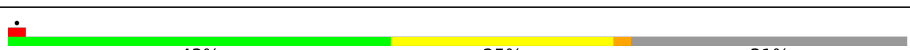
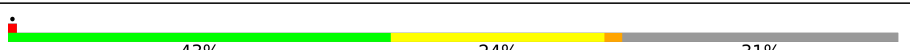
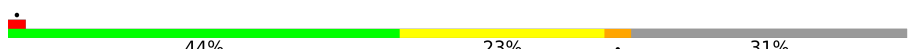
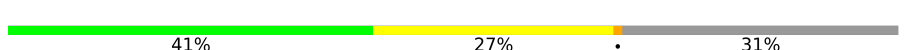
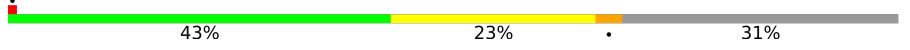
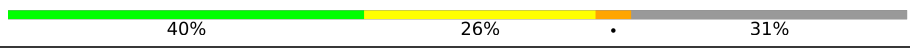
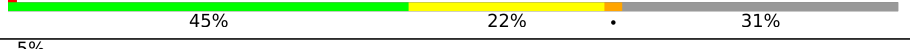
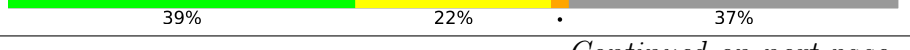
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Mol	Chain	Length	Quality of chain
1	VG	1048	 97%
1	WG	1048	 97%
1	XG	1048	 97%
1	YG	1048	 97%
1	ZG	1048	 97%
2	AH	361	 44% 24% • 29%
2	BH	361	 43% 26% • 29%
2	CH	361	 48% 22% • 29%
2	DH	361	 45% 25% • 29%
2	EH	361	 41% 26% • 29%
2	FH	361	 45% 25% • 29%
2	GH	361	 47% 23% • 29%
2	HH	361	 47% 23% • 29%
2	IH	361	 47% 22% • 29%
2	JH	361	 49% 19% • 29%
2	KH	361	 47% 21% • 29%
2	LH	361	 44% 24% • 29%
2	MH	361	 45% 24% • 29%
2	VH	361	 12% 41% 24% • 23%
2	WH	361	 23% 44% 20% • 13%
2	XH	361	 9% 43% 22% • 26%
2	YH	361	 23% 41% 24% • 12%
2	ZH	361	 14% 44% 22% • 20%
3	AK	189	 52% 26% • 20%
3	BK	189	 53% 25% • 20%

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Mol	Chain	Length	Quality of chain
3	CK	189	
3	DK	189	
3	EK	189	
3	FK	189	
3	GK	189	
3	HK	189	
3	IK	189	
3	JK	189	
3	KK	189	
3	LK	189	
3	MK	189	
4	AL	249	
4	BL	249	
4	CL	249	
4	DL	249	
4	EL	249	
4	FL	249	
4	GL	249	
4	HL	249	
4	IL	249	
4	JL	249	
4	KL	249	
4	LL	249	
4	ML	249	
5	AN	124	

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Mol	Chain	Length	Quality of chain
5	BN	124	
5	CN	124	
5	DN	124	
5	EN	124	
5	FN	124	
5	GN	124	
5	HN	124	
5	IN	124	
5	JN	124	
5	KN	124	
5	LN	124	
5	MN	124	
6	AU	9	
6	BU	9	
6	CU	9	
6	DU	9	
6	EU	9	
6	FU	9	
6	GU	9	
6	HU	9	
6	IU	9	
6	JU	9	
6	KU	9	
6	LU	9	
6	MU	9	

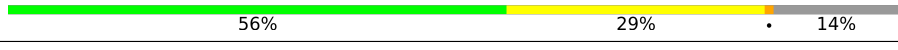
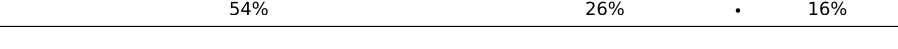
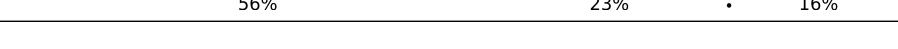
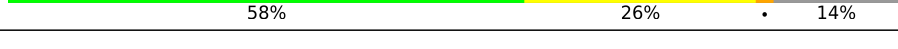
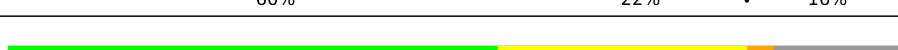
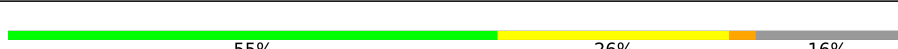
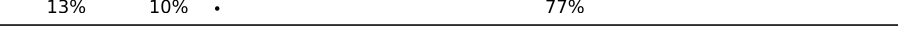
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Mol	Chain	Length	Quality of chain
7	AX	48	31% 92% 8%
7	BX	48	35% 85% 15%
7	CX	48	40% 96% .
7	DX	48	42% 98% .
7	EX	48	46% 100% .
7	FX	48	29% 100% .
7	GX	48	33% 100% .
7	HX	48	44% 96% .
7	IX	48	46% 98% .
7	JX	48	31% 96% .
7	KX	48	42% 98% .
7	LX	48	29% 100% .
7	MX	48	31% 90% 10%
7	VX	48	42% 98% .
7	WX	48	40% 96% .
7	XX	48	29% 98% .
7	YX	48	40% 88% 12%
7	ZX	48	38% 96% .
8	AD	163	51% 33% 14%
8	Ad	163	52% 30% 16%
8	BD	163	58% 27% 14%
8	Bd	163	56% 26% 16%
8	CD	163	63% 22% 14%
8	Cd	163	54% 28% 16%
8	DD	163	55% 29% 14%

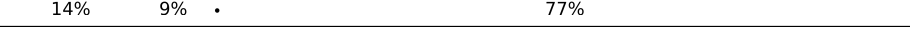
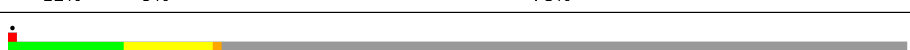
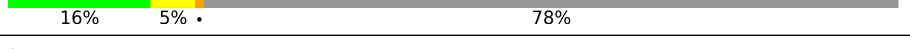
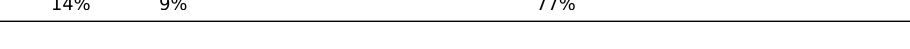
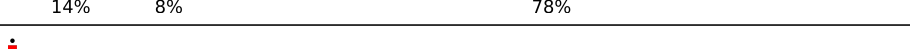
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Mol	Chain	Length	Quality of chain
8	Dd	163	
8	ED	163	
8	Ed	163	
8	FD	163	
8	Fd	163	
8	GD	163	
8	Gd	163	
8	HD	163	
8	Hd	163	
8	ID	163	
8	Id	163	
8	JD	163	
8	Jd	163	
8	KD	163	
8	Kd	163	
8	LD	163	
8	Ld	163	
8	MD	163	
8	Md	163	
9	AF	269	
9	Af	269	
9	BF	269	
9	Bf	269	
9	CF	269	
9	Cf	269	

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Mol	Chain	Length	Quality of chain
9	DF	269	 13% 10% 77%
9	Df	269	 13% 8% 78%
9	EF	269	 14% 9% 77%
9	Ef	269	 12% 9% 78%
9	FF	269	 13% 10% 77%
9	Ff	269	 12% 10% 78%
9	GF	269	 13% 9% 77%
9	Gf	269	 16% 5% 78%
9	HF	269	 14% 9% 77%
9	Hf	269	 16% 5% 78%
9	IF	269	 16% 7% 77%
9	If	269	 12% 8% 78%
9	JF	269	 14% 9% 77%
9	Jf	269	 15% 6% 78%
9	KF	269	 13% 10% 77%
9	Kf	269	 12% 10% 78%
9	LF	269	 15% 7% 77%
9	Lf	269	 13% 8% 78%
9	MF	269	 14% 8% 77%
9	Mf	269	 14% 8% 78%
9	VF	269	 13% 10% 77%
9	WF	269	 13% 10% 77%
9	XF	269	 12% 12% 77%
9	YF	269	 13% 9% 77%
9	ZF	269	17% 6% 77%

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Mol	Chain	Length	Quality of chain
10	AC	303	
10	BC	303	
10	CC	303	
10	DC	303	
10	EC	303	
10	FC	303	
10	GC	303	
10	HC	303	
10	IC	303	
10	JC	303	
10	KC	303	
10	LC	303	
10	MC	303	
11	AM	320	
11	BM	320	
11	CM	320	
11	DM	320	
11	EM	320	
11	FM	320	
11	GM	320	
11	HM	320	
11	IM	320	
11	JM	320	
11	KM	320	
11	LM	320	

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Mol	Chain	Length	Quality of chain
11	MM	320	

2 Entry composition [i](#)

There are 11 unique types of molecules in this entry. The entry contains 192370 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 IcmE protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	AG	165	Total 1229	C 780	N 203	O 242	S 4	0	0
1	Fg	34	Total 276	C 168	N 47	O 60	S 1	0	0
1	Gg	34	Total 276	C 168	N 47	O 60	S 1	0	0
1	BG	165	Total 1229	C 780	N 203	O 242	S 4	0	0
1	Hg	34	Total 276	C 168	N 47	O 60	S 1	0	0
1	Bg	34	Total 276	C 168	N 47	O 60	S 1	0	0
1	Ig	34	Total 276	C 168	N 47	O 60	S 1	0	0
1	CG	165	Total 1229	C 780	N 203	O 242	S 4	0	0
1	Jg	34	Total 276	C 168	N 47	O 60	S 1	0	0
1	Kg	34	Total 276	C 168	N 47	O 60	S 1	0	0
1	DG	165	Total 1229	C 780	N 203	O 242	S 4	0	0
1	Lg	34	Total 276	C 168	N 47	O 60	S 1	0	0
1	Mg	34	Total 276	C 168	N 47	O 60	S 1	0	0
1	EG	165	Total 1229	C 780	N 203	O 242	S 4	0	0
1	VG	34	Total 276	C 168	N 47	O 60	S 1	0	0
1	WG	34	Total 276	C 168	N 47	O 60	S 1	0	0
1	XG	34	Total 276	C 168	N 47	O 60	S 1	0	0

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Mol	Chain	Residues	Atoms					AltConf	Trace
1	YG	34	Total	C	N	O	S	0	0
			276	168	47	60	1		
1	ZG	34	Total	C	N	O	S	0	0
			276	168	47	60	1		
1	FG	165	Total	C	N	O	S	0	0
			1229	780	203	242	4		
1	Cg	34	Total	C	N	O	S	0	0
			276	168	47	60	1		
1	GG	165	Total	C	N	O	S	0	0
			1229	780	203	242	4		
1	Dg	34	Total	C	N	O	S	0	0
			276	168	47	60	1		
1	HG	165	Total	C	N	O	S	0	0
			1229	780	203	242	4		
1	IG	165	Total	C	N	O	S	0	0
			1229	780	203	242	4		
1	JG	165	Total	C	N	O	S	0	0
			1229	780	203	242	4		
1	KG	165	Total	C	N	O	S	0	0
			1229	780	203	242	4		
1	LG	165	Total	C	N	O	S	0	0
			1229	780	203	242	4		
1	MG	165	Total	C	N	O	S	0	0
			1229	780	203	242	4		
1	NG	165	Total	C	N	O	S	0	0
			1229	780	203	242	4		
1	OG	165	Total	C	N	O	S	0	0
			1229	780	203	242	4		
1	PG	165	Total	C	N	O	S	0	0
			1229	780	203	242	4		
1	Ag	34	Total	C	N	O	S	0	0
			276	168	47	60	1		
1	Eg	34	Total	C	N	O	S	0	0
			276	168	47	60	1		

- Molecule 2 is a protein called Type IV secretion protein IcmK.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	EH	258	Total	C	N	O	S	0	0
			1983	1268	336	371	8		
2	FH	258	Total	C	N	O	S	0	0
			1983	1268	336	371	8		

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Mol	Chain	Residues	Atoms					AltConf	Trace
2	GH	258	Total	C	N	O	S	0	0
			1983	1268	336	371	8		
2	HH	258	Total	C	N	O	S	0	0
			1983	1268	336	371	8		
2	IH	258	Total	C	N	O	S	0	0
			1983	1268	336	371	8		
2	JH	258	Total	C	N	O	S	0	0
			1983	1268	336	371	8		
2	AH	258	Total	C	N	O	S	0	0
			1983	1268	336	371	8		
2	KH	258	Total	C	N	O	S	0	0
			1983	1268	336	371	8		
2	LH	258	Total	C	N	O	S	0	0
			1983	1268	336	371	8		
2	MH	258	Total	C	N	O	S	0	0
			1983	1268	336	371	8		
2	VH	243	Total	C	N	O	S	0	0
			1875	1201	319	348	7		
2	WH	243	Total	C	N	O	S	0	0
			1875	1201	319	348	7		
2	XH	243	Total	C	N	O	S	0	0
			1875	1201	319	348	7		
2	YH	243	Total	C	N	O	S	0	0
			1875	1201	319	348	7		
2	ZH	243	Total	C	N	O	S	0	0
			1875	1201	319	348	7		
2	BH	258	Total	C	N	O	S	0	0
			1983	1268	336	371	8		
2	CH	258	Total	C	N	O	S	0	0
			1983	1268	336	371	8		
2	DH	258	Total	C	N	O	S	0	0
			1983	1268	336	371	8		

- Molecule 3 is a protein called Inner membrane lipoprotein YiaD.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	EK	151	Total	C	N	O	S	0	0
			1175	747	209	215	4		
3	FK	151	Total	C	N	O	S	0	0
			1175	747	209	215	4		
3	GK	151	Total	C	N	O	S	0	0
			1175	747	209	215	4		

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Mol	Chain	Residues	Atoms					AltConf	Trace
3	HK	151	Total	C	N	O	S	0	0
			1175	747	209	215	4		
3	IK	151	Total	C	N	O	S	0	0
			1175	747	209	215	4		
3	JK	151	Total	C	N	O	S	0	0
			1175	747	209	215	4		
3	KK	151	Total	C	N	O	S	0	0
			1175	747	209	215	4		
3	LK	151	Total	C	N	O	S	0	0
			1175	747	209	215	4		
3	AK	151	Total	C	N	O	S	0	0
			1175	747	209	215	4		
3	MK	151	Total	C	N	O	S	0	0
			1175	747	209	215	4		
3	BK	151	Total	C	N	O	S	0	0
			1175	747	209	215	4		
3	CK	151	Total	C	N	O	S	0	0
			1175	747	209	215	4		
3	DK	151	Total	C	N	O	S	0	0
			1175	747	209	215	4		

- Molecule 4 is a protein called Outer membrane protein, OmpA family protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	EL	173	Total	C	N	O	S	0	0
			1388	877	253	253	5		
4	FL	173	Total	C	N	O	S	0	0
			1388	877	253	253	5		
4	GL	173	Total	C	N	O	S	0	0
			1388	877	253	253	5		
4	HL	173	Total	C	N	O	S	0	0
			1388	877	253	253	5		
4	IL	173	Total	C	N	O	S	0	0
			1388	877	253	253	5		
4	JL	173	Total	C	N	O	S	0	0
			1388	877	253	253	5		
4	KL	173	Total	C	N	O	S	0	0
			1388	877	253	253	5		
4	LL	173	Total	C	N	O	S	0	0
			1388	877	253	253	5		
4	ML	173	Total	C	N	O	S	0	0
			1388	877	253	253	5		

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Mol	Chain	Residues	Atoms					AltConf	Trace
4	AL	173	Total	C	N	O	S	0	0
			1388	877	253	253	5		
4	BL	173	Total	C	N	O	S	0	0
			1388	877	253	253	5		
4	CL	173	Total	C	N	O	S	0	0
			1388	877	253	253	5		
4	DL	173	Total	C	N	O	S	0	0
			1388	877	253	253	5		

- Molecule 5 is a protein called Neurogenic locus notch.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	EN	78	Total	C	N	O	S	0	0
			582	357	99	113	13		
5	FN	78	Total	C	N	O	S	0	0
			582	357	99	113	13		
5	GN	78	Total	C	N	O	S	0	0
			582	357	99	113	13		
5	HN	78	Total	C	N	O	S	0	0
			582	357	99	113	13		
5	IN	78	Total	C	N	O	S	0	0
			582	357	99	113	13		
5	JN	78	Total	C	N	O	S	0	0
			582	357	99	113	13		
5	KN	78	Total	C	N	O	S	0	0
			582	357	99	113	13		
5	LN	78	Total	C	N	O	S	0	0
			582	357	99	113	13		
5	MN	78	Total	C	N	O	S	0	0
			582	357	99	113	13		
5	AN	78	Total	C	N	O	S	0	0
			582	357	99	113	13		
5	BN	78	Total	C	N	O	S	0	0
			582	357	99	113	13		
5	CN	78	Total	C	N	O	S	0	0
			582	357	99	113	13		
5	DN	78	Total	C	N	O	S	0	0
			582	357	99	113	13		

- Molecule 6 is a protein called Unknown protein fragment.

Mol	Chain	Residues	Atoms				AltConf	Trace
6	EU	9	Total	C	N	O	0	0
			45	27	9	9		
6	FU	9	Total	C	N	O	0	0
			45	27	9	9		
6	GU	9	Total	C	N	O	0	0
			45	27	9	9		
6	HU	9	Total	C	N	O	0	0
			45	27	9	9		
6	IU	9	Total	C	N	O	0	0
			45	27	9	9		
6	JU	9	Total	C	N	O	0	0
			45	27	9	9		
6	KU	9	Total	C	N	O	0	0
			45	27	9	9		
6	LU	9	Total	C	N	O	0	0
			45	27	9	9		
6	MU	9	Total	C	N	O	0	0
			45	27	9	9		
6	AU	9	Total	C	N	O	0	0
			45	27	9	9		
6	BU	9	Total	C	N	O	0	0
			45	27	9	9		
6	CU	9	Total	C	N	O	0	0
			45	27	9	9		
6	DU	9	Total	C	N	O	0	0
			45	27	9	9		

- Molecule 7 is a protein called Unknown protein fragment.

Mol	Chain	Residues	Atoms				AltConf	Trace
7	EX	48	Total	C	N	O	0	0
			240	144	48	48		
7	FX	48	Total	C	N	O	0	0
			240	144	48	48		
7	GX	48	Total	C	N	O	0	0
			240	144	48	48		
7	HX	48	Total	C	N	O	0	0
			240	144	48	48		
7	IX	48	Total	C	N	O	0	0
			240	144	48	48		
7	JX	48	Total	C	N	O	0	0
			240	144	48	48		
7	KX	48	Total	C	N	O	0	0
			240	144	48	48		

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Mol	Chain	Residues	Atoms				AltConf	Trace
7	LX	48	Total	C	N	O	0	0
			240	144	48	48		
7	MX	48	Total	C	N	O	0	0
			240	144	48	48		
7	VX	48	Total	C	N	O	0	0
			240	144	48	48		
7	WX	48	Total	C	N	O	0	0
			240	144	48	48		
7	XX	48	Total	C	N	O	0	0
			240	144	48	48		
7	YX	48	Total	C	N	O	0	0
			240	144	48	48		
7	ZX	48	Total	C	N	O	0	0
			240	144	48	48		
7	AX	48	Total	C	N	O	0	0
			240	144	48	48		
7	BX	48	Total	C	N	O	0	0
			240	144	48	48		
7	CX	48	Total	C	N	O	0	0
			240	144	48	48		
7	DX	48	Total	C	N	O	0	0
			240	144	48	48		

- Molecule 8 is a protein called DotD.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	Ed	137	Total	C	N	O	S	0	0
			1058	672	182	202	2		
8	FD	140	Total	C	N	O	S	0	0
			1086	692	185	206	3		
8	Fd	137	Total	C	N	O	S	0	0
			1058	672	182	202	2		
8	GD	140	Total	C	N	O	S	0	0
			1086	692	185	206	3		
8	Gd	137	Total	C	N	O	S	0	0
			1058	672	182	202	2		
8	HD	140	Total	C	N	O	S	0	0
			1086	692	185	206	3		
8	Hd	137	Total	C	N	O	S	0	0
			1058	672	182	202	2		
8	ID	140	Total	C	N	O	S	0	0
			1086	692	185	206	3		

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Mol	Chain	Residues	Atoms					AltConf	Trace
8	Id	137	Total	C	N	O	S	0	0
			1058	672	182	202	2		
8	JD	140	Total	C	N	O	S	0	0
			1086	692	185	206	3		
8	Jd	137	Total	C	N	O	S	0	0
			1058	672	182	202	2		
8	KD	140	Total	C	N	O	S	0	0
			1086	692	185	206	3		
8	Kd	137	Total	C	N	O	S	0	0
			1058	672	182	202	2		
8	LD	140	Total	C	N	O	S	0	0
			1086	692	185	206	3		
8	CD	140	Total	C	N	O	S	0	0
			1086	692	185	206	3		
8	Ld	137	Total	C	N	O	S	0	0
			1058	672	182	202	2		
8	MD	140	Total	C	N	O	S	0	0
			1086	692	185	206	3		
8	Md	137	Total	C	N	O	S	0	0
			1058	672	182	202	2		
8	Ad	137	Total	C	N	O	S	0	0
			1058	672	182	202	2		
8	BD	140	Total	C	N	O	S	0	0
			1086	692	185	206	3		
8	Bd	137	Total	C	N	O	S	0	0
			1058	672	182	202	2		
8	Cd	137	Total	C	N	O	S	0	0
			1058	672	182	202	2		
8	DD	140	Total	C	N	O	S	0	0
			1086	692	185	206	3		
8	Dd	137	Total	C	N	O	S	0	0
			1058	672	182	202	2		
8	AD	140	Total	C	N	O	S	0	0
			1086	692	185	206	3		
8	ED	140	Total	C	N	O	S	0	0
			1086	692	185	206	3		

- Molecule 9 is a protein called DotF.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	Ef	59	Total	C	N	O	S	0	0
			449	290	77	81	1		

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Mol	Chain	Residues	Atoms					AltConf	Trace
9	FF	63	Total 483	C 308	N 84	O 90	S 1	0	0
9	Ff	59	Total 449	C 290	N 77	O 81	S 1	0	0
9	AF	63	Total 483	C 308	N 84	O 90	S 1	0	0
9	GF	63	Total 483	C 308	N 84	O 90	S 1	0	0
9	Gf	59	Total 449	C 290	N 77	O 81	S 1	0	0
9	HF	63	Total 483	C 308	N 84	O 90	S 1	0	0
9	Hf	59	Total 449	C 290	N 77	O 81	S 1	0	0
9	IF	63	Total 483	C 308	N 84	O 90	S 1	0	0
9	If	59	Total 449	C 290	N 77	O 81	S 1	0	0
9	JF	63	Total 483	C 308	N 84	O 90	S 1	0	0
9	Jf	59	Total 449	C 290	N 77	O 81	S 1	0	0
9	KF	63	Total 483	C 308	N 84	O 90	S 1	0	0
9	Kf	59	Total 449	C 290	N 77	O 81	S 1	0	0
9	LF	63	Total 483	C 308	N 84	O 90	S 1	0	0
9	Lf	59	Total 449	C 290	N 77	O 81	S 1	0	0
9	MF	63	Total 483	C 308	N 84	O 90	S 1	0	0
9	Mf	59	Total 449	C 290	N 77	O 81	S 1	0	0
9	VF	63	Total 483	C 308	N 84	O 90	S 1	0	0
9	CF	63	Total 483	C 308	N 84	O 90	S 1	0	0
9	WF	63	Total 483	C 308	N 84	O 90	S 1	0	0
9	XF	63	Total 483	C 308	N 84	O 90	S 1	0	0

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Mol	Chain	Residues	Atoms					AltConf	Trace
9	YF	63	Total 483	C 308	N 84	O 90	S 1	0	0
9	ZF	63	Total 483	C 308	N 84	O 90	S 1	0	0
9	Af	59	Total 449	C 290	N 77	O 81	S 1	0	0
9	BF	63	Total 483	C 308	N 84	O 90	S 1	0	0
9	Bf	59	Total 449	C 290	N 77	O 81	S 1	0	0
9	Cf	59	Total 449	C 290	N 77	O 81	S 1	0	0
9	DF	63	Total 483	C 308	N 84	O 90	S 1	0	0
9	Df	59	Total 449	C 290	N 77	O 81	S 1	0	0
9	EF	63	Total 483	C 308	N 84	O 90	S 1	0	0

- Molecule 10 is a protein called DotC.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	BC	243	Total 1921	C 1216	N 340	O 357	S 8	0	0
10	CC	209	Total 1667	C 1061	N 292	O 309	S 5	0	0
10	DC	209	Total 1667	C 1061	N 292	O 309	S 5	0	0
10	EC	209	Total 1667	C 1061	N 292	O 309	S 5	0	0
10	FC	243	Total 1921	C 1216	N 340	O 357	S 8	0	0
10	GC	209	Total 1667	C 1061	N 292	O 309	S 5	0	0
10	HC	209	Total 1667	C 1061	N 292	O 309	S 5	0	0
10	IC	209	Total 1667	C 1061	N 292	O 309	S 5	0	0
10	JC	243	Total 1921	C 1216	N 340	O 357	S 8	0	0
10	KC	243	Total 1921	C 1216	N 340	O 357	S 8	0	0

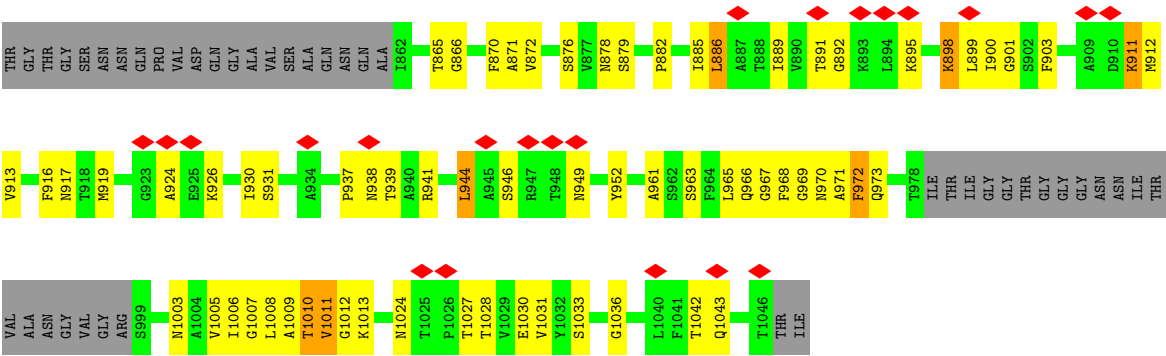
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Mol	Chain	Residues	Atoms					AltConf	Trace
10	LC	209	Total	C	N	O	S	0	0
			1667	1061	292	309	5		
10	MC	209	Total	C	N	O	S	0	0
			1667	1061	292	309	5		
10	AC	243	Total	C	N	O	S	0	0
			1921	1216	340	357	8		

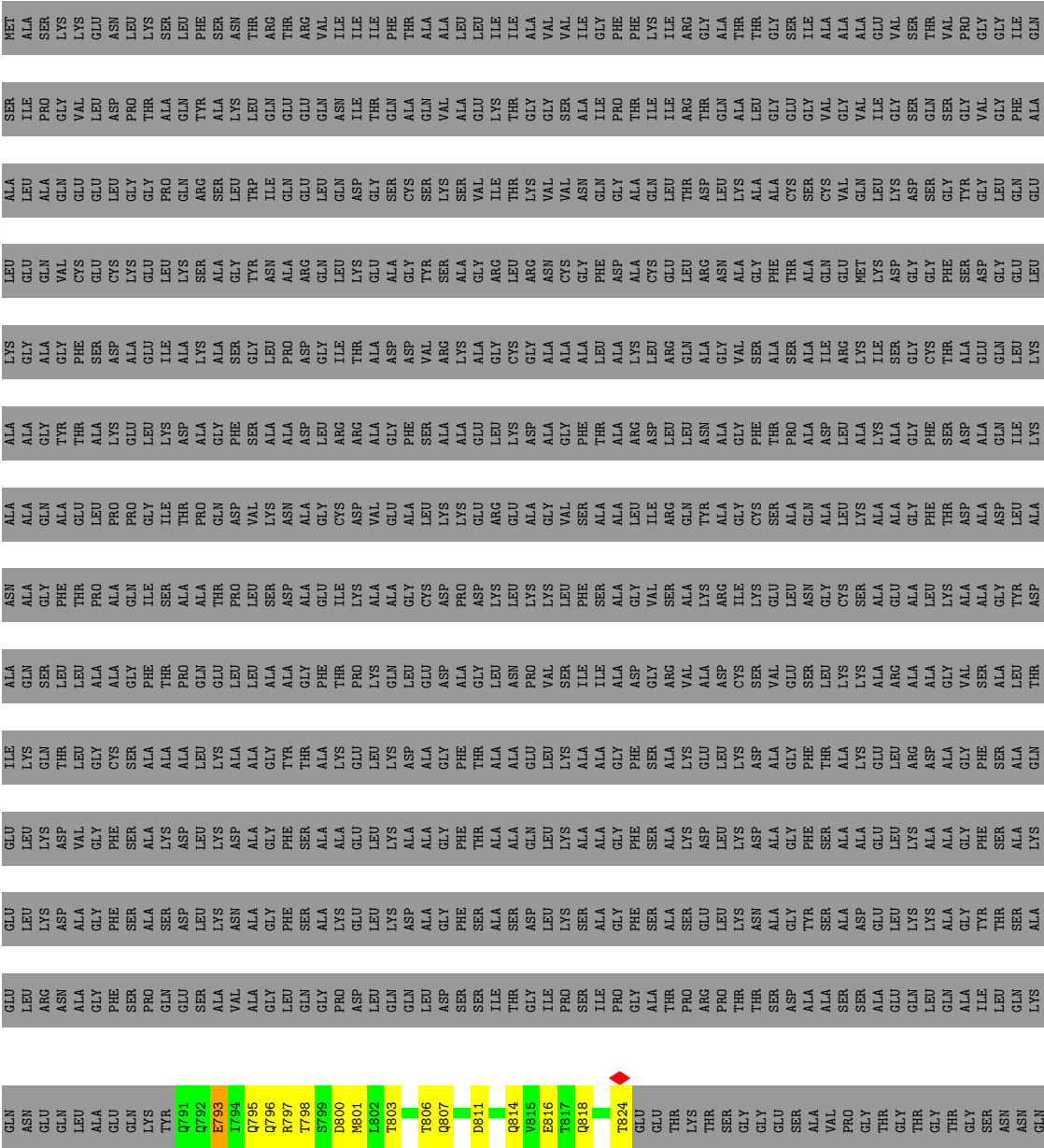
- Molecule 11 is a protein called DUF2807 domain-containing protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
11	HM	208	Total	C	N	O	S	0	0
			1650	1046	293	308	3		
11	IM	208	Total	C	N	O	S	0	0
			1650	1046	293	308	3		
11	KM	208	Total	C	N	O	S	0	0
			1650	1046	293	308	3		
11	LM	208	Total	C	N	O	S	0	0
			1650	1046	293	308	3		
11	MM	208	Total	C	N	O	S	0	0
			1650	1046	293	308	3		
11	AM	208	Total	C	N	O	S	0	0
			1650	1046	293	308	3		
11	BM	208	Total	C	N	O	S	0	0
			1650	1046	293	308	3		
11	EM	208	Total	C	N	O	S	0	0
			1650	1046	293	308	3		
11	FM	208	Total	C	N	O	S	0	0
			1650	1046	293	308	3		
11	GM	208	Total	C	N	O	S	0	0
			1650	1046	293	308	3		
11	JM	208	Total	C	N	O	S	0	0
			1650	1046	293	308	3		
11	CM	208	Total	C	N	O	S	0	0
			1650	1046	293	308	3		
11	DM	208	Total	C	N	O	S	0	0
			1650	1046	293	308	3		



● Molecule 1: IcmE protein

Chain Fg: .. 97%



[illegible]

- Molecule 1: IcmE protein

Chain Gg: 97%

[illegible]

THR THR VAL GLU VAL TYR SER GLY THR GLY LEU GLY ILE LEU PHE THR GLN ASP VAL THR THR ILE

- Molecule 1: IcmE protein



SER	ILE	PRO	GLY	VAL	LEU	ASP	PRO	THR	ALA	THR	GLN	TYR	ALA	LYS	LEU	GLN	GLU	GLU	GLN	ASN	ILE	THR	ILE	GLN	ALA	ALA	GLN	VAL	ALA	ALA	GLU	LYS	THR	GLY	GLY	SER	ALA	ILE	PRO	THR	ILE	ILE	ILE	ARG	THR	GLN	ALA	LEU	GLU	GLY	VAL	VAL	ILE	ILE	GLY	SER	GLN	SER	GLY	VAL	GLY	PHE	ALA
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ALA	LEU	ALA	ALA	GLN	GLY	GLU	LEU	GLY	PRO	GLN	ARG	SER	LEU	TRP	ILE	GLN	GLU	LEU	GLN	ASP	GLY	CYS	SER	LYS	SER	VAL	THR	LYS	VAL	ASN	GLY	ALA	GLN	LEU	THR	ASP	LEU	LYS	ALA	ALA	CYS	CYS	VAL	GLN	LEU	LYS	GLY	TYR	GLY	GLN	THR
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LEU GLU GLN GLN VAL CYS GLU CYS LYS LEU GLU LYS SER ALA GLY TYR GLY ASN ALA ALA ARG GLN LEU LYS GLU LYS GLY TYR SER ALA ALA GLY GLY PHE ASP LEU ARG ASN CYS GLY GLY PHE ASP ALA ALA CYS GLU GLU LEU ARG ASN GLY GLY THR THR ALA ALA GLN GLU GLU MET LYS ASP ASP ASP GLY GLY PHE SER SER ASP ASP GLY GLY LEU

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

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

ALA	ALA	ALA	GLN	ALA	ALA	GLU	LEU	PRO	PRO	GLY	ILE	THR	PRO	GLN	ASP	VAL	LYS	ASN	ALA	GLY	CYS	ASP	VAL	GLU	ALA	LEU	LYS	LYS	ARG	GLU	ALA	ALA	ALA	LEU	ILE	ARG	GLN	TYR	ALA	ALA	CYS	GLY	SER	ALA	GLN	GLN	LEU	LYS	ALA	ALA	GLY	PHE	THR	ASP	ASP	LEU
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ASN	ALA	GLY	PHE	THR	PRO	ALA	GLN	ILE	SER	ALA	ALA	THR	PRO	LEU	SER	ASP	ALA	GLU	ILE	LYS	ALA	ALA	ALA	GLY	CYS	ASP	ASP	ASP	LYS	LYS	LEU	LYS	PHE	SER	ALA	GLY	VAL	SER	ALA	LYS	ARG	ILE	LYS	GLU	LEU	ASN	GLY	CYS	SER	ALA	GLU	LEU	ALA	ALA	GLY	TYR	ASP
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ALA	GLN	SER	SER	LEU	LEU	ALA	ALA	GLY	PHE	THR	PRO	GLN	GLU	LEU	ALA	ALA	GLY	PHE	THR	PRO	LYS	GLN	GLU	ASP	ASN	PRO	VAL	SER	ILE	ILE	ALA	ASP	CYS	SER	GLU	VAL	GLU	SER	LEU	LYS	LYS	ALA	ALA	ARG	ALA	ALA	GLY	VAL	SER	SER	ALA	LEU	LEU
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ILE	LYS	LYS	GLN	THR	LEU	THR	GLY	CYS	SER	ALA	ALA	ALA	ALA	ALA	ALA	LYS	LEU	LEU	LYS	ALA	ALA	GLY	THR	THR	THR	GLY	PHE	ASP	LYS	LYS	GLU	GLU	LEU	GLY	ASP	ASP	ALA	ALA	GLY	PHE	SER	SER	ALA	ILE
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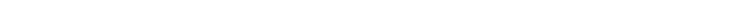
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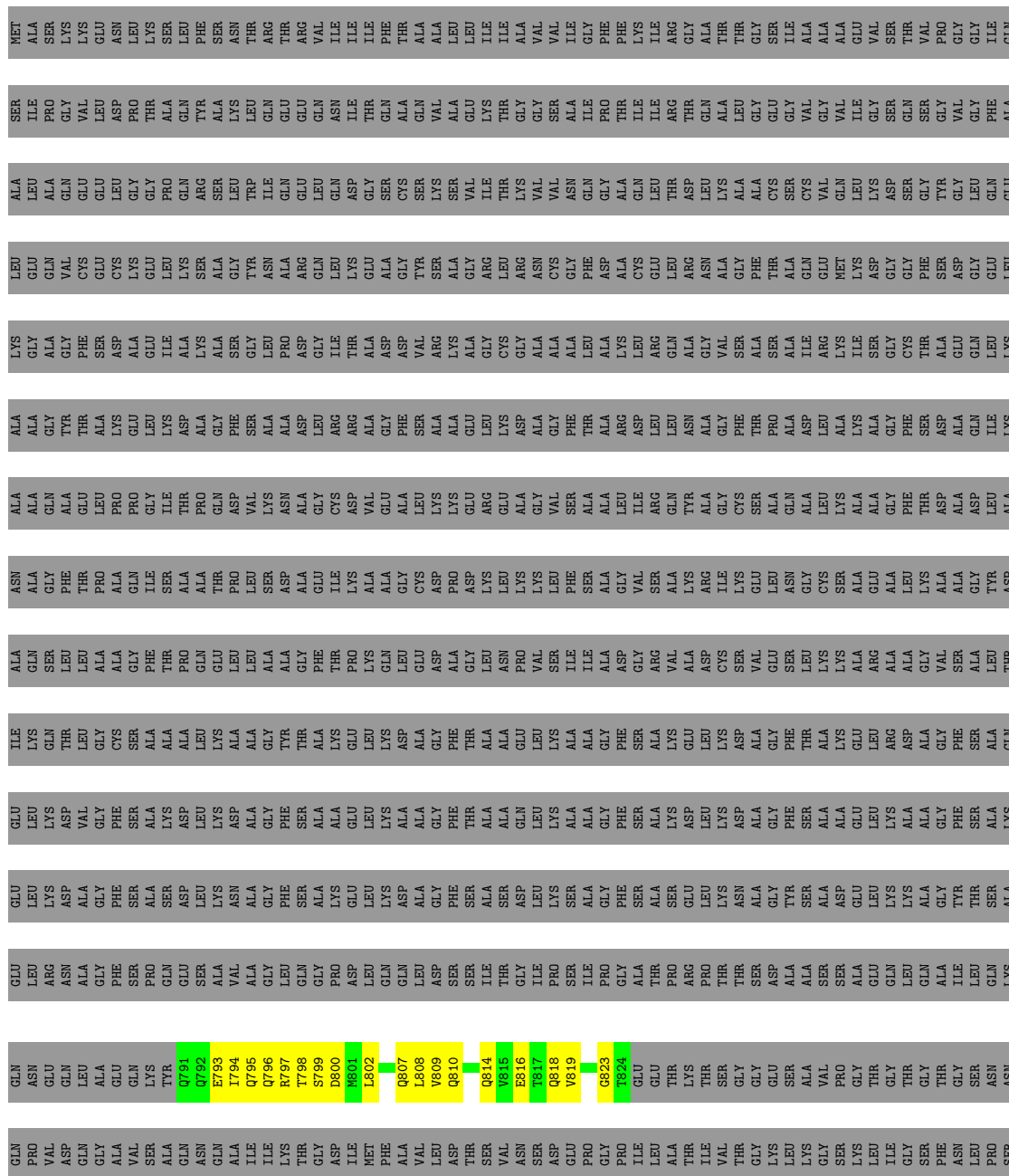
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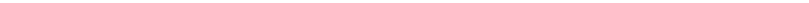
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THR
GLY
THR
GLY
SER
ASN
ASN
GLN
PRO
VAL
ASP
GLN
GLY
ALA
VAL
SER
ALA
GLN
ASN
GLN
ALA
1862
1863
C866
D867
1868
M869
F870
A871
W872
L873
S876
B877
N878
S879
P882
L886
A887
T888
I889
V890
T891
G892
K893
L894
K898
L899
I900
G901
S902
F903
N904
L905
H906

- Molecule 1: IcmE protein

Chain Hg:  97%

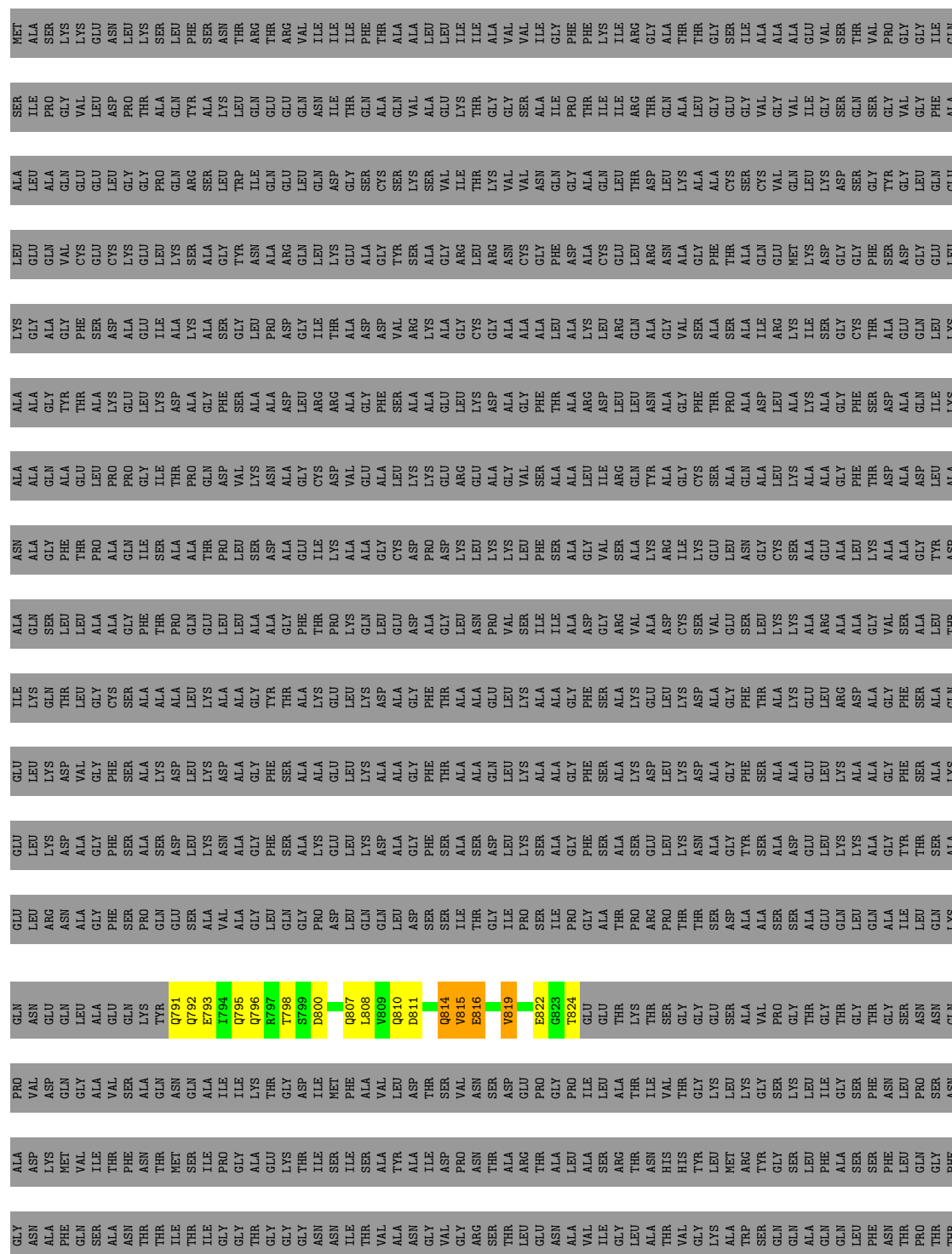


Chain Kg:  97%



- Molecule 1: IcmE protein

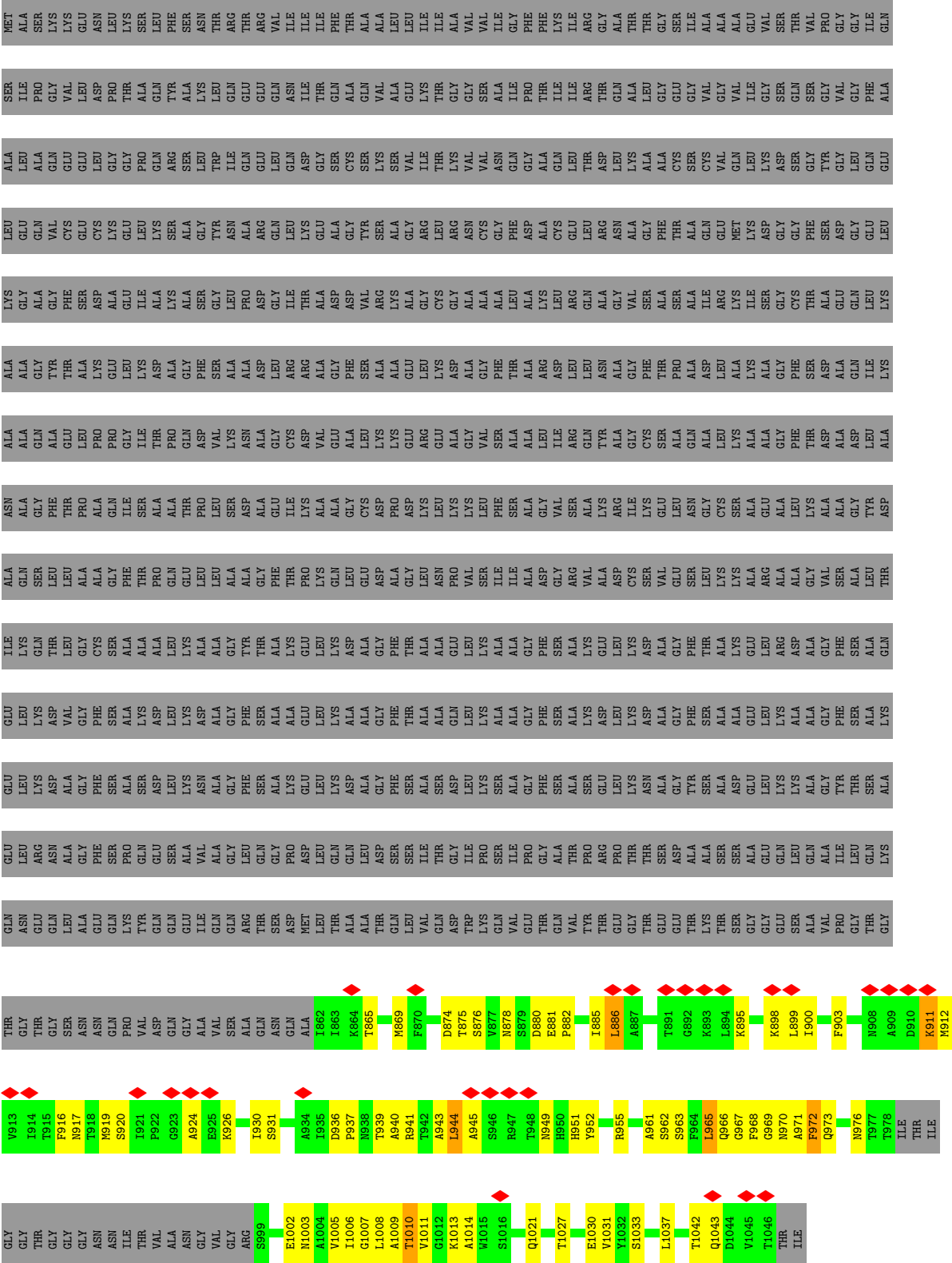
Chain Lg: 97%



Chain Mg: .. 97%

VAL
TYR
SER
GLY
THR
GLY
LEU
GLY
ILE
LEU
PHE
THR
GLN
ASP
VAL
THR
THR
ILE

● Molecule 1: IcmE protein



● Molecule 1: IcmE protein

97%

- Molecule 1: IcmE protein

97%





WORLDWIDE
PDB
PROTEIN DATA BANK

- Molecule 1: IcmE protein

- Molecule 1: IcmE protein

V1045	THR
T1046	ILE

Chain IG: 8% 7% 84%





- Molecule 1: IcmE protein

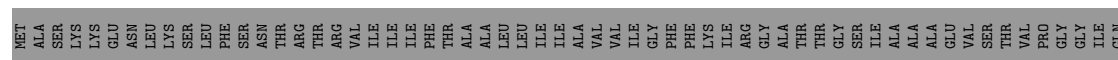
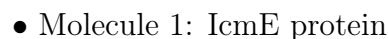
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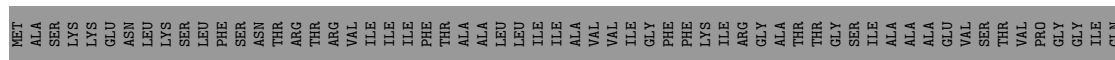
- Molecule 1: IcmE protein



MET	ALA	SER	LYS	GLU	ASN	LYS	SER	LEU	PHE	SER	THR	THR	ARG	ARG	VAL	ILE	ILE	ILE	PHE	THR	ALA	ALA	LEU	LEU	LEU	ILE	ILE	ILE	GLY	PHE	PHE	LYS	ILE	ILE	ARG	GLY	ALA	ALA	THR	THR	THR	SER	ILE	ALA	ALA	ALA	GLU	VAL	VAL	THR	THR	PRO	GLY	GLY	GLY	ILE.
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- Molecule 1: IcmE protein



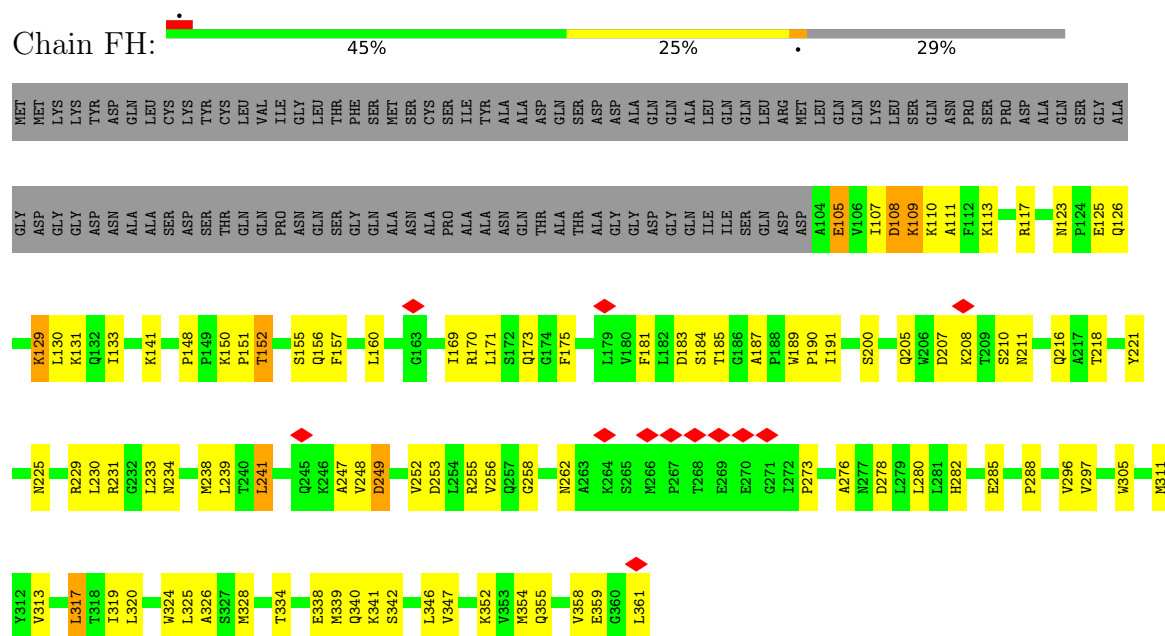




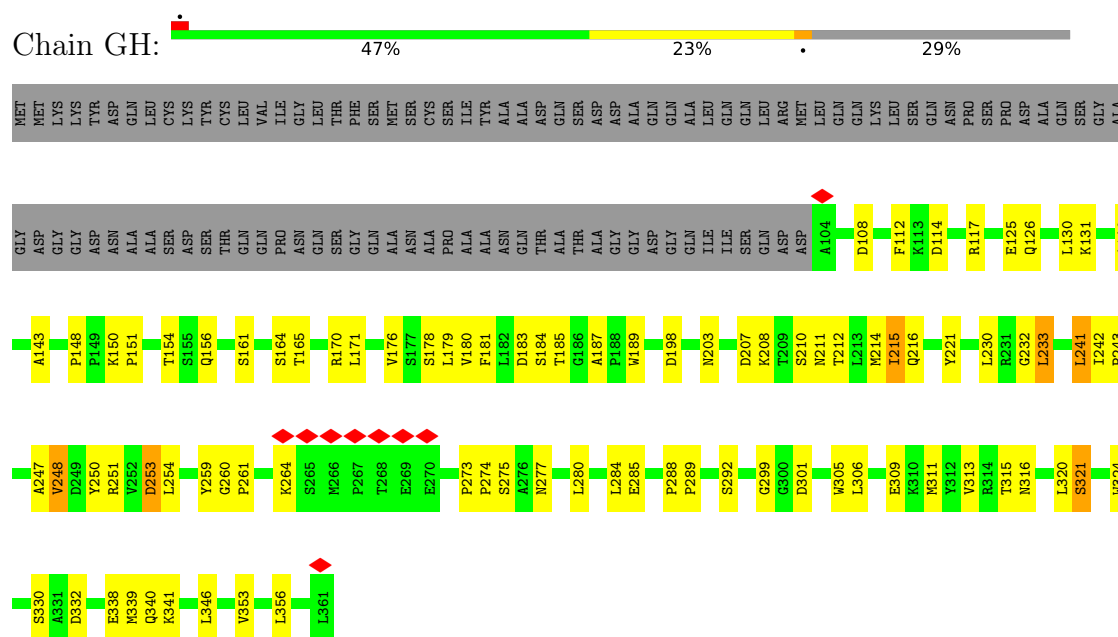




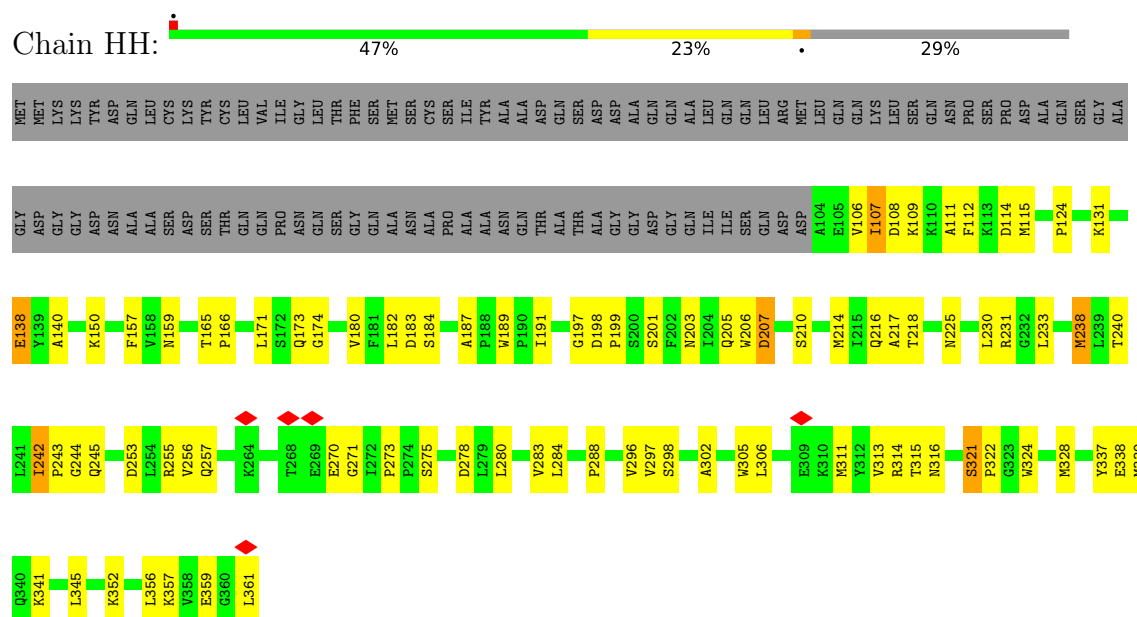
- Molecule 2: Type IV secretion protein IcmK



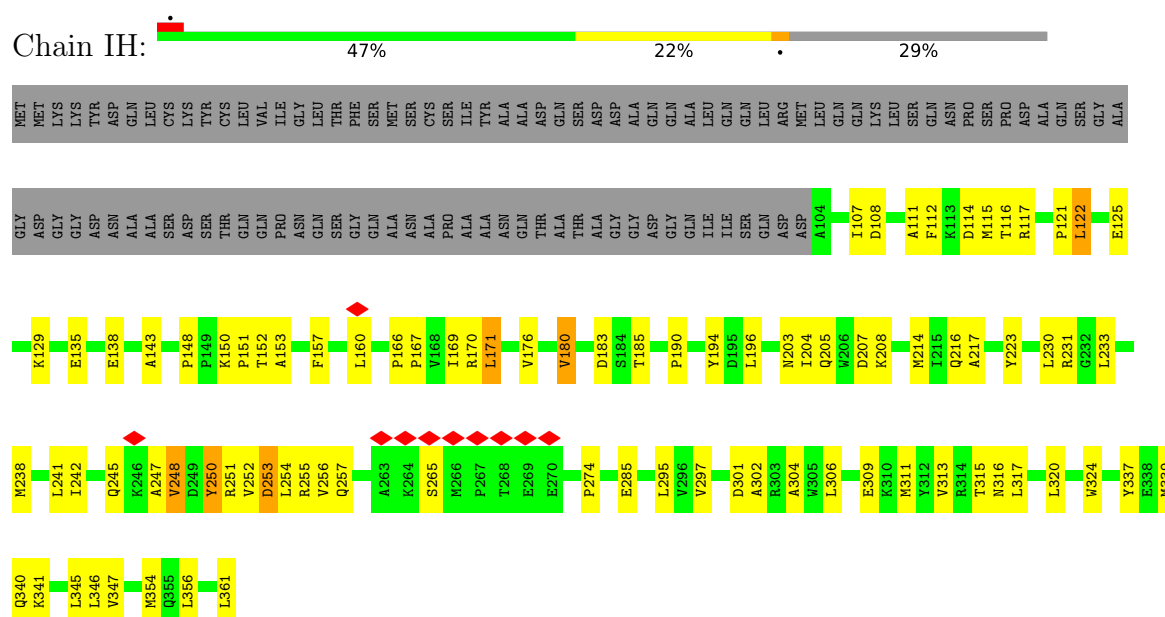
- Molecule 2: Type IV secretion protein IcmK



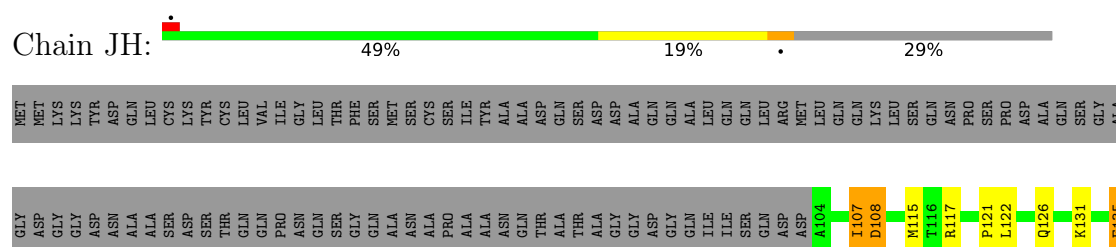
- Molecule 2: Type IV secretion protein IcmK

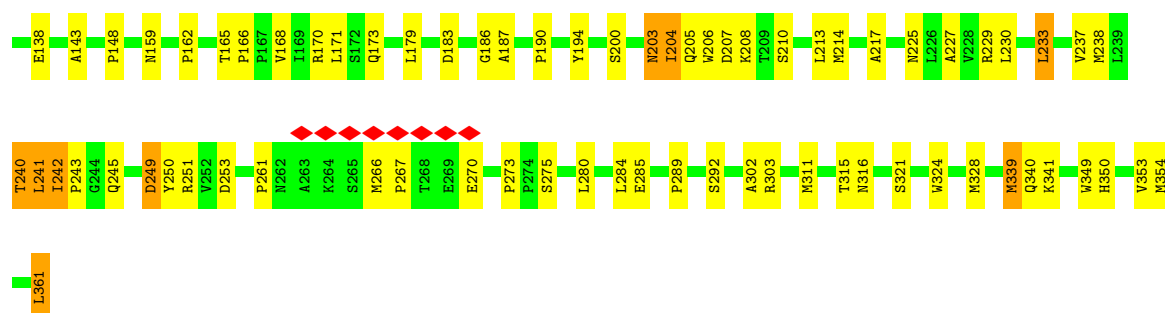


- Molecule 2: Type IV secretion protein IcmK

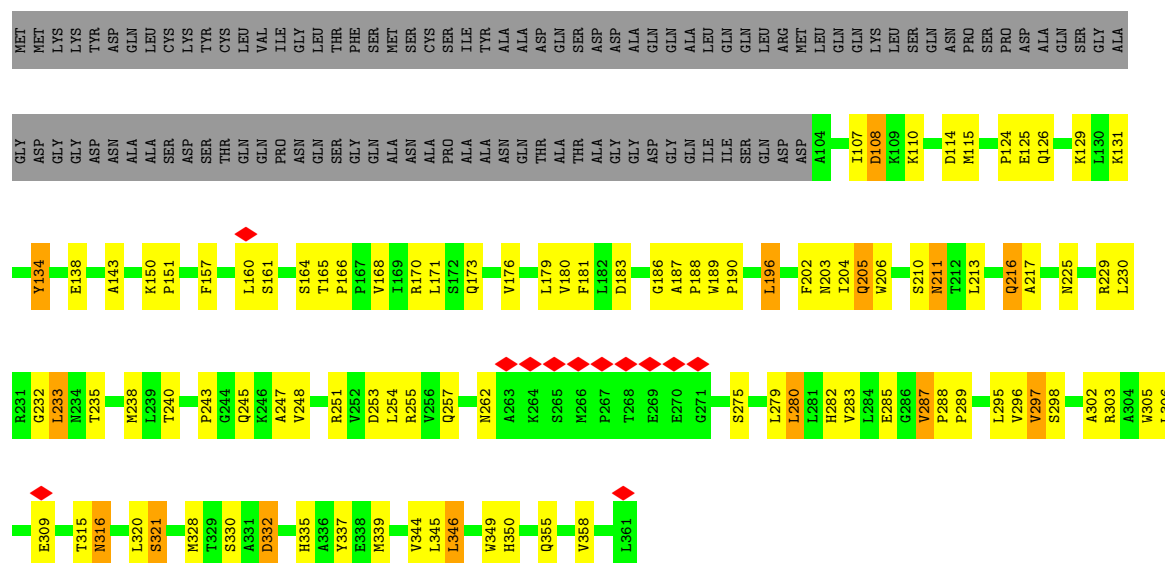
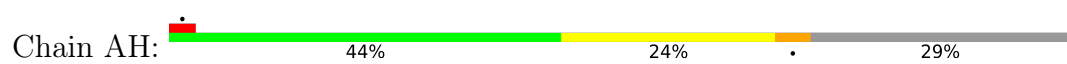


- Molecule 2: Type IV secretion protein IcmK

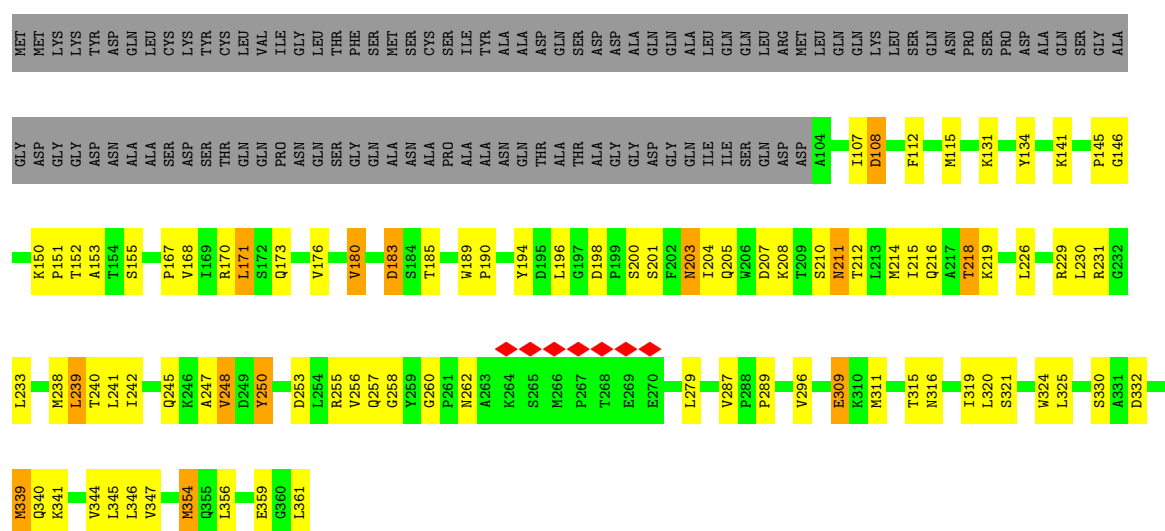




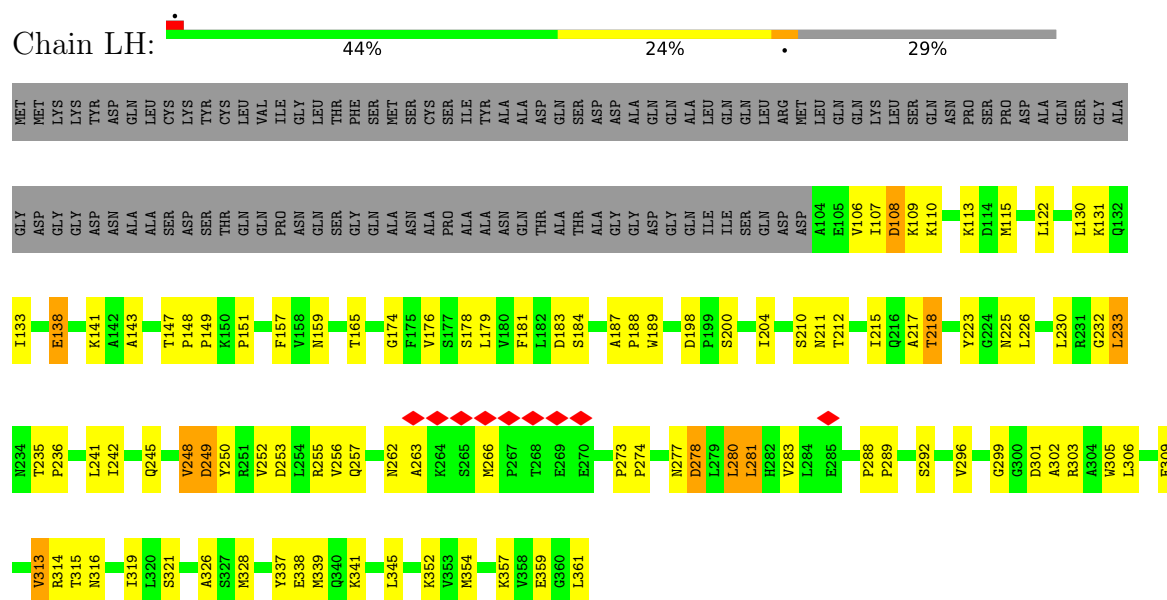
• Molecule 2: Type IV secretion protein IcmK



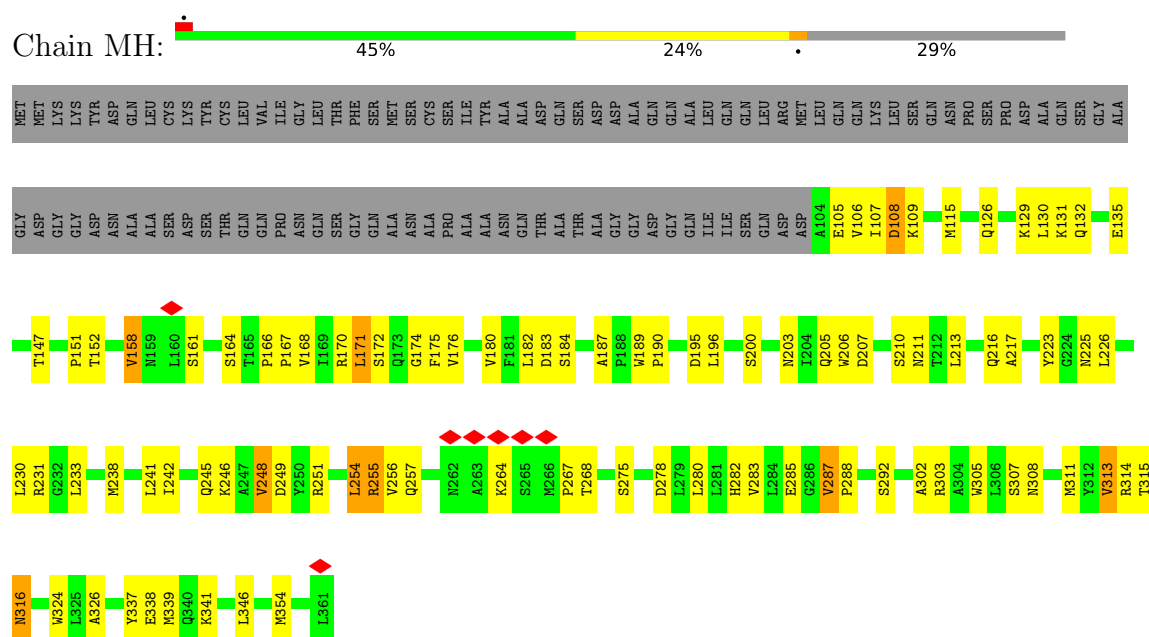
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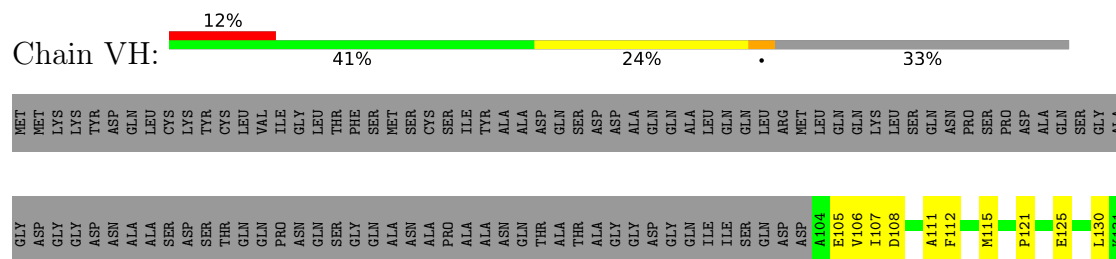
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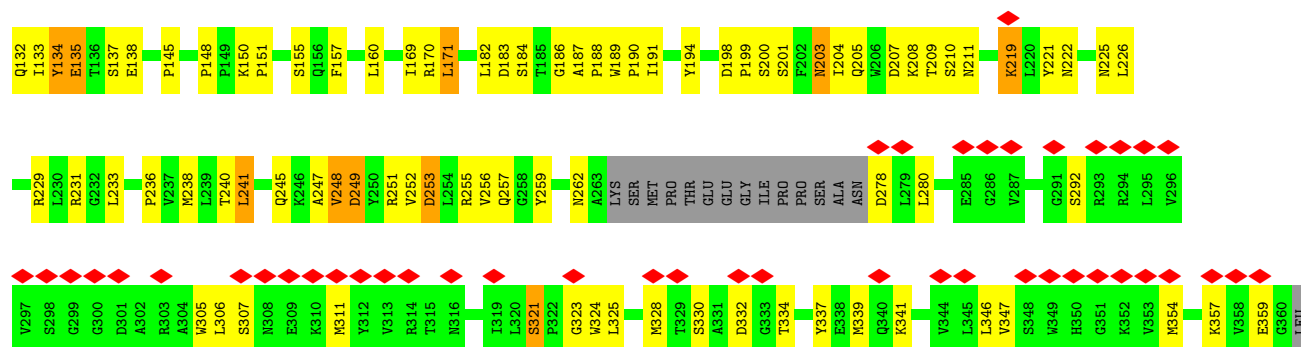


- Molecule 2: Type IV secretion protein IcmK

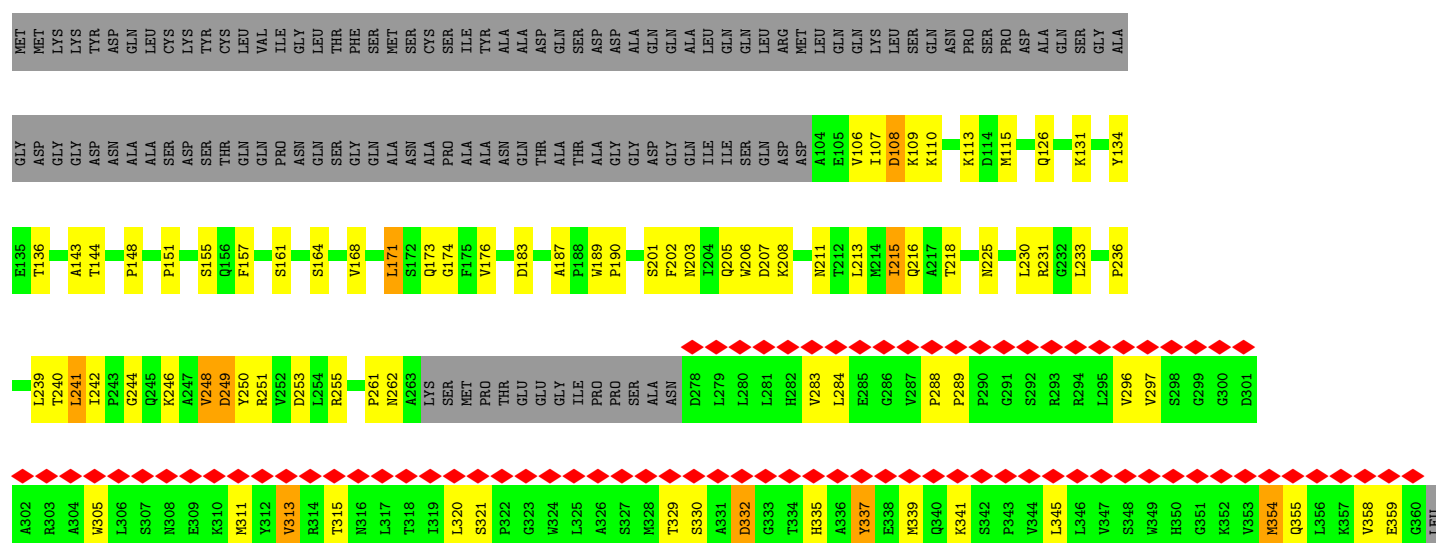


- Molecule 2: Type IV secretion protein IcmK

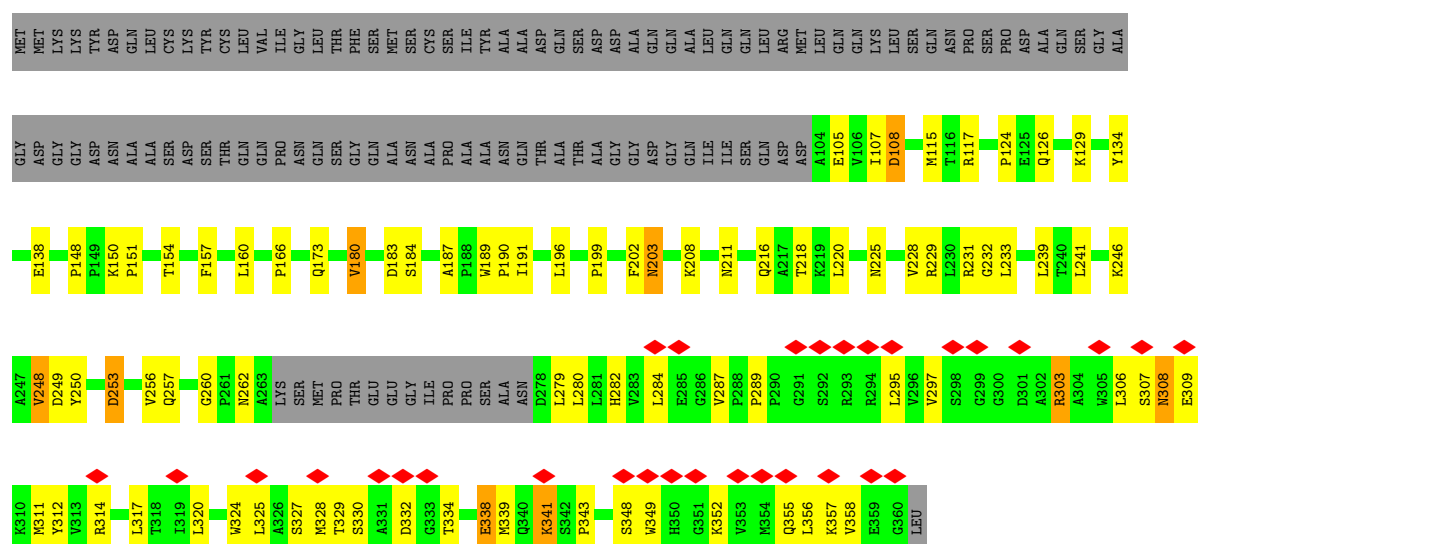




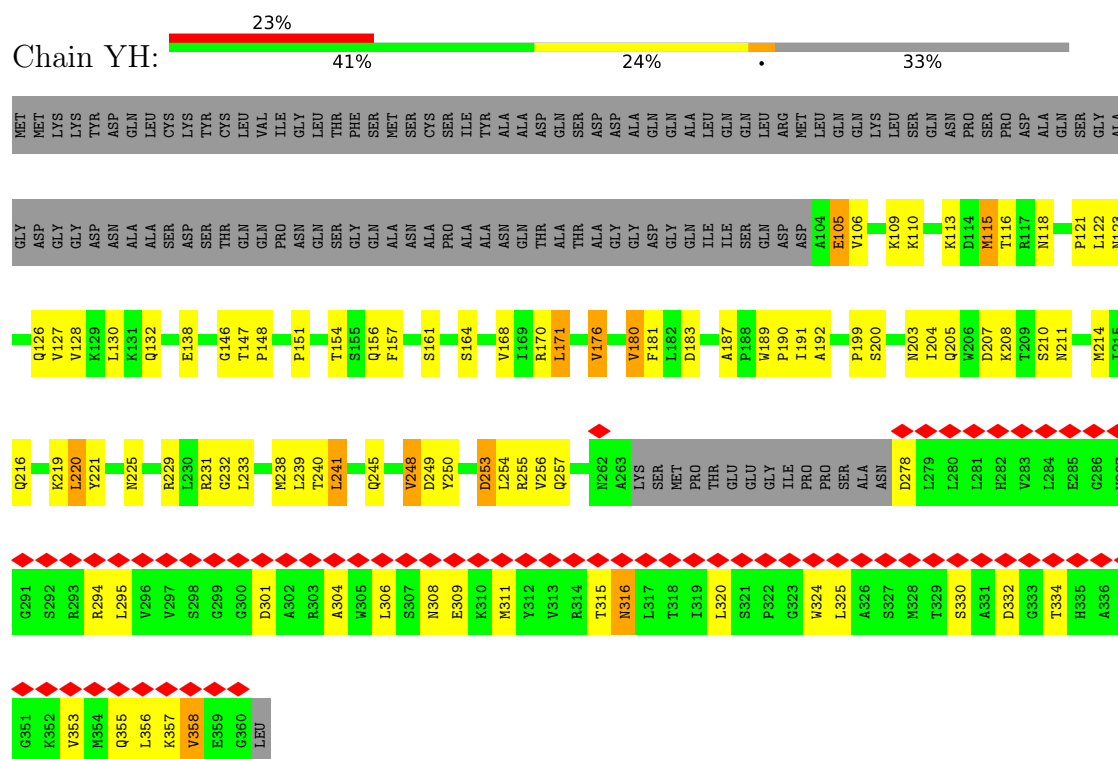
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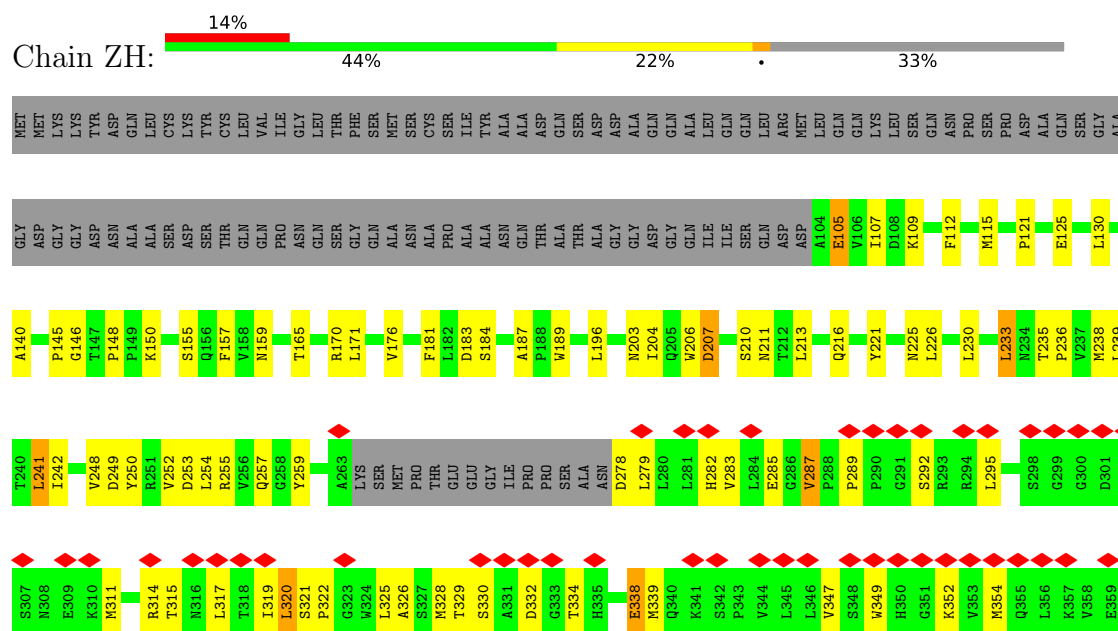
• Molecule 2: Type IV secretion protein IcmK



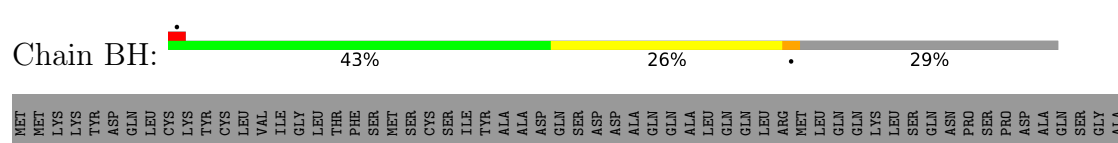
- Molecule 2: Type IV secretion protein IcmK



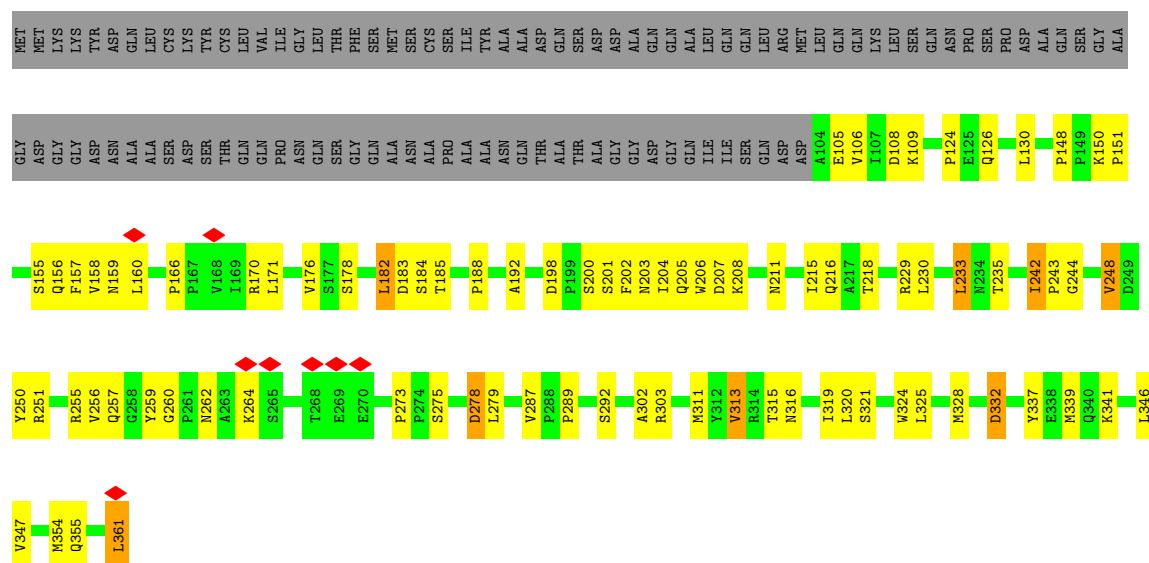
- Molecule 2: Type IV secretion protein IcmK



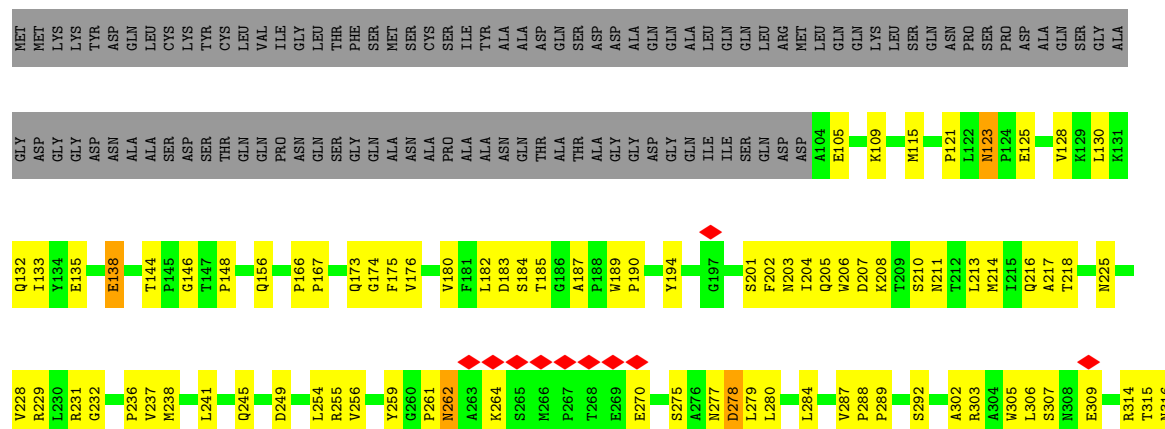
- Molecule 2: Type IV secretion protein IcmK

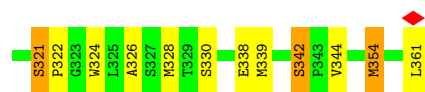


- Molecule 2: Type IV secretion protein IcmK

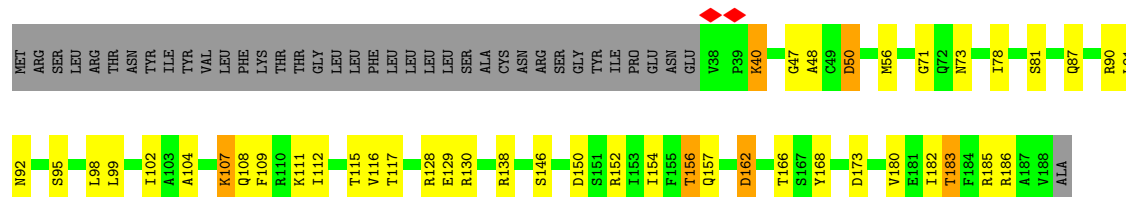


- Molecule 2: Type IV secretion protein IcmK

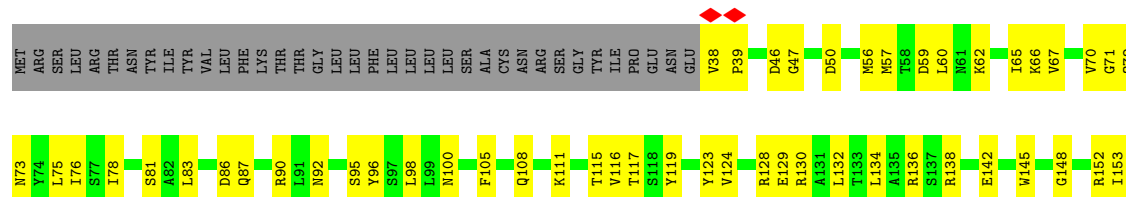




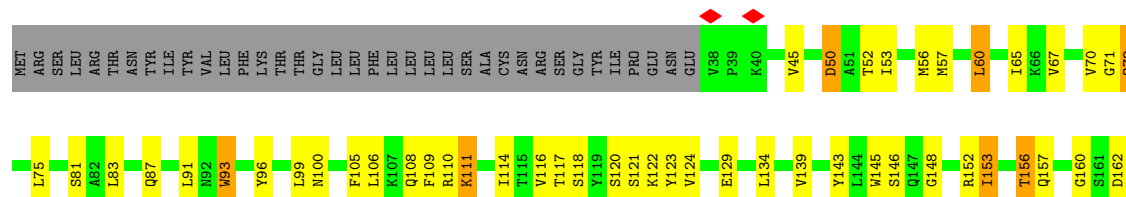
• Molecule 3: Inner membrane lipoprotein YiaD



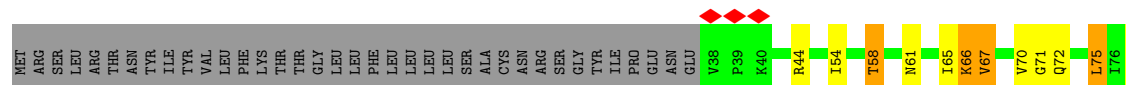
• Molecule 3: Inner membrane lipoprotein YiaD

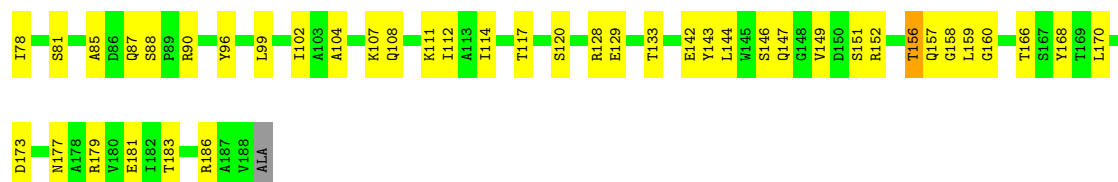


• Molecule 3: Inner membrane lipoprotein YiaD

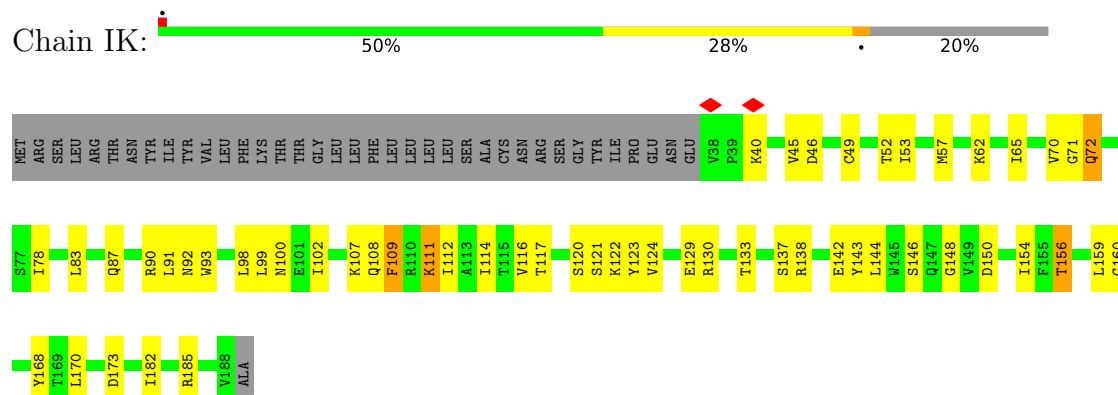


• Molecule 3: Inner membrane lipoprotein YiaD

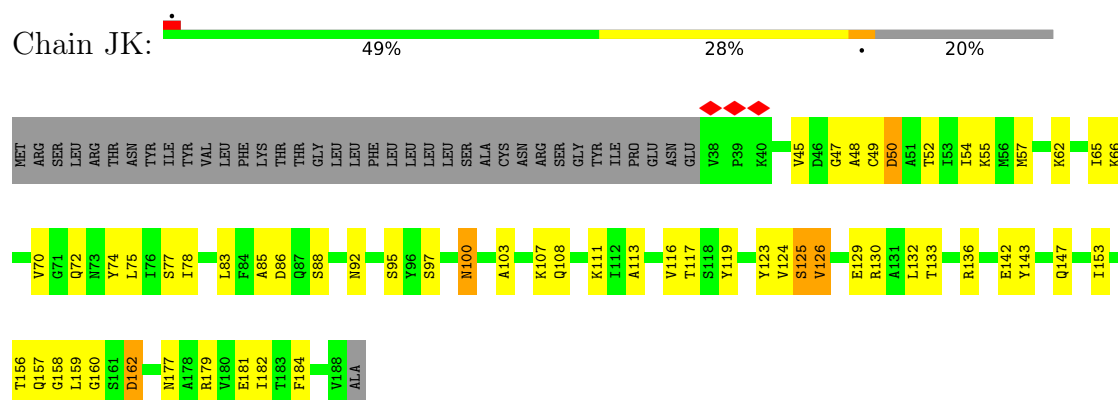




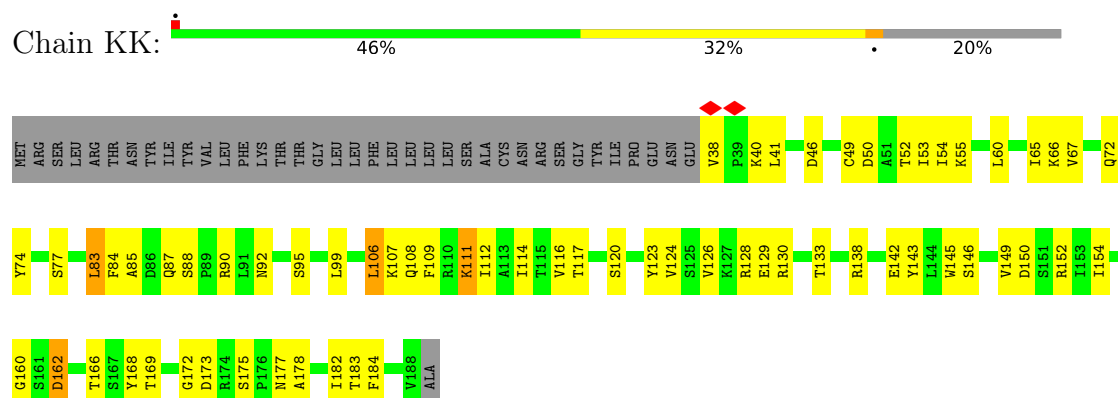
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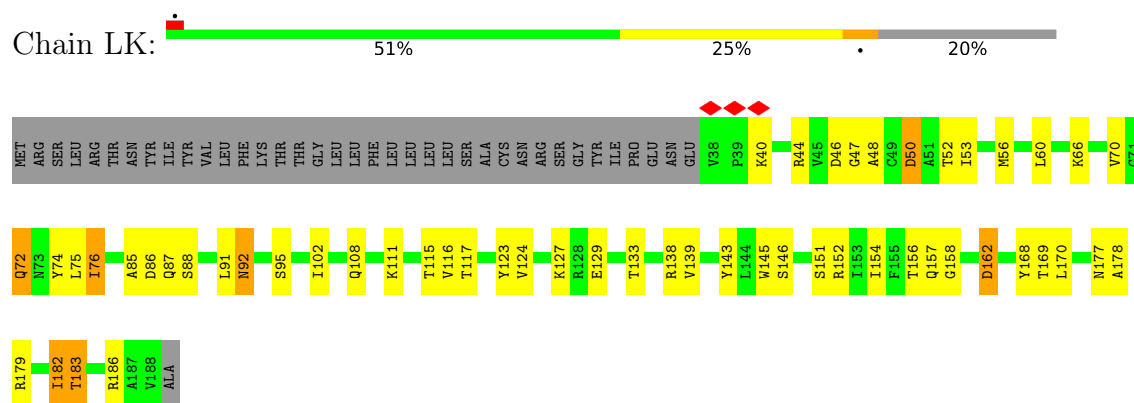
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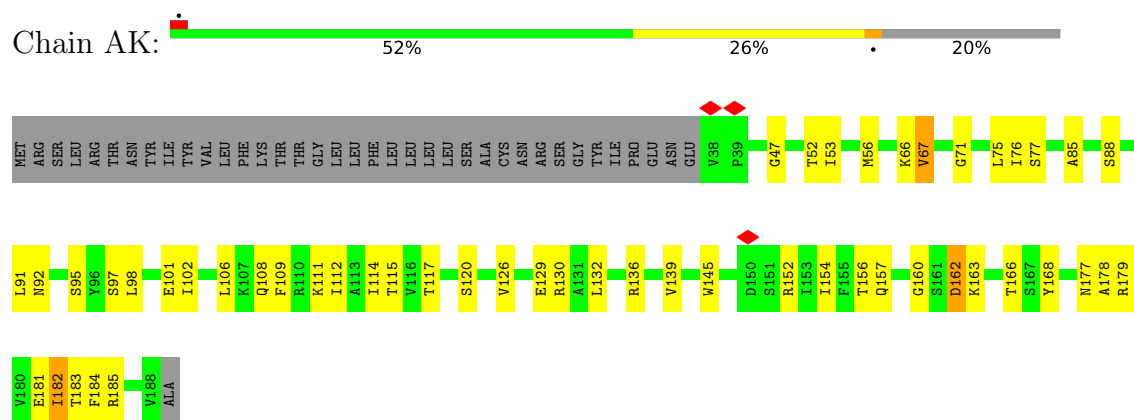
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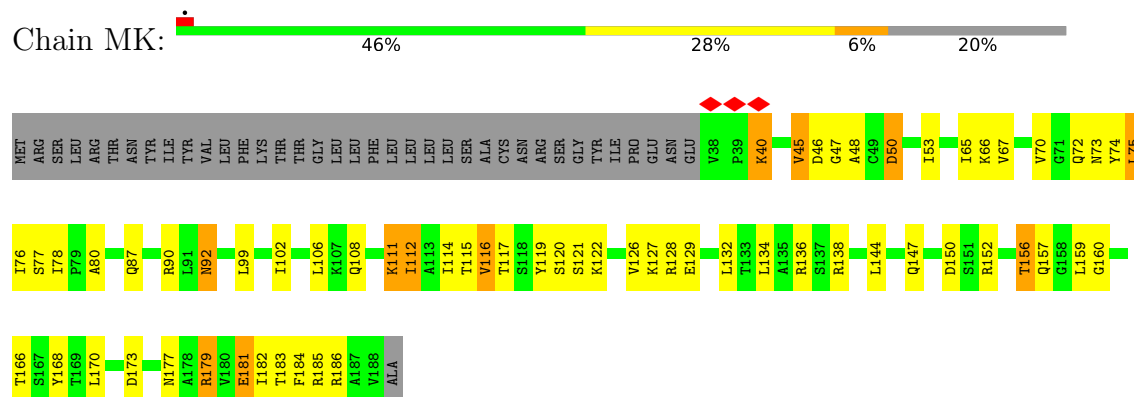
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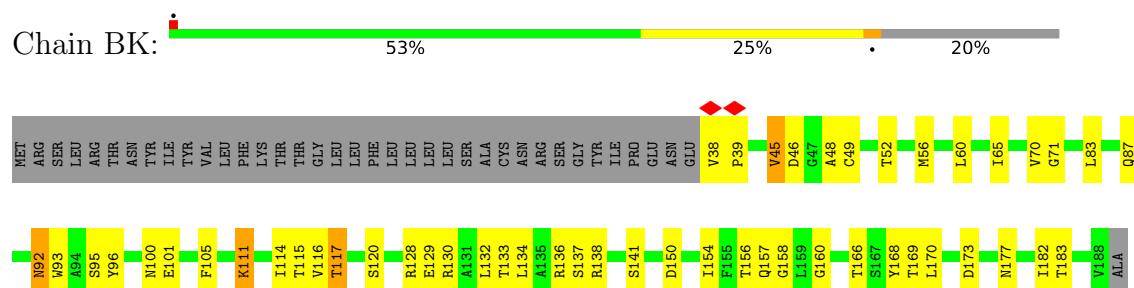
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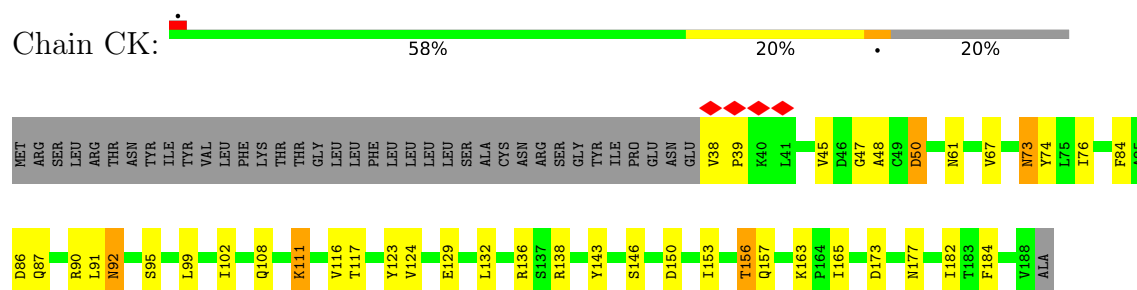
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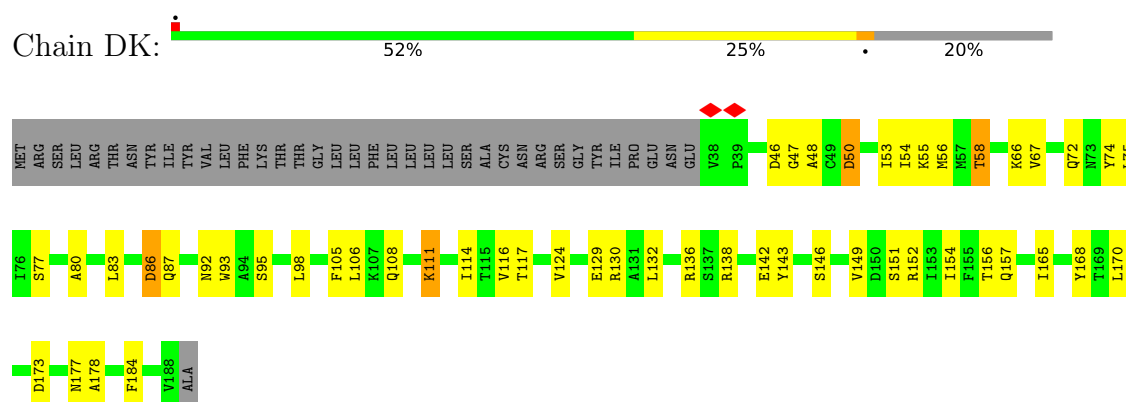
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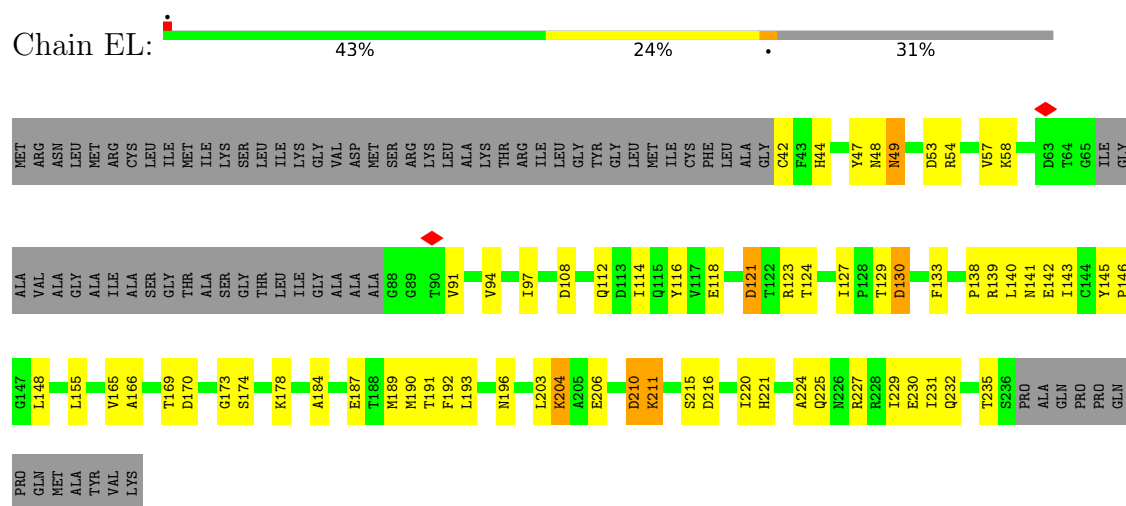
- Molecule 3: Inner membrane lipoprotein YiaD



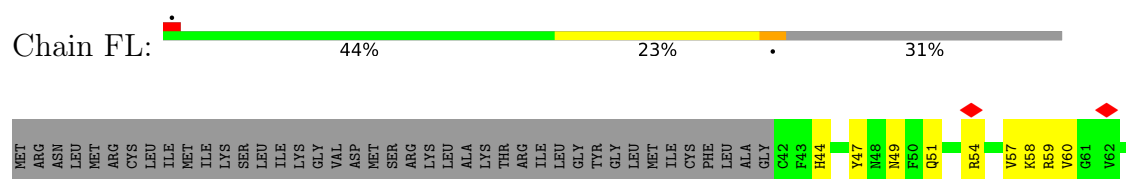
- Molecule 3: Inner membrane lipoprotein YiaD

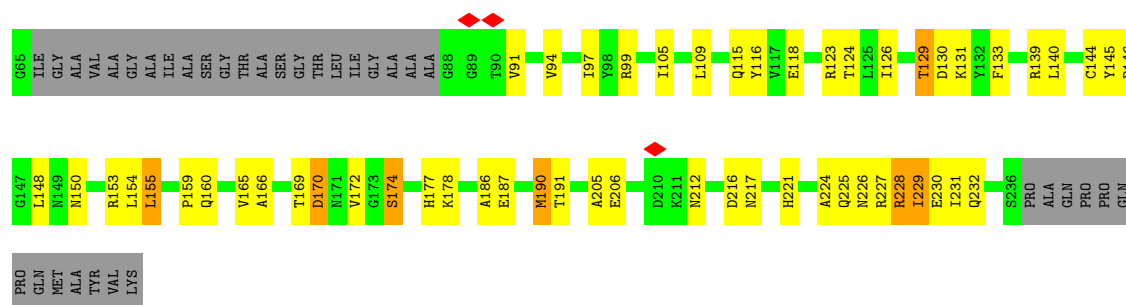


- Molecule 4: Outer membrane protein, OmpA family protein



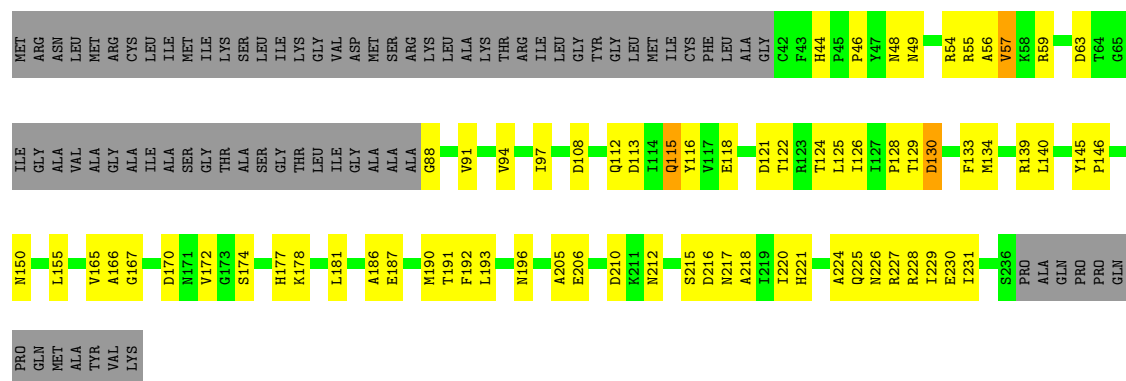
- Molecule 4: Outer membrane protein, OmpA family protein





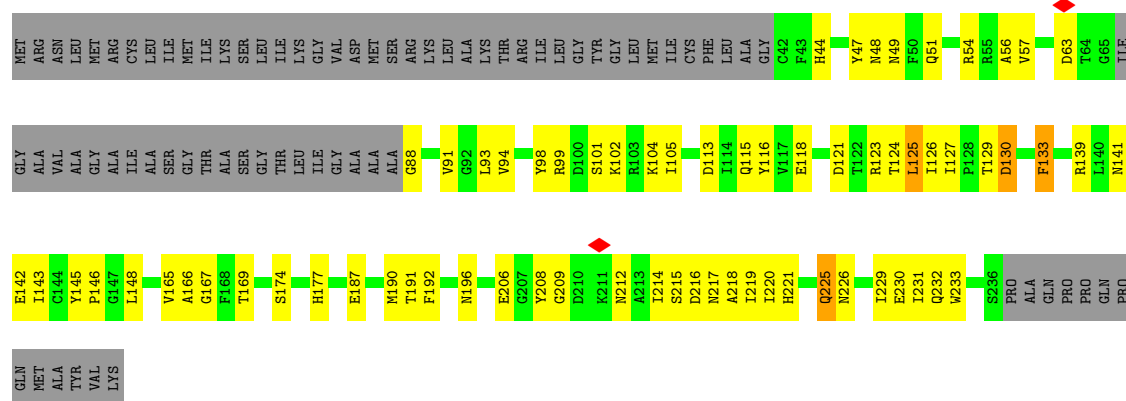
- Molecule 4: Outer membrane protein, OmpA family protein

Chain GL: 41% 27% • 31%



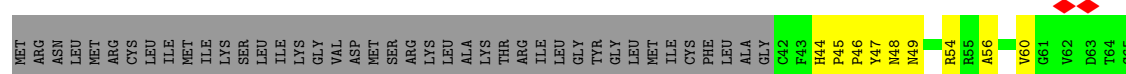
- Molecule 4: Outer membrane protein, OmpA family protein

Chain HL: 42% 26% • 31%

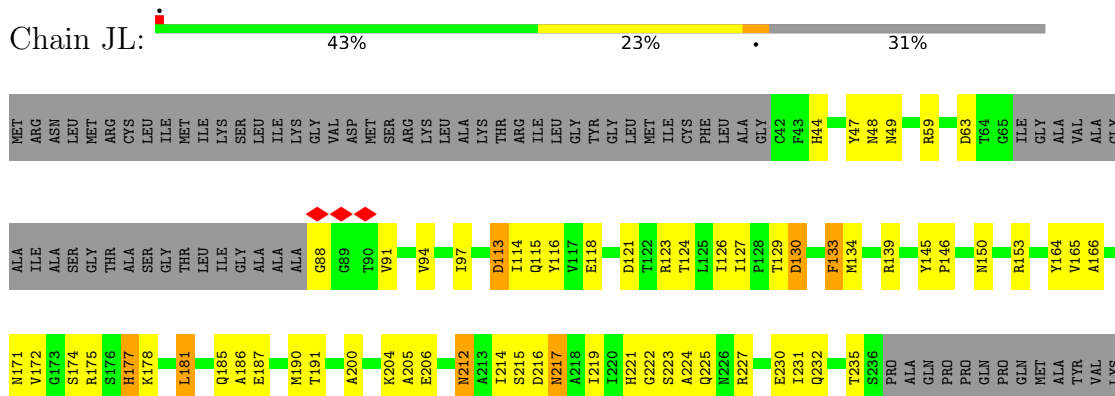


- Molecule 4: Outer membrane protein, OmpA family protein

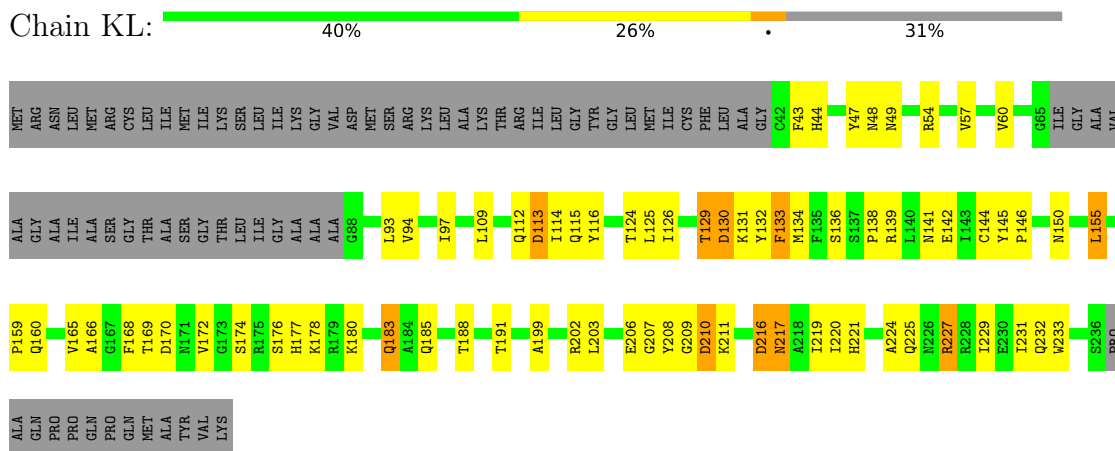
Chain IL: 43% 24% • 31%



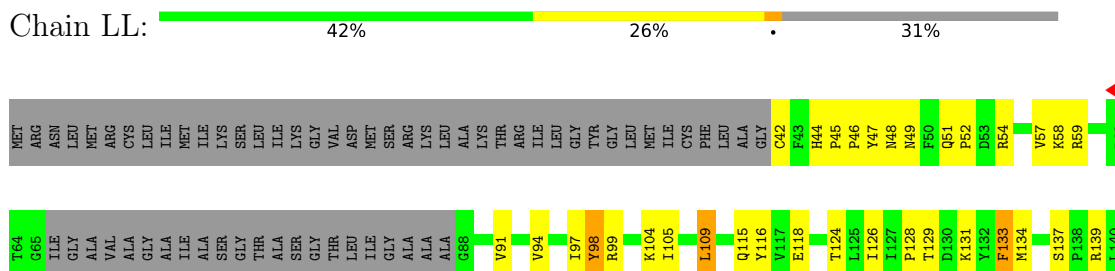
- Molecule 4: Outer membrane protein, OmpA family protein



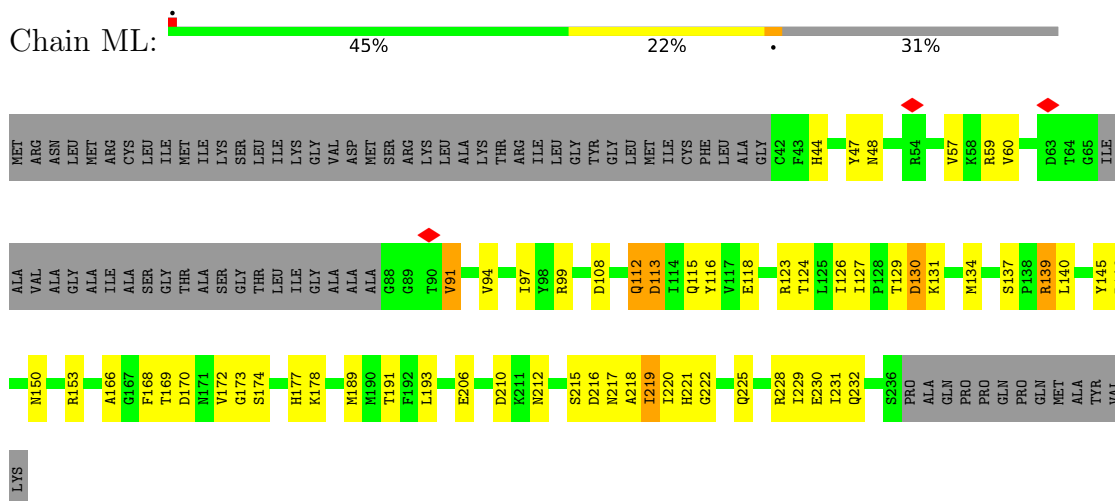
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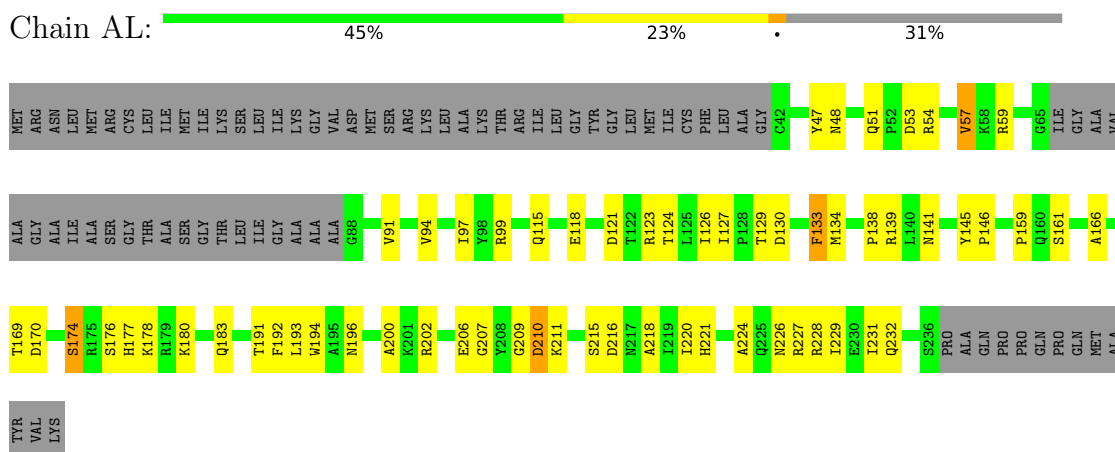
- Molecule 4: Outer membrane protein, OmpA family protein



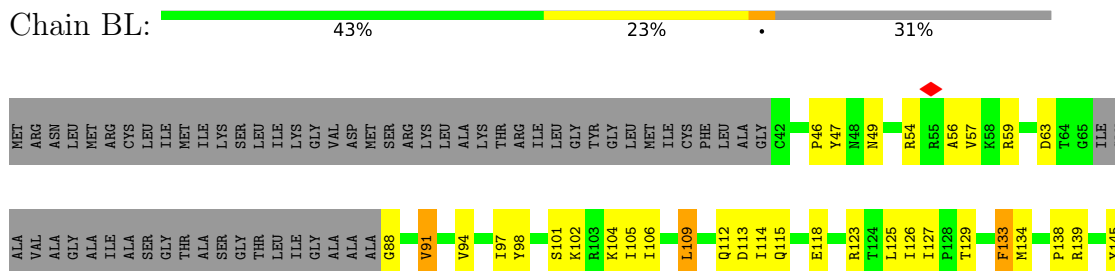
- Molecule 4: Outer membrane protein, OmpA family protein

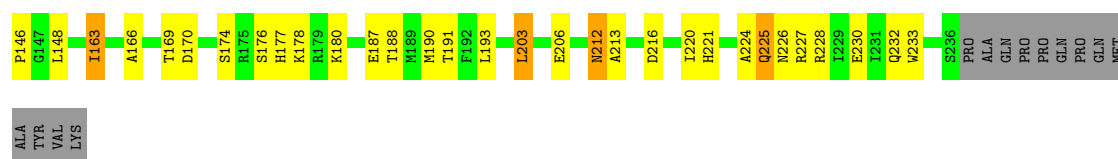


- Molecule 4: Outer membrane protein, OmpA family protein

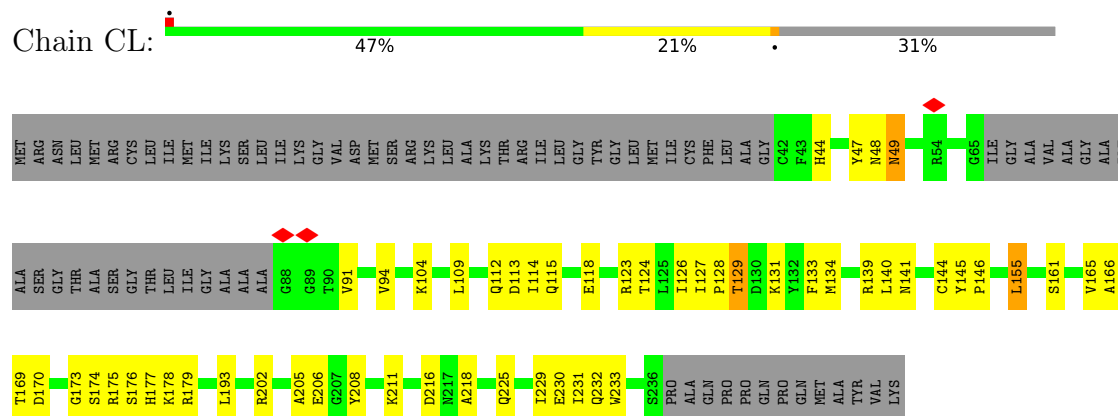


- Molecule 4: Outer membrane protein, OmpA family protein

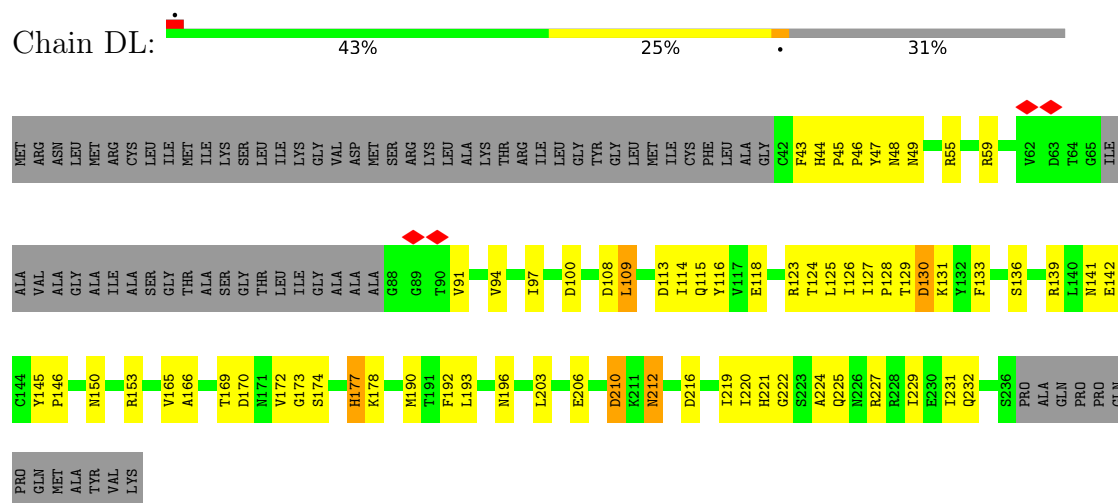




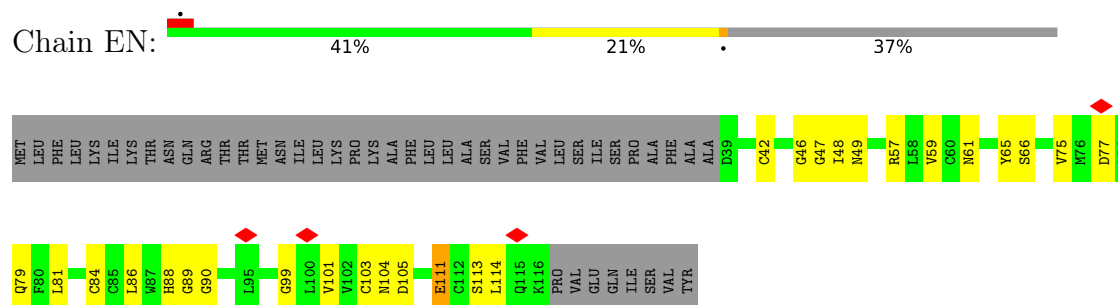
- Molecule 4: Outer membrane protein, OmpA family protein



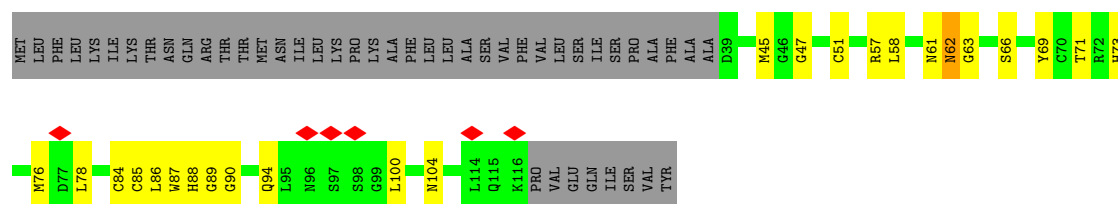
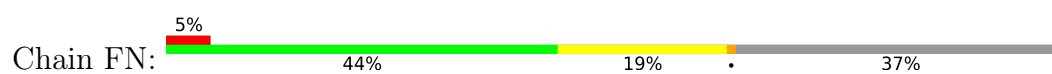
- Molecule 4: Outer membrane protein, OmpA family protein



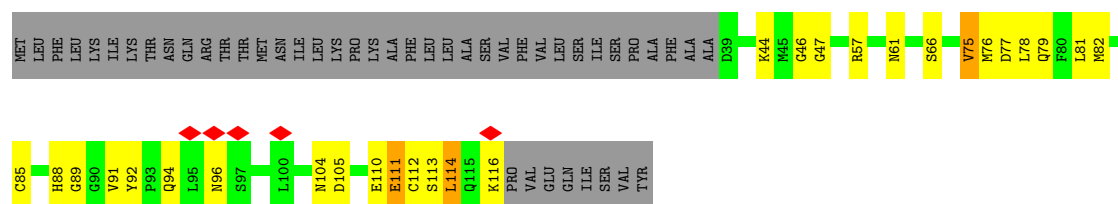
- Molecule 5: Neurogenic locus notch



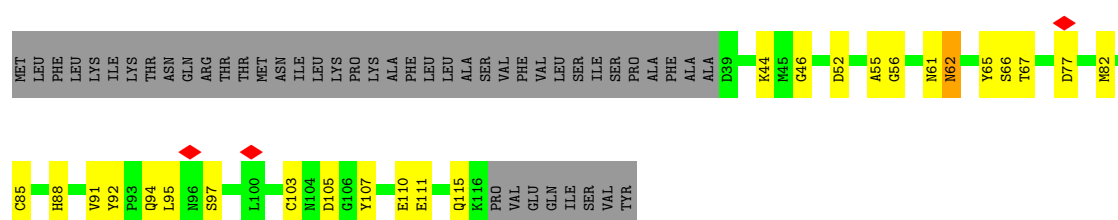
- Molecule 5: Neurogenic locus notch



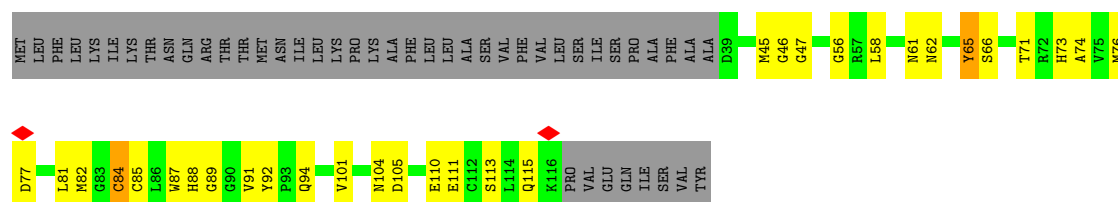
- Molecule 5: Neurogenic locus notch



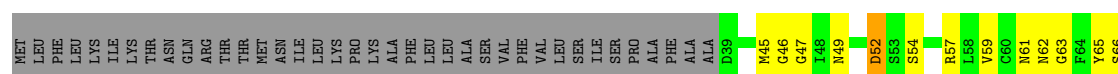
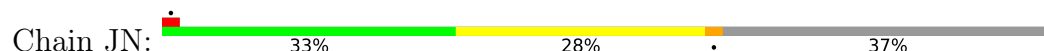
- Molecule 5: Neurogenic locus notch

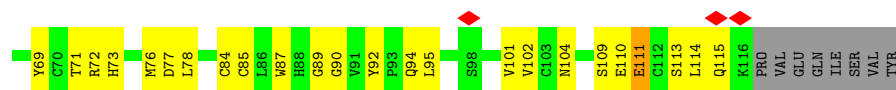


- Molecule 5: Neurogenic locus notch

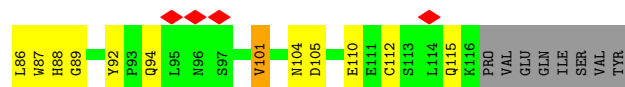
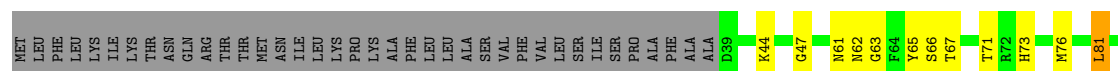


- Molecule 5: Neurogenic locus notch

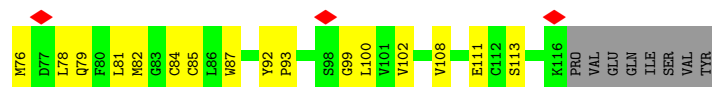
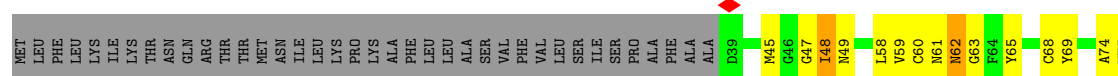




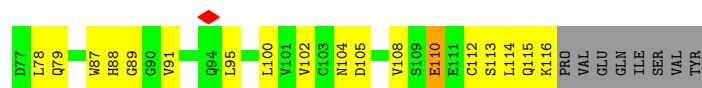
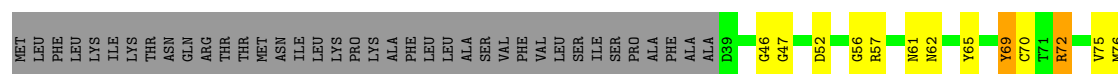
- Molecule 5: Neurogenic locus notch



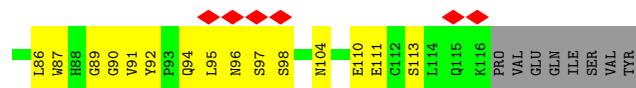
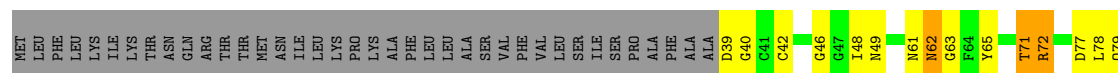
- Molecule 5: Neurogenic locus notch



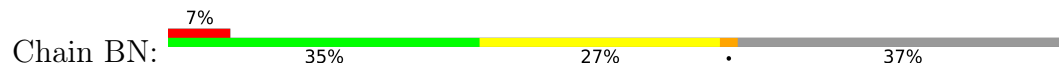
- Molecule 5: Neurogenic locus notch

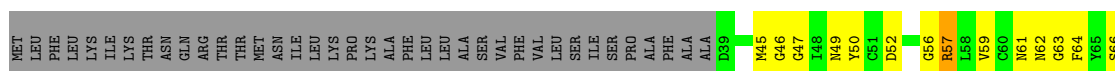


- Molecule 5: Neurogenic locus notch



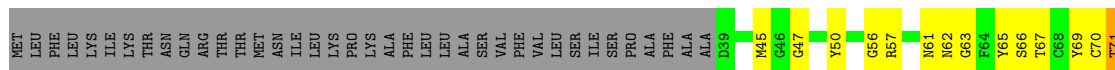
- Molecule 5: Neurogenic locus notch





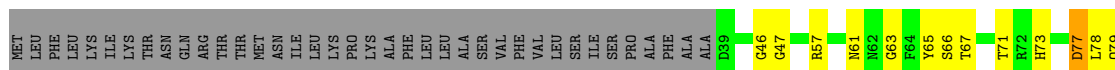
- Molecule 5: Neurogenic locus notch

Chain CN: 36% 26% 37%



- Molecule 5: Neurogenic locus notch

Chain DN: 44% 18% 37%



- Molecule 6: Unknown protein fragment

Chain EU: 89% 11%



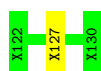
- Molecule 6: Unknown protein fragment

Chain FU: 78% 22%



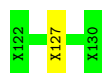
- Molecule 6: Unknown protein fragment

Chain GU: 89% 11%



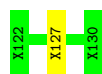
- Molecule 6: Unknown protein fragment

Chain HU:  89% 11%



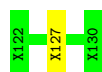
- Molecule 6: Unknown protein fragment

Chain IU:  89% 11%



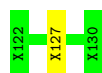
- Molecule 6: Unknown protein fragment

Chain JU:  89% 11%



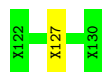
- Molecule 6: Unknown protein fragment

Chain KU:  89% 11%

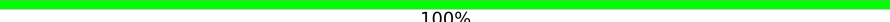


- Molecule 6: Unknown protein fragment

Chain LU:  89% 11%



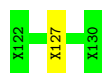
- Molecule 6: Unknown protein fragment

Chain MU:  100%

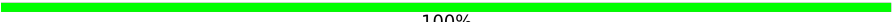
There are no outlier residues recorded for this chain.

- Molecule 6: Unknown protein fragment

Chain AU:  89% 11%

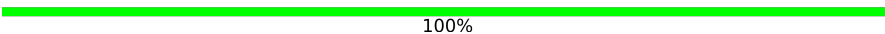


- Molecule 6: Unknown protein fragment

Chain BU:  100%

There are no outlier residues recorded for this chain.

- Molecule 6: Unknown protein fragment

Chain CU:  100%

There are no outlier residues recorded for this chain.

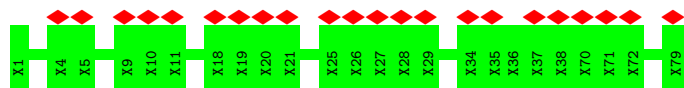
- Molecule 6: Unknown protein fragment

Chain DU:  89% 11%



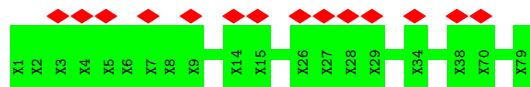
- Molecule 7: Unknown protein fragment

Chain EX:  46% 100%



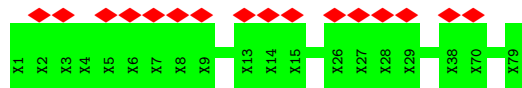
- Molecule 7: Unknown protein fragment

Chain FX:  29% 100%



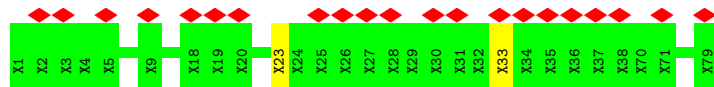
- Molecule 7: Unknown protein fragment

Chain GX:  33% 100%



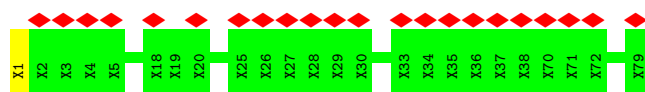
- Molecule 7: Unknown protein fragment

Chain HX:  44% 96%

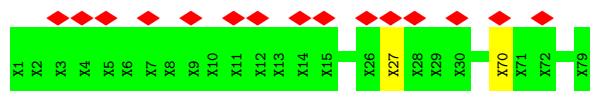


- Molecule 7: Unknown protein fragment

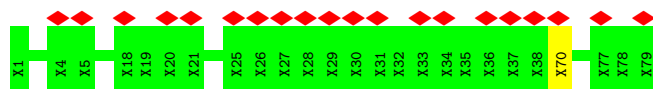
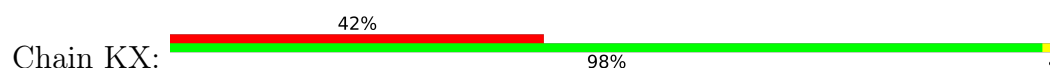
Chain IX:  46% 98%



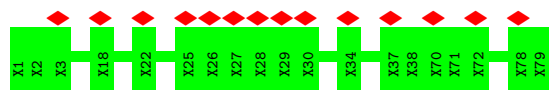
- Molecule 7: Unknown protein fragment



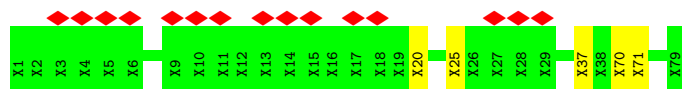
- Molecule 7: Unknown protein fragment



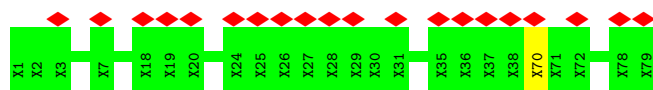
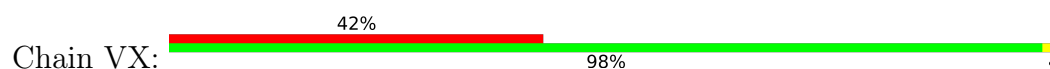
- Molecule 7: Unknown protein fragment



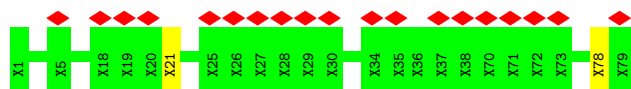
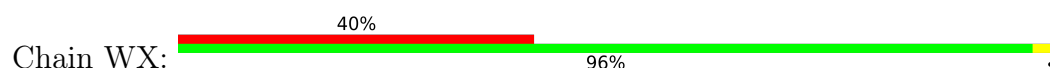
- Molecule 7: Unknown protein fragment



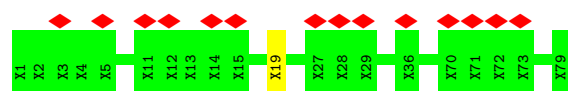
- Molecule 7: Unknown protein fragment



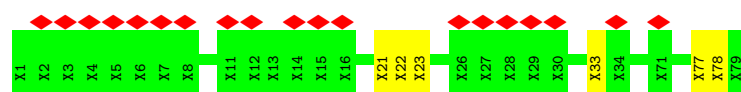
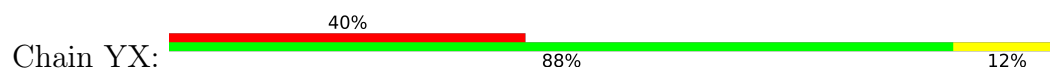
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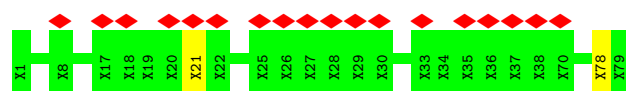
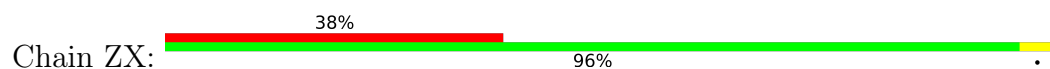
- Molecule 7: Unknown protein fragment



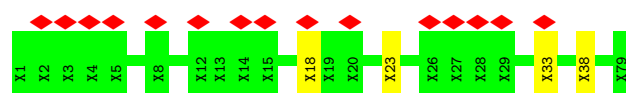
- Molecule 7: Unknown protein fragment



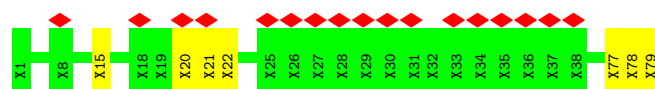
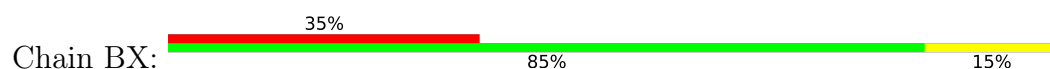
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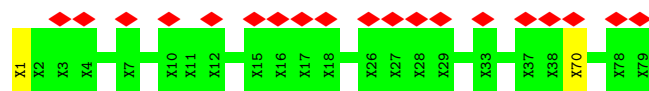
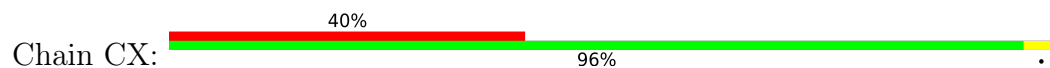
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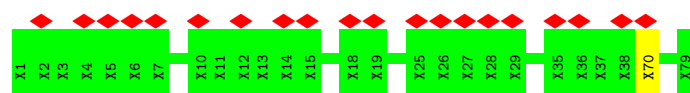
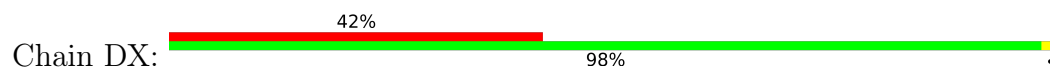
- Molecule 7: Unknown protein fragment



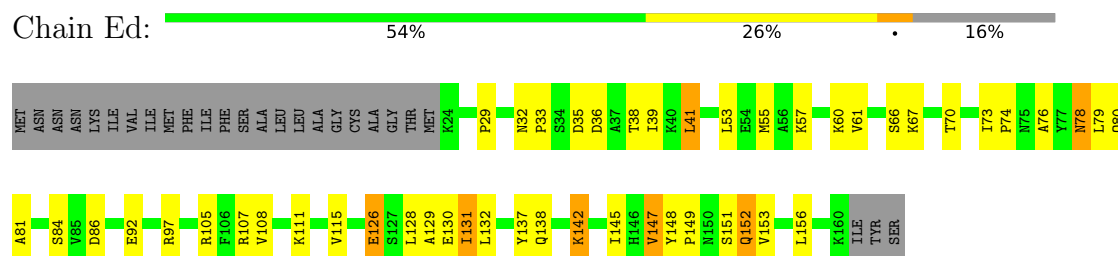
- Molecule 7: Unknown protein fragment



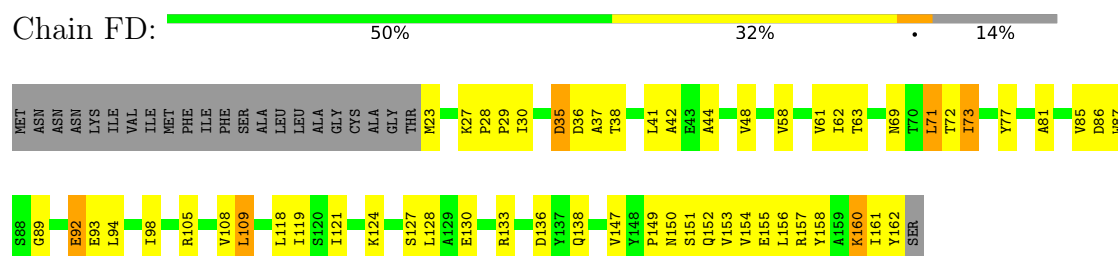
- Molecule 7: Unknown protein fragment



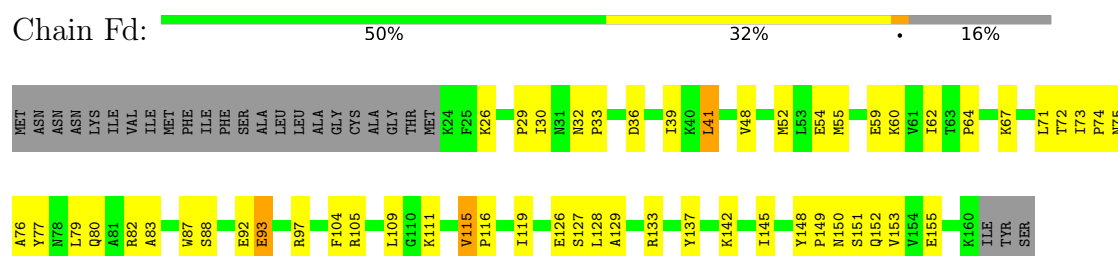
- Molecule 8: DotD



- Molecule 8: DotD



- Molecule 8: DotD



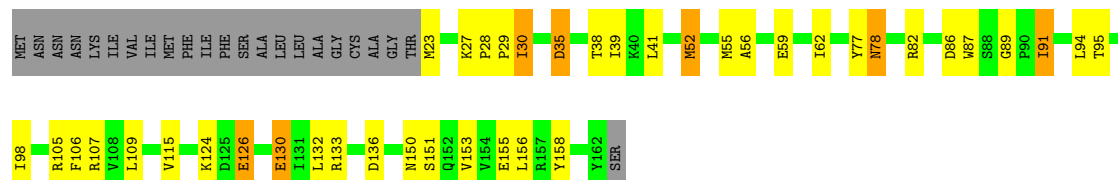
- Molecule 8: DotD



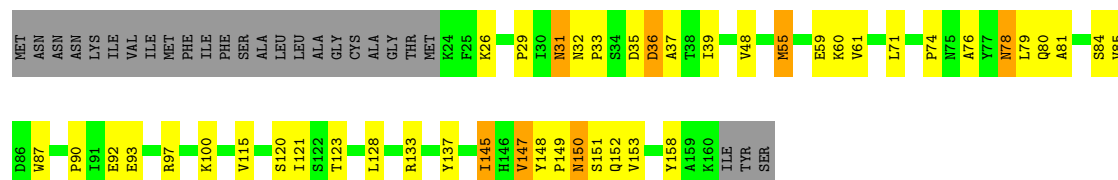
- Molecule 8: DotD



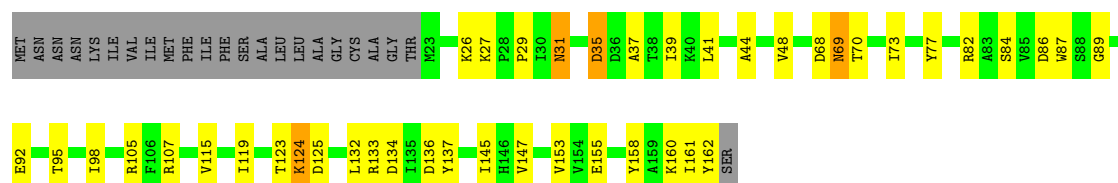
• Molecule 8: DotD

Chain HD:  61% 21% 14%

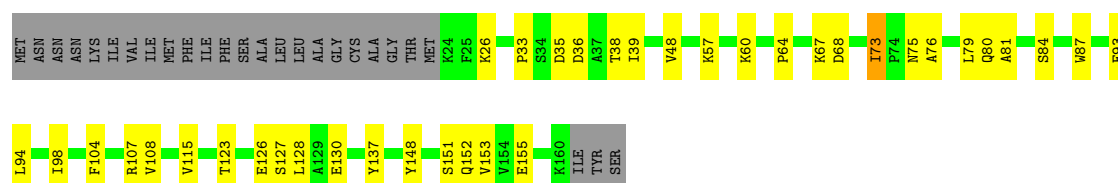
• Molecule 8: DotD

Chain Hd:  56% 23% 16%

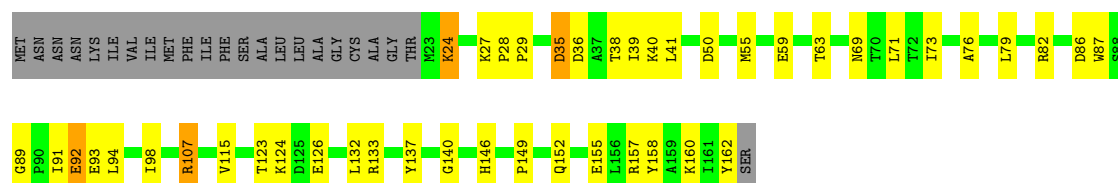
• Molecule 8: DotD

Chain ID:  60% 24% 14%

• Molecule 8: DotD

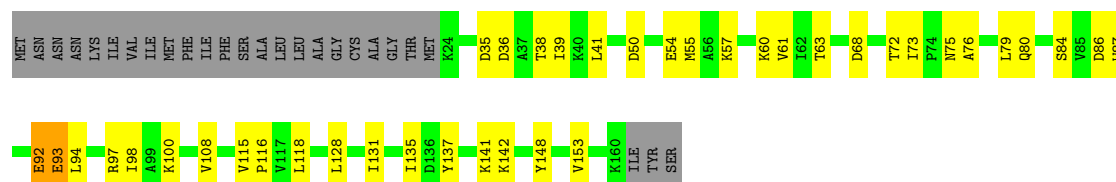
Chain Id:  61% 23% 16%

• Molecule 8: DotD

Chain JD:  58% 25% 14%

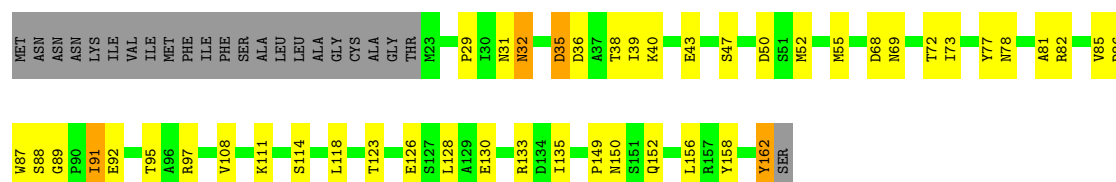
- Molecule 8: DotD

Chain Jd: 



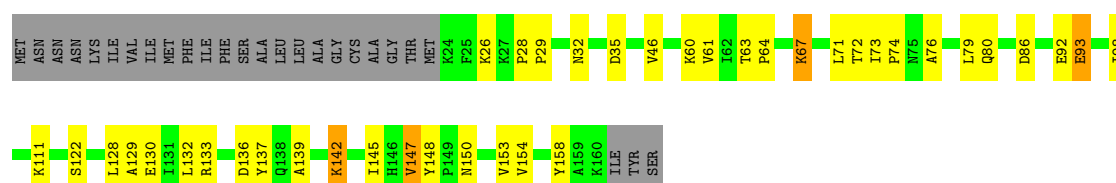
- Molecule 8: DotD

Chain KD: 



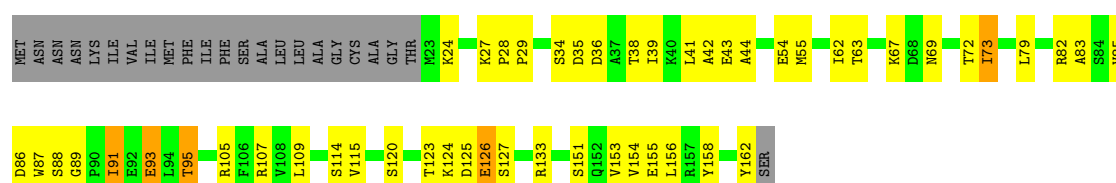
- Molecule 8: DotD

Chain Kd: 



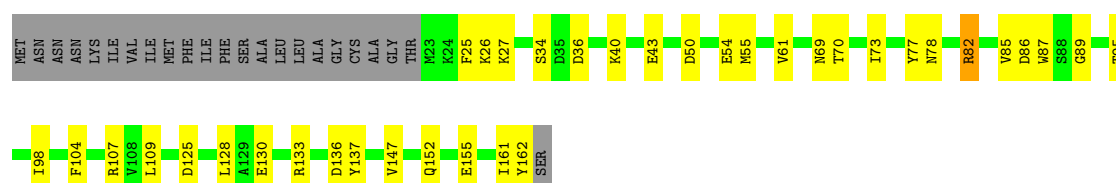
- Molecule 8: DotD

Chain LD: 



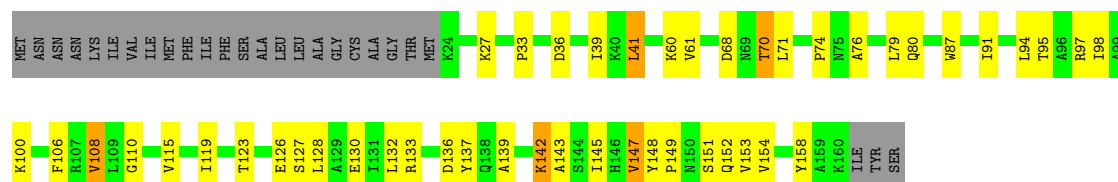
- Molecule 8: DotD

Chain CD: 



- Molecule 8: DotD

Chain Ld: 



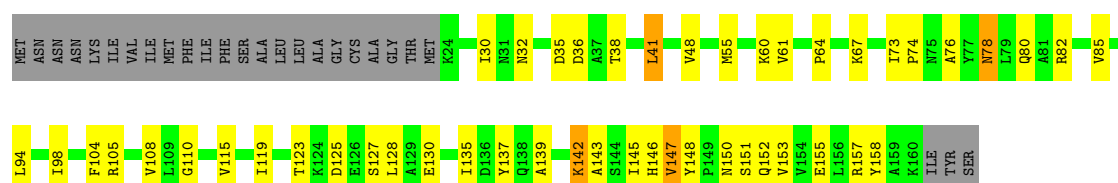
- Molecule 8: DotD

Chain MD: 



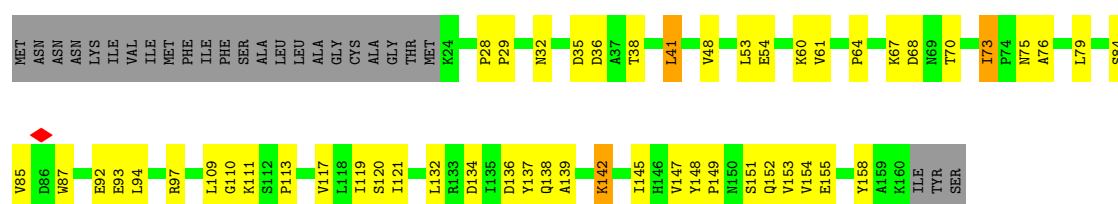
- Molecule 8: DotD

Chain Md: 



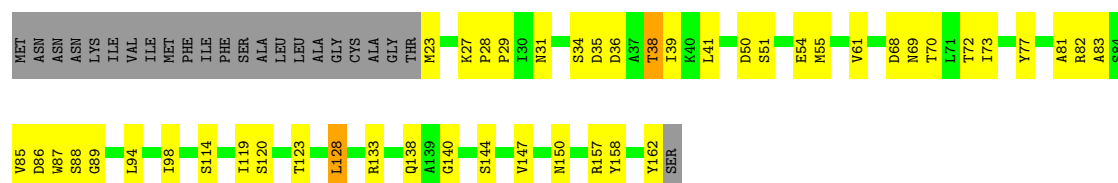
- Molecule 8: DotD

Chain Ad: 



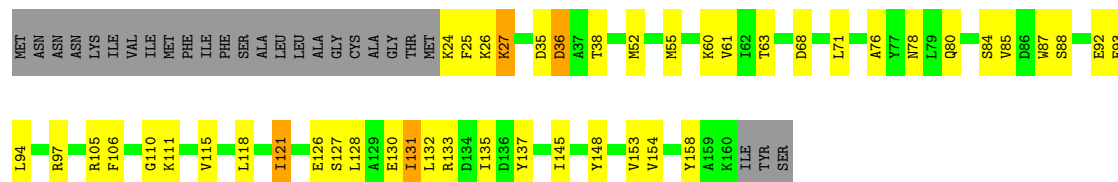
- Molecule 8: DotD

Chain BD: 



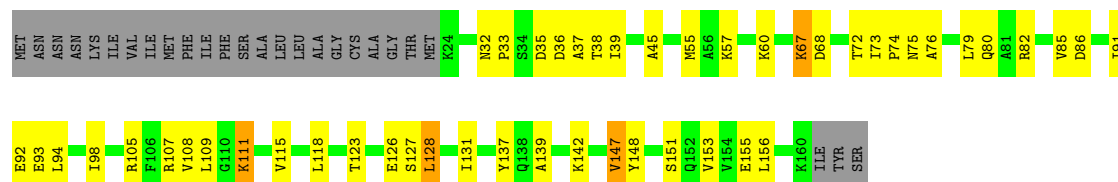
- Molecule 8: DotD

Chain Bd: 



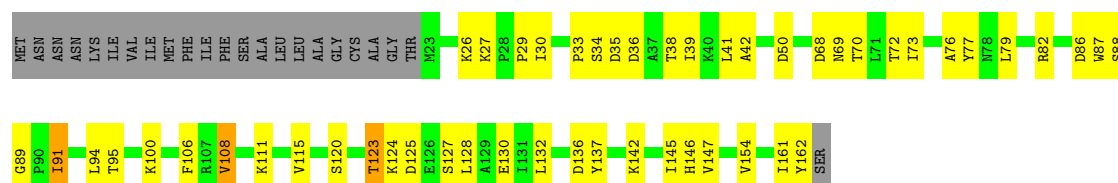
- Molecule 8: DotD

Chain Cd: 



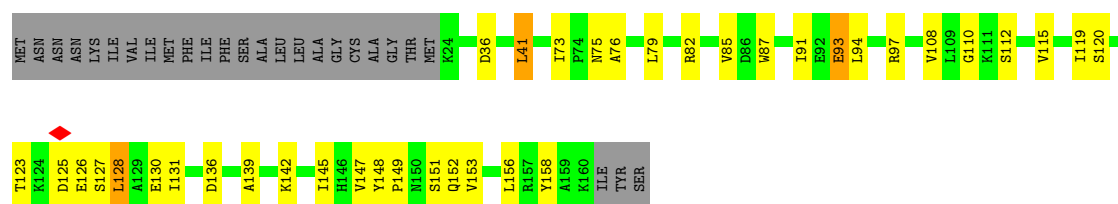
- Molecule 8: DotD

Chain DD: 



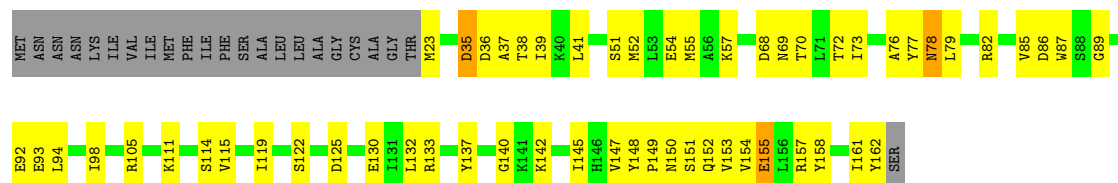
- Molecule 8: DotD

Chain Dd: 

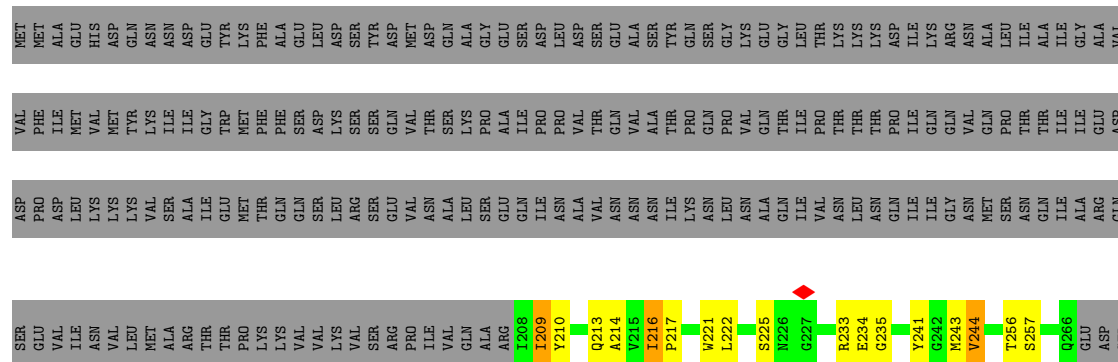
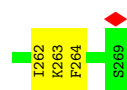
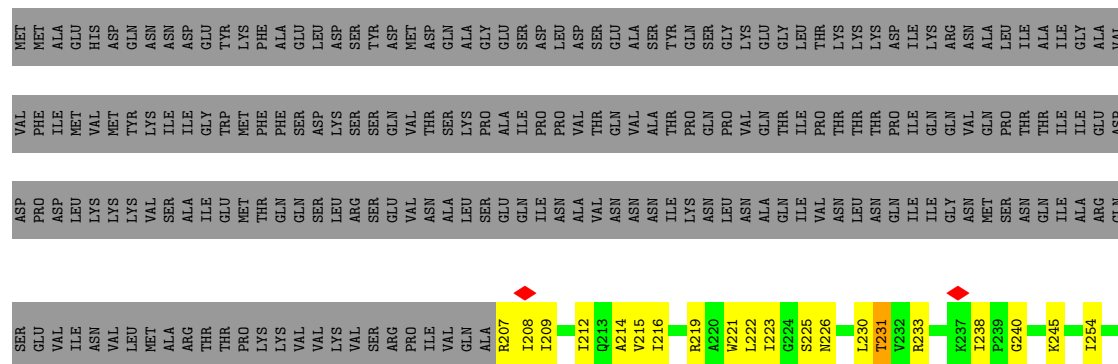
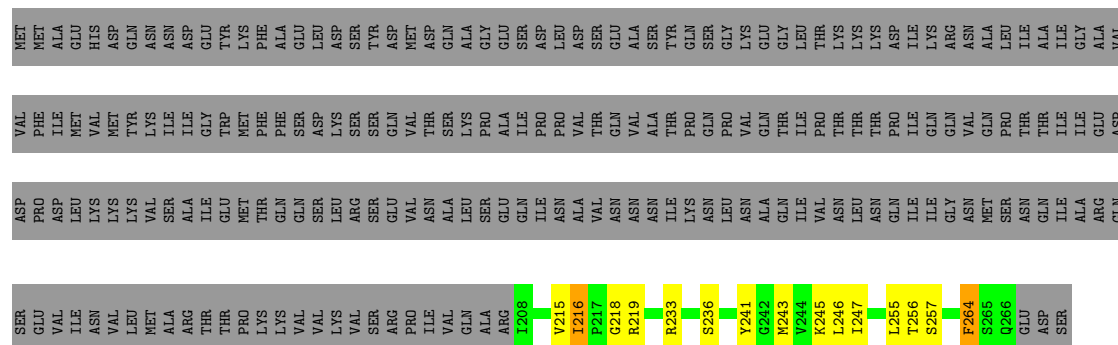


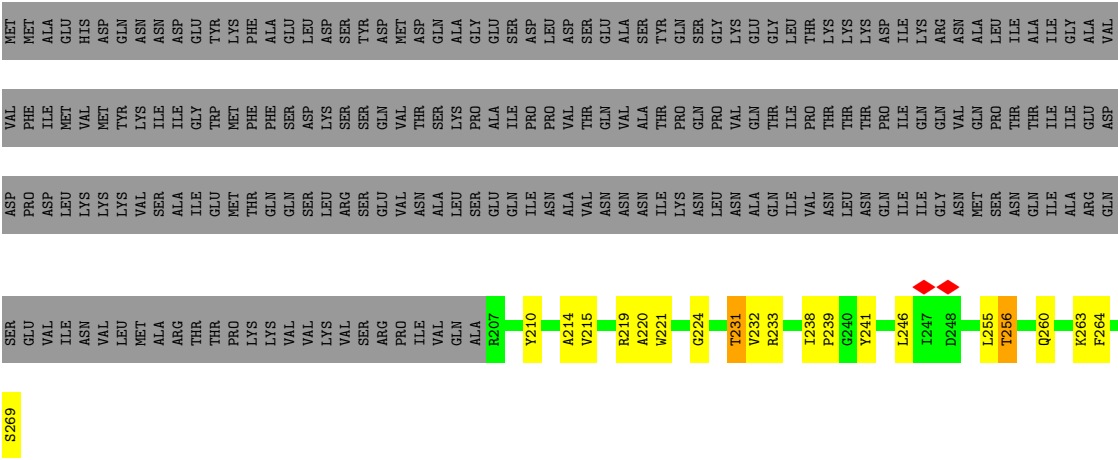
- Molecule 8: DotD

Chain AD: 

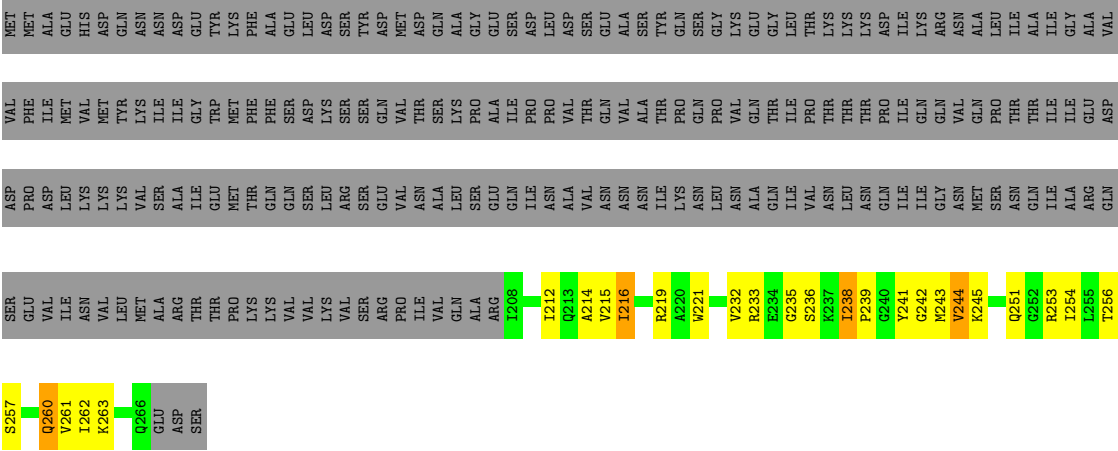




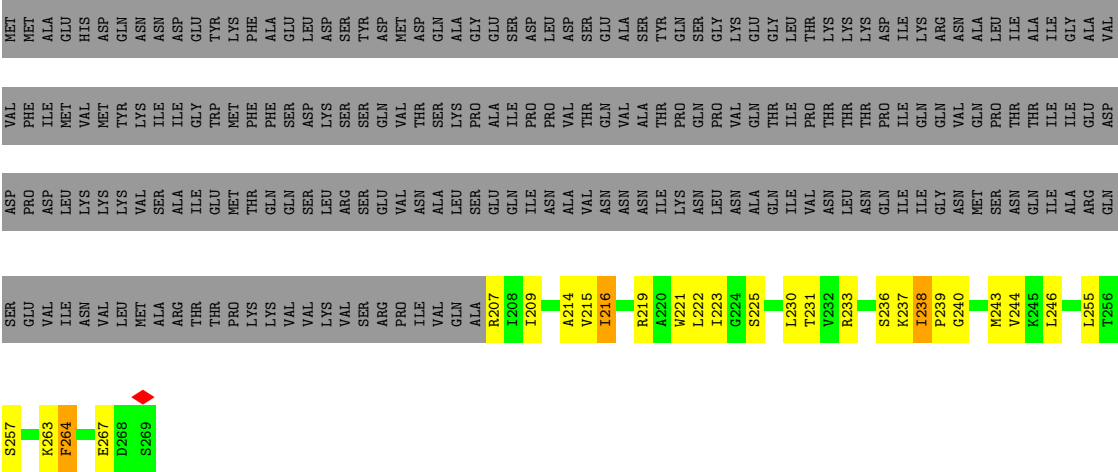




• Molecule 9: DotF



• Molecule 9: DotF

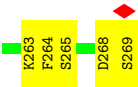


• Molecule 9: DotF

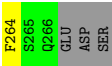
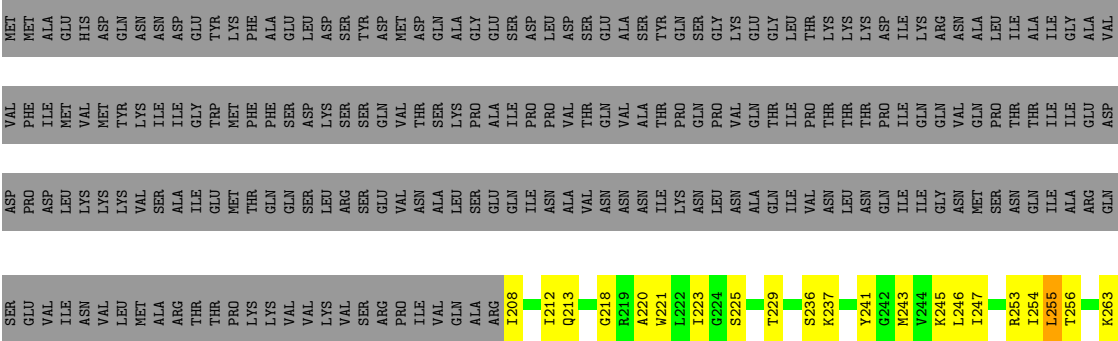
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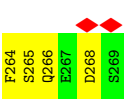
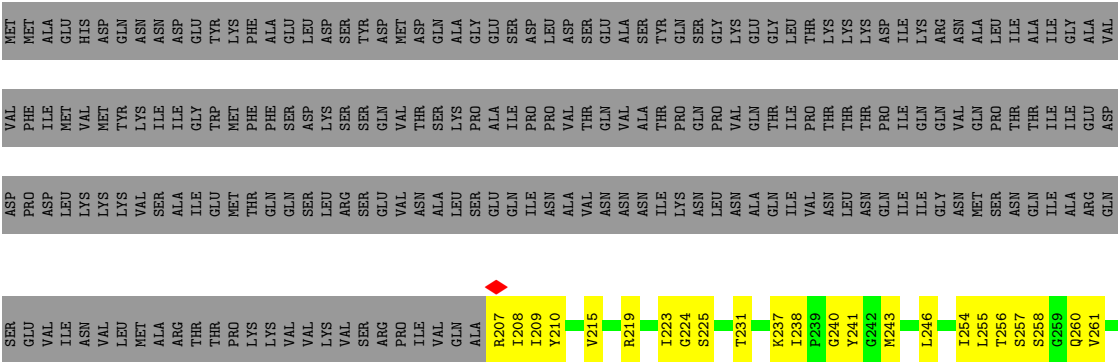
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S258	ASN	LYS	VAL
G259	VAL	LYS	MET
Q260	MET	VAL	TYR
S265	ALA	SER	ASN
Q266	ARG	ALA	ILE
GLU	THR	ILE	GLY
ASP	THR	GLU	TRP
SER	THR	GLU	TRP
	PRO	MET	LYS
	LYS	THR	PHE
	LYS	GLN	PHE
	VAL	GLN	ALA
	VAL	GLN	SER
	VAL	SER	ASP
	LYS	LEU	LYS
	VAL	ARG	ASP
	SER	SER	SER
	ARG	GLU	GLN
	PRO	VAL	VAL
	ILE	ASN	THR
	VAL	ALA	SER
	GLN	LEU	ALA
	ALA	SER	GLY
	ARG	GLU	ALA
	I208	GLN	ILE
	I209	ILE	PRO
	Y210	ASN	PRO
	Y211	ALA	VAL
		VAL	THR
	A214	ASN	GLN
	V215	ASN	ALA
	T216	ASN	ALA
	T217	ILE	THR
	G218	LYS	PRO
	R219	ASN	GLN
	A220	LEU	PRO
	W221	ASN	LYS
	L222	ALA	GLU
	L223	GLN	GLY
	G224	ILE	THR
	S225	VAL	ILE
		VAL	PRO
		ASN	THR
	T229	LEU	LYS
		ASN	THR
		GLN	THR
	V232	ILE	PRO
	R233	THR	ASP
	K237	ILE	ILE
	T238	GLN	LYS
		ASN	ARG
		MET	ASN
		GLN	ALA
	Y241	SER	LEU
	G242	ASN	ILE
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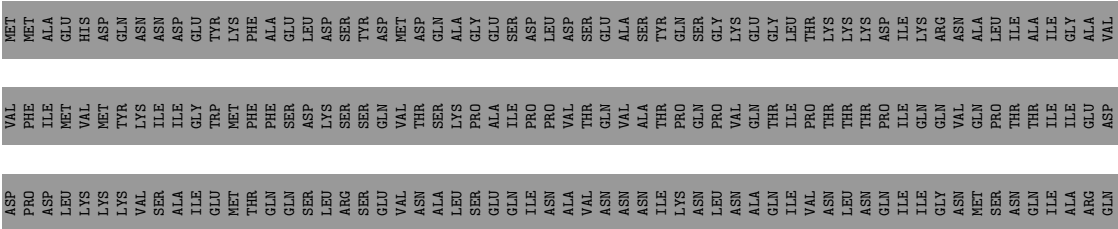
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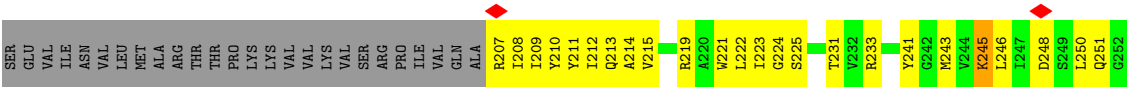


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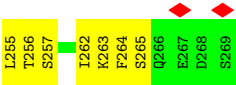
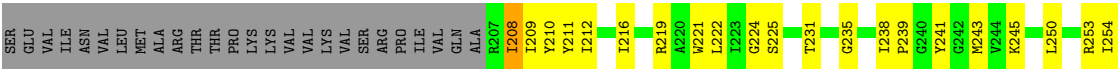
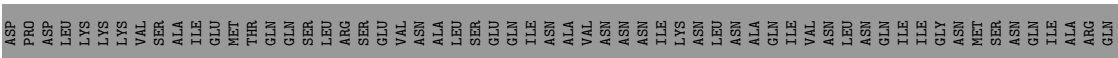
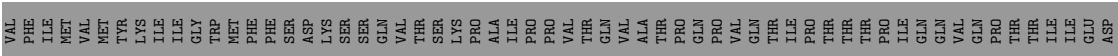
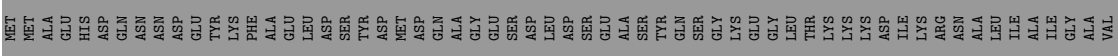


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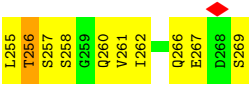
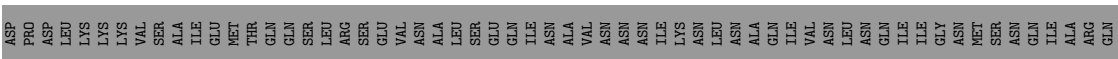
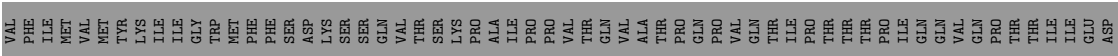
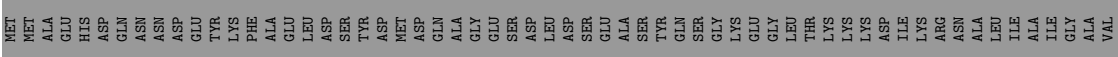




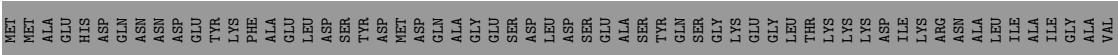
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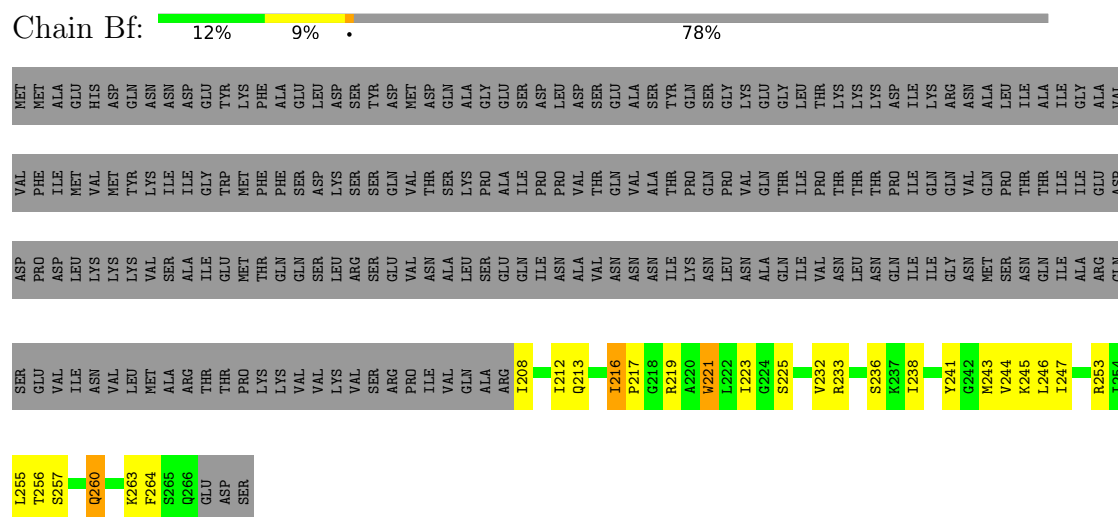
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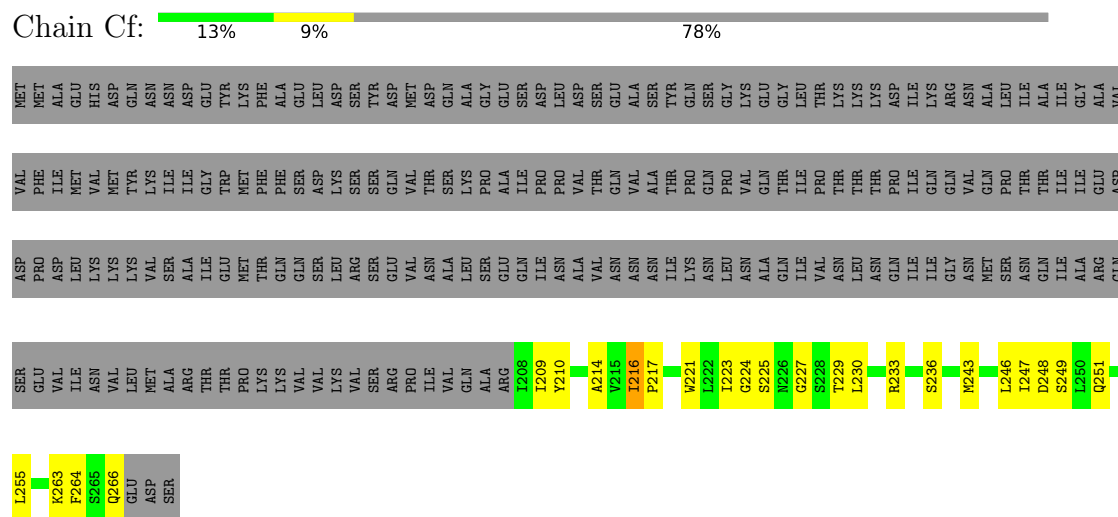
• Molecule 9: DotF



- Molecule 9: DotF



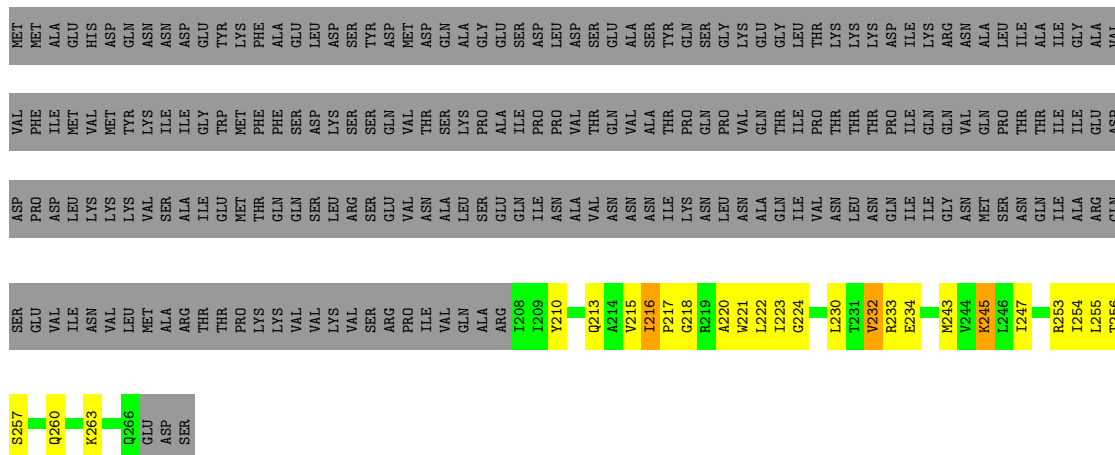
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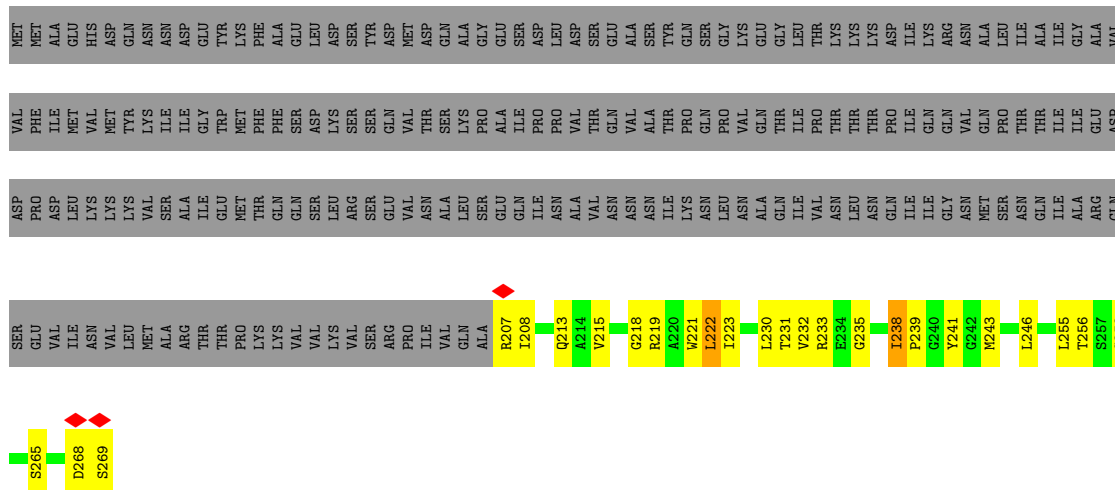
- Molecule 9: DotF



- Molecule 9: DotF



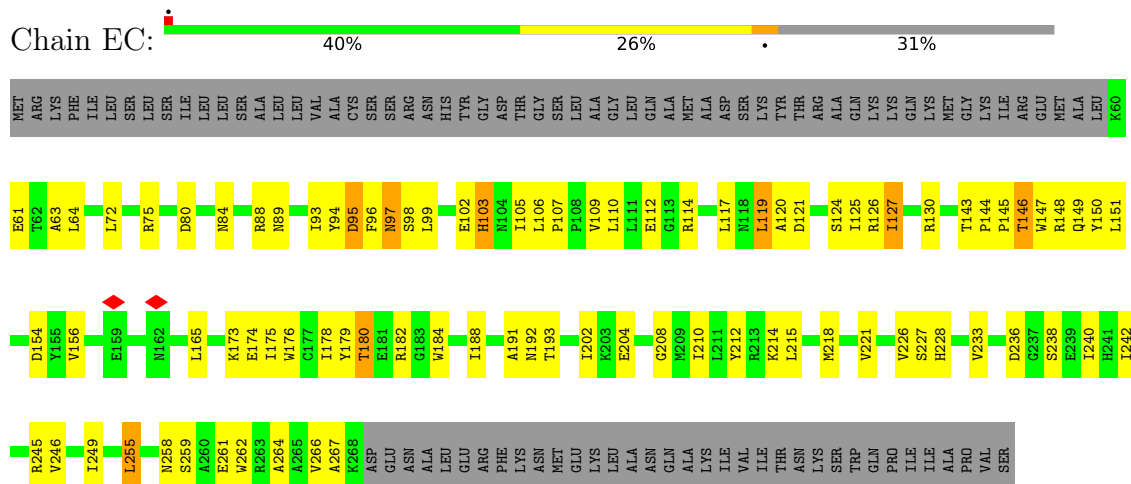
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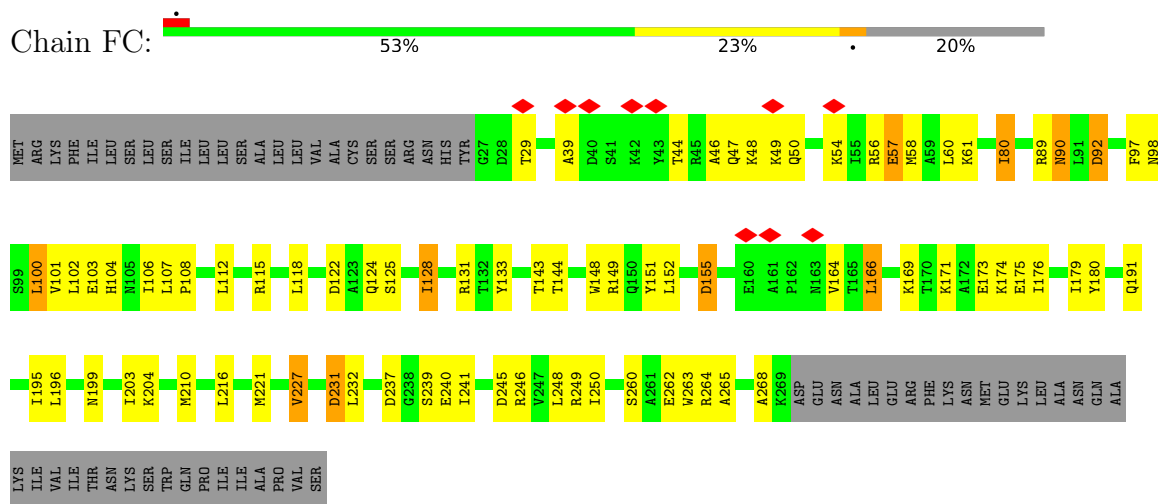
Chain DC:

Amino Acid	Percentage
MET	42%
ARG	25%
VAL	31%

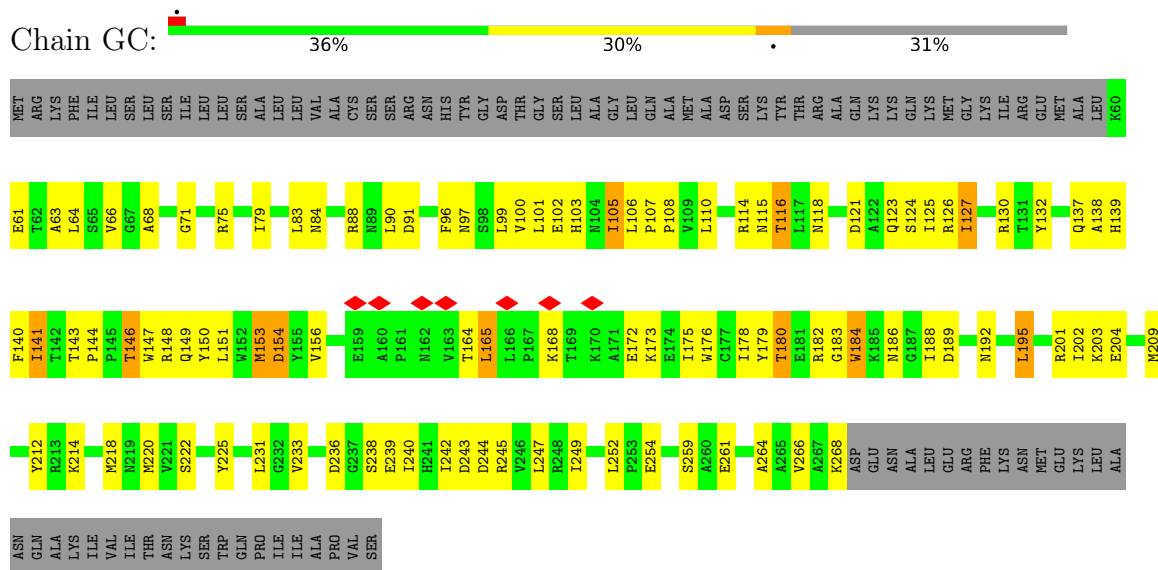
- Molecule 10: DotC



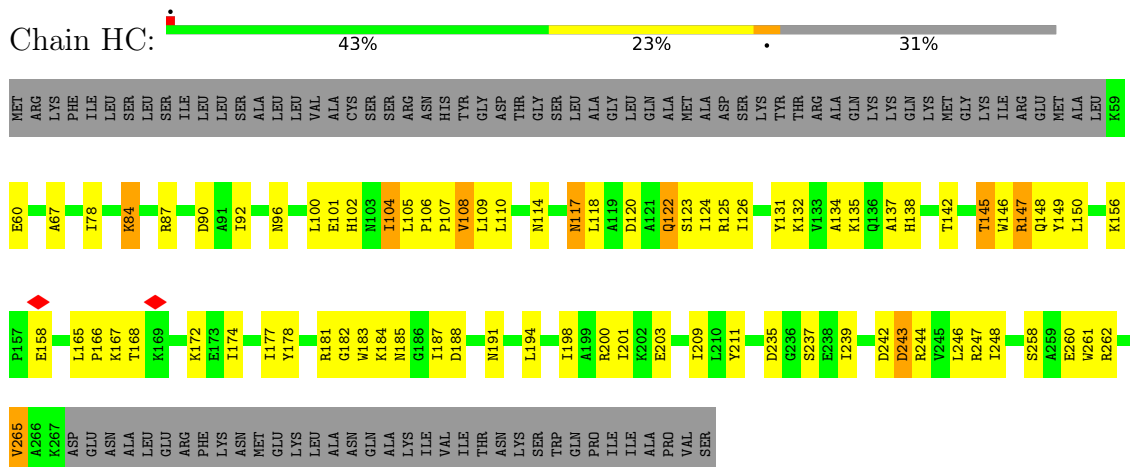
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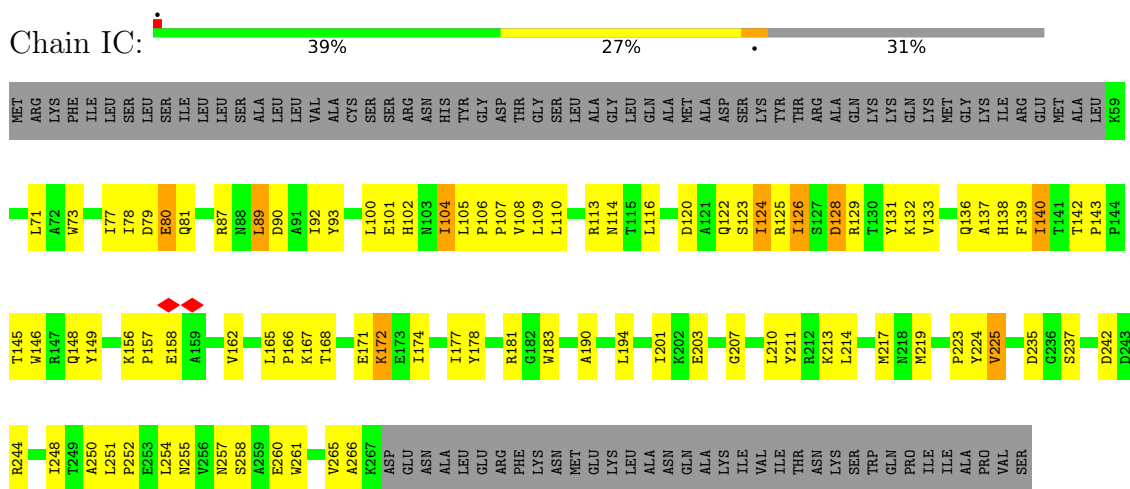
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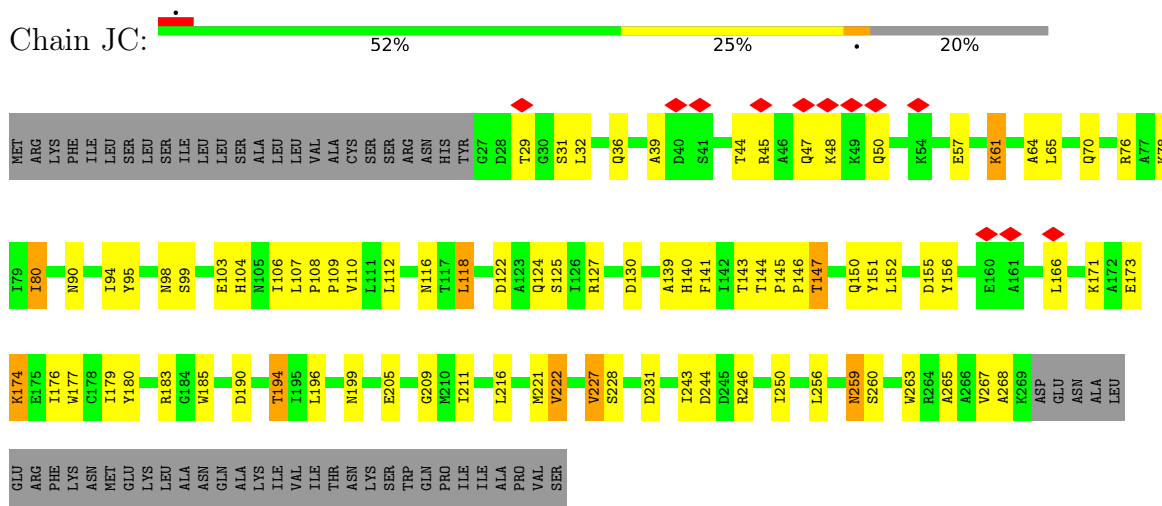
- Molecule 10: DotC



- Molecule 10: DotC

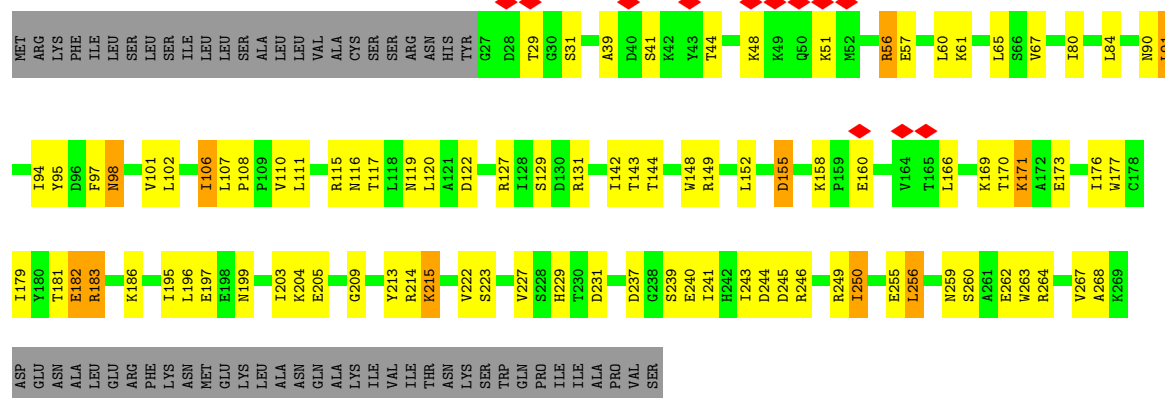


- Molecule 10: DotC

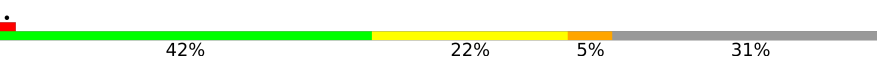


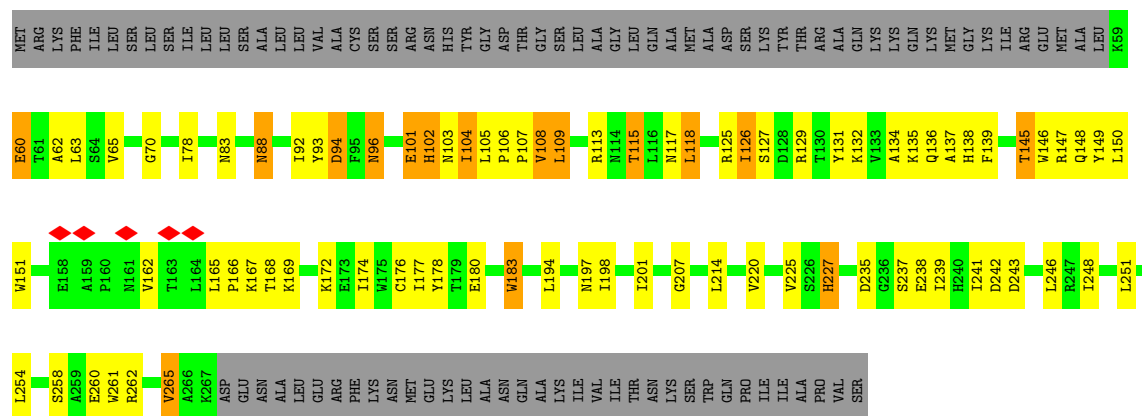
- Molecule 10: DotC

Chain KC: 

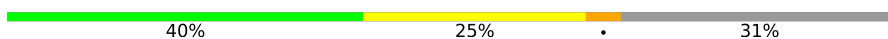


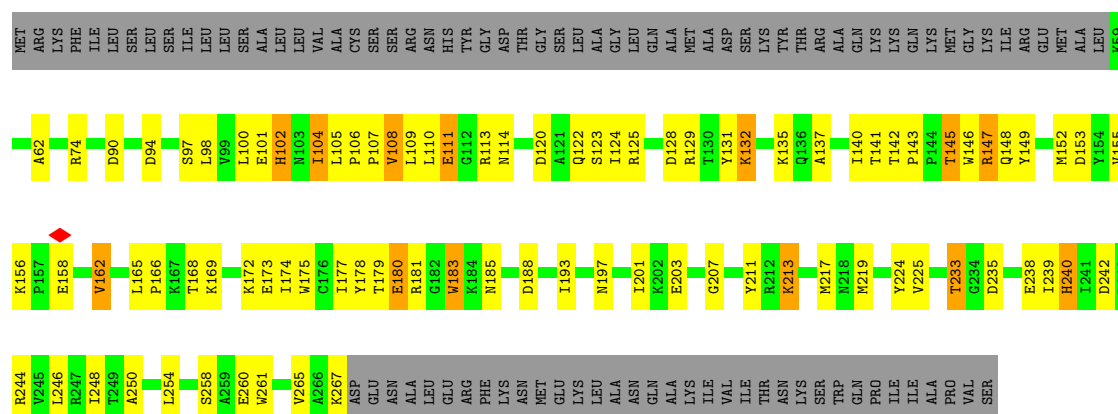
• Molecule 10: DotC

Chain LC: 



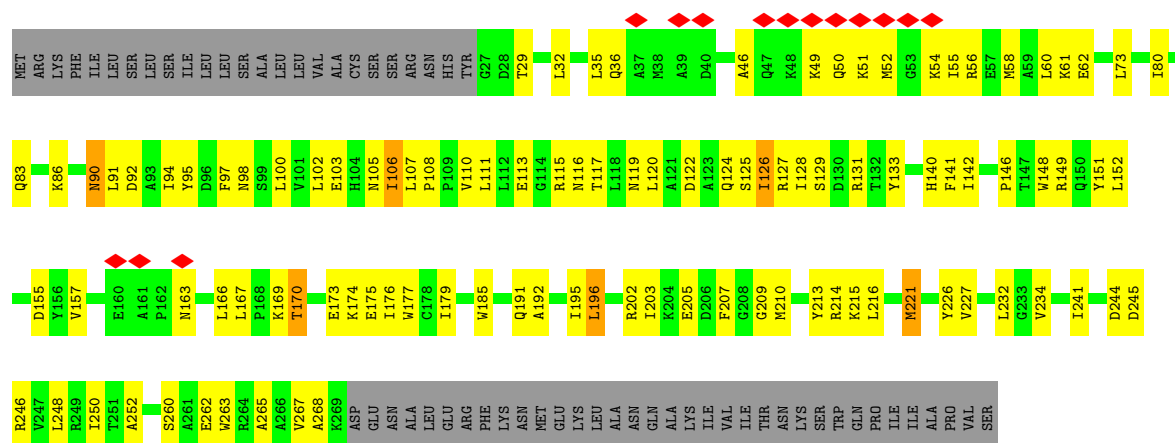
• Molecule 10: DotC

Chain MC: 

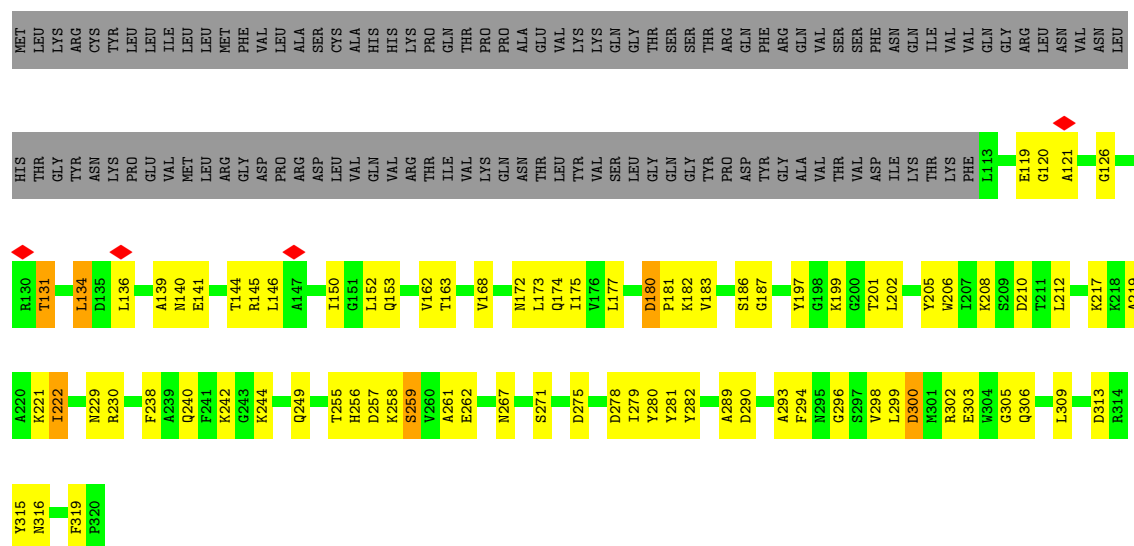


• Molecule 10: DotC

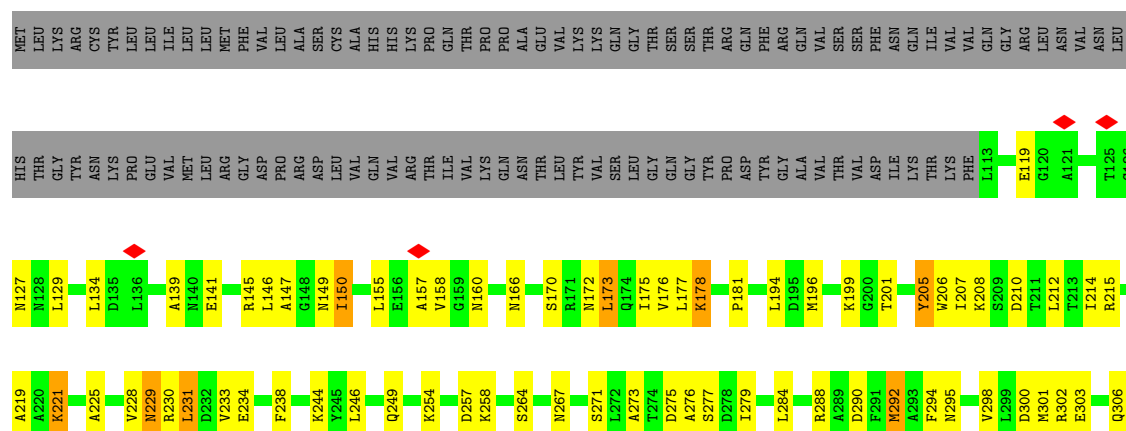
Chain AC: 



- Molecule 11: DUF2807 domain-containing protein

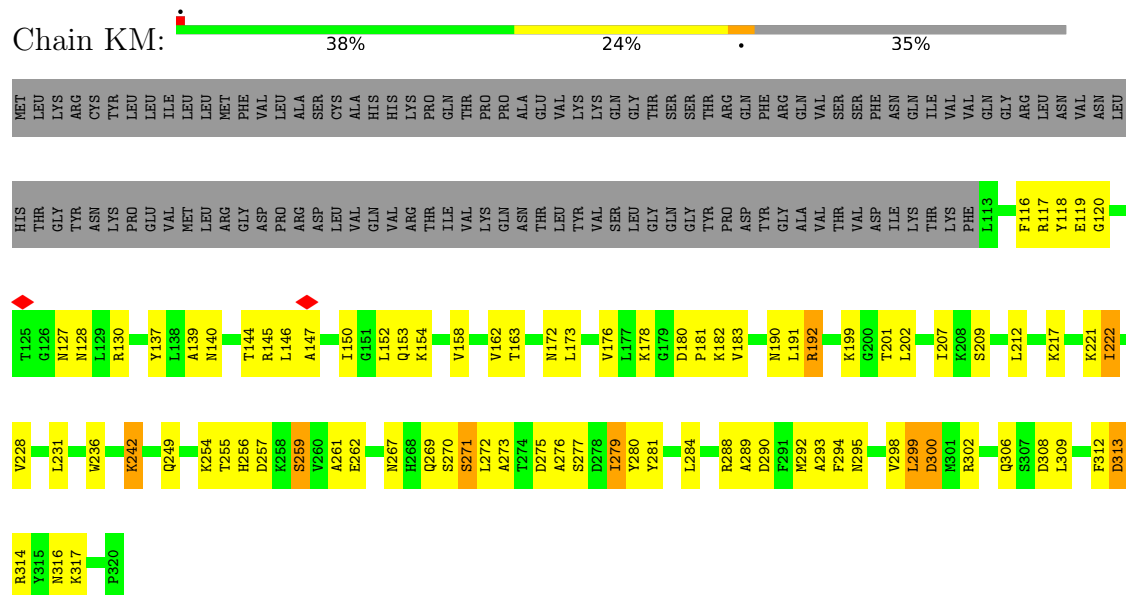


- Molecule 11: DUF2807 domain-containing protein

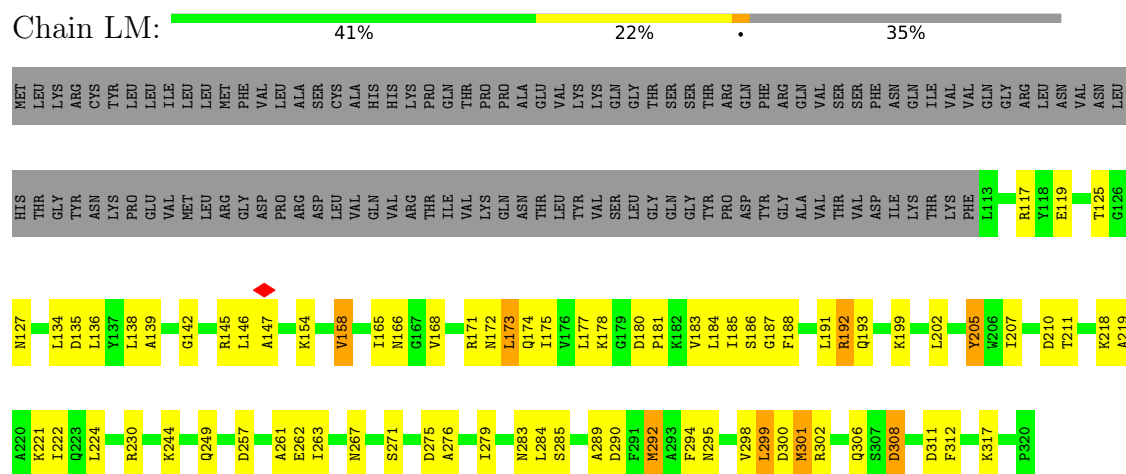




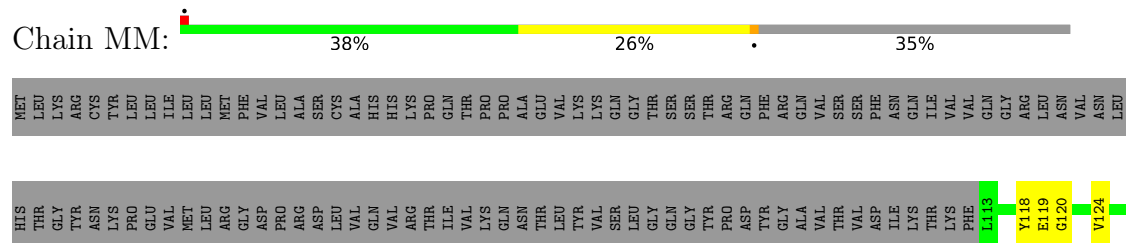
- Molecule 11: DUF2807 domain-containing protein

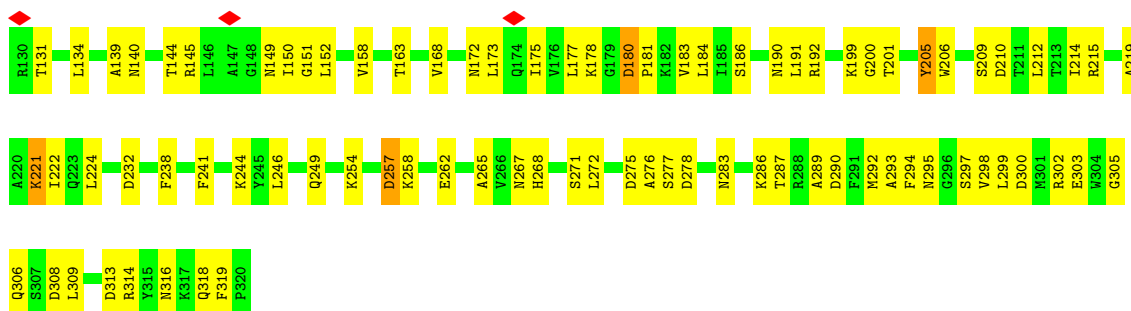


- Molecule 11: DUF2807 domain-containing protein

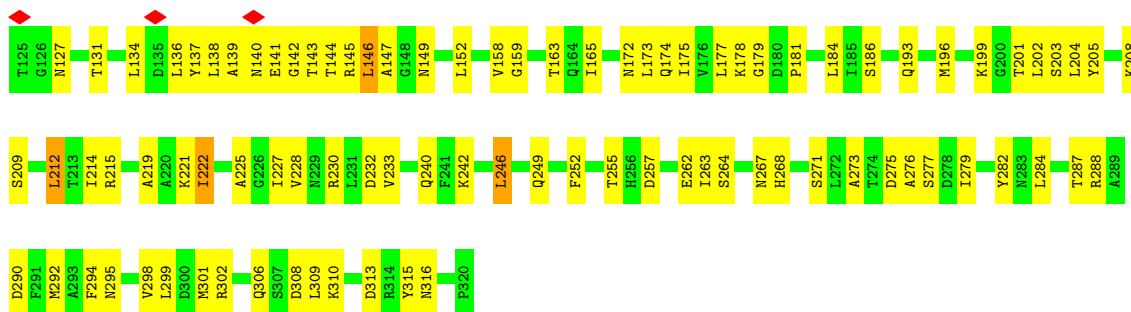
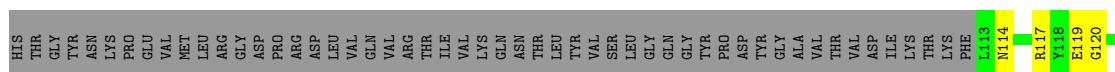
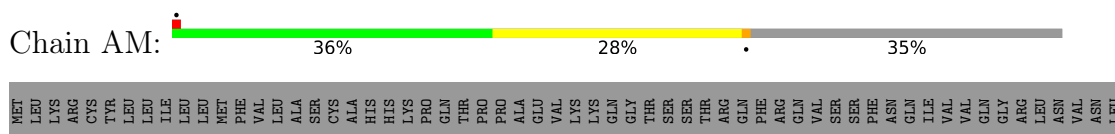


- Molecule 11: DUF2807 domain-containing protein

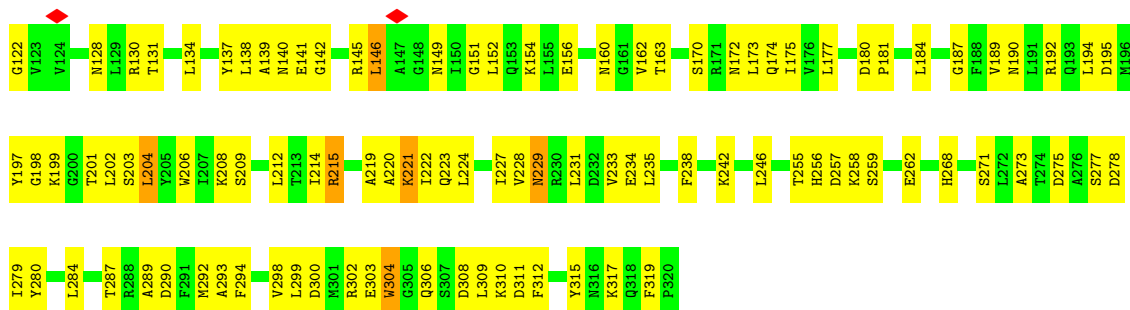
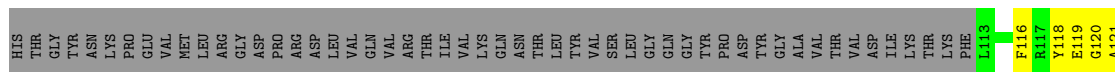
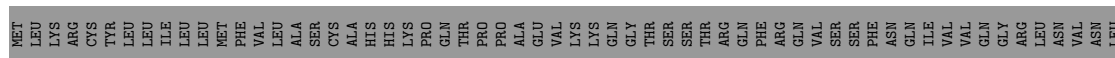




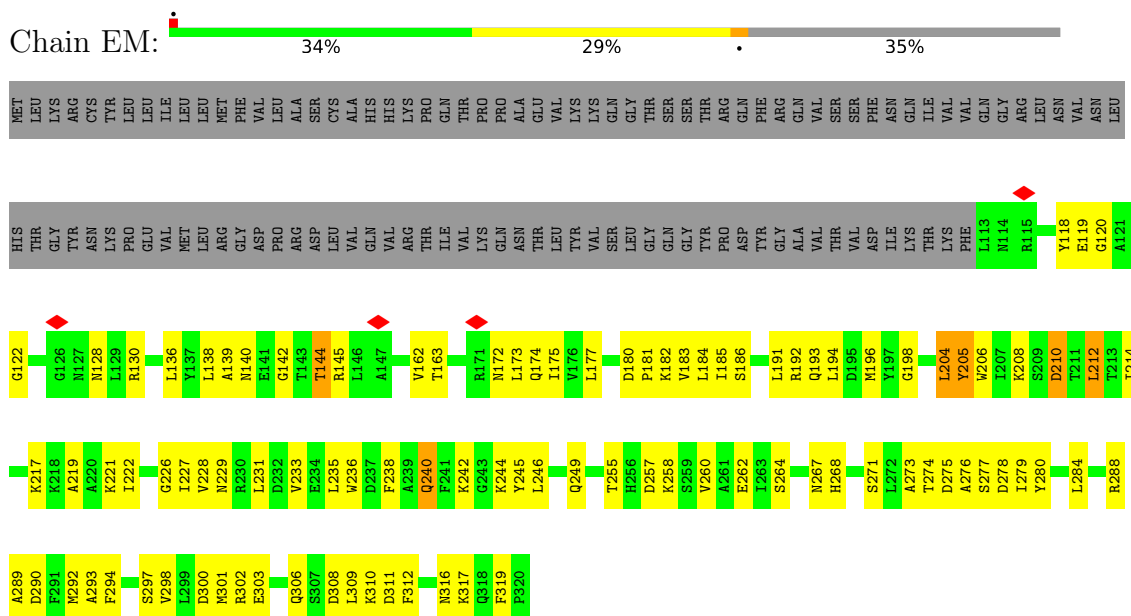
- Molecule 11: DUF2807 domain-containing protein



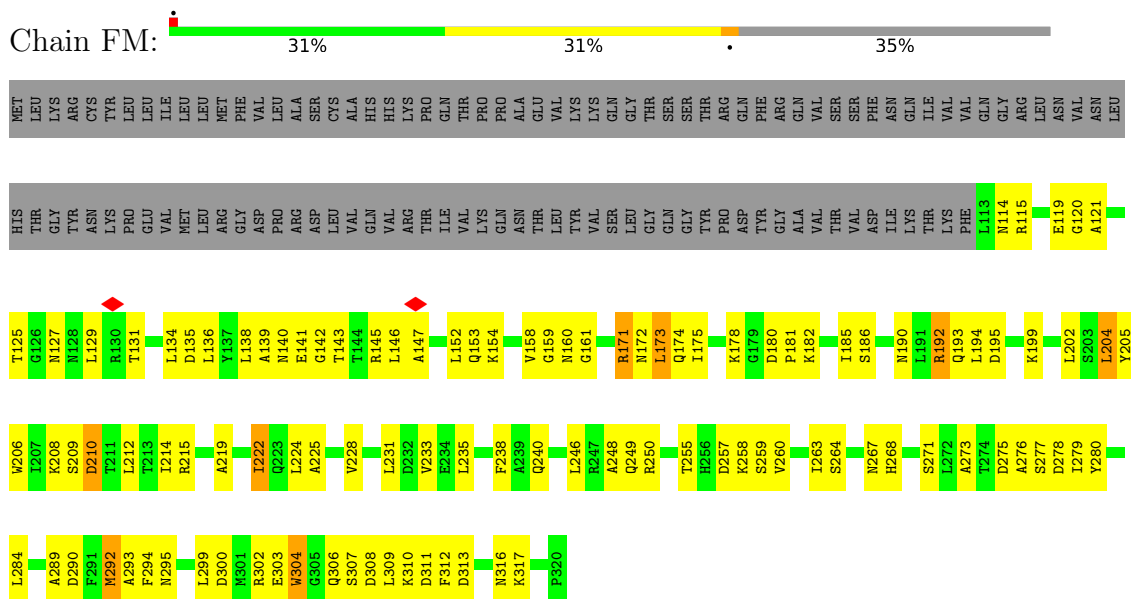
- Molecule 11: DUF2807 domain-containing protein



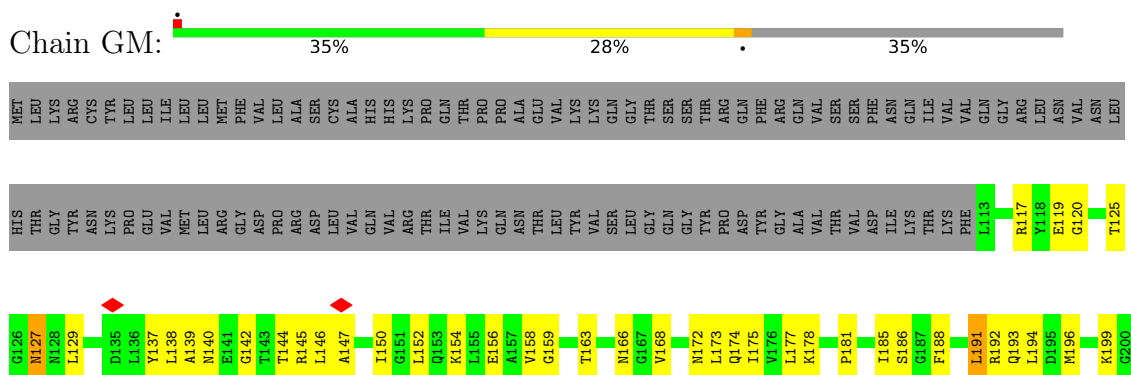
- Molecule 11: DUF2807 domain-containing protein

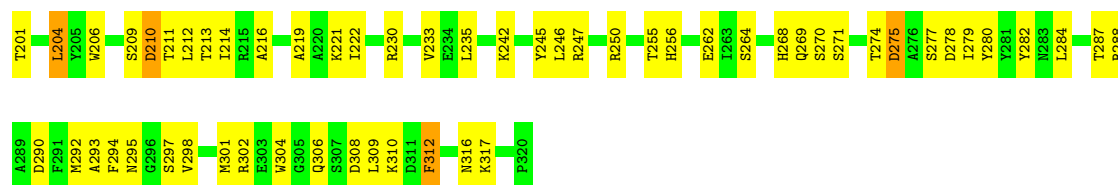


- Molecule 11: DUF2807 domain-containing protein

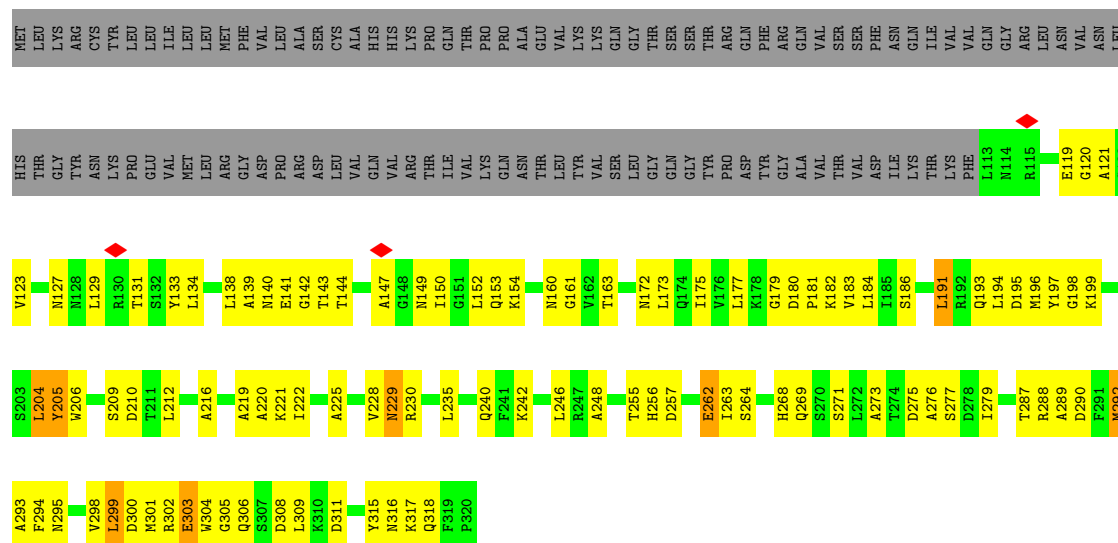


- Molecule 11: DUF2807 domain-containing protein

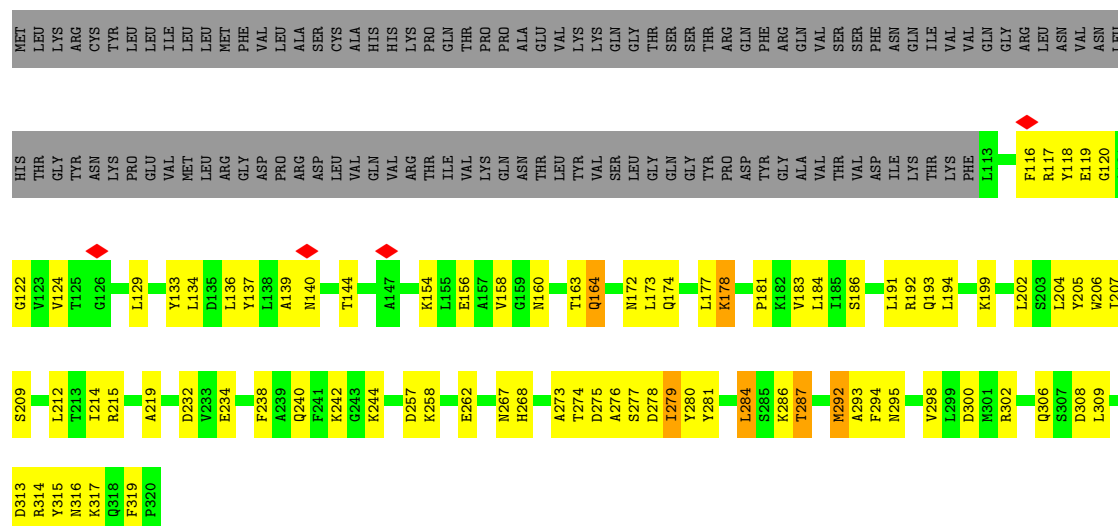




- Molecule 11: DUF2807 domain-containing protein



- Molecule 11: DUF2807 domain-containing protein



- Molecule 11: DUF2807 domain-containing protein





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	64698	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 ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	9.229	Depositor
Minimum map value	-3.621	Depositor
Average map value	0.091	Depositor
Map value standard deviation	0.546	Depositor
Recommended contour level	2.25	Depositor
Map size (Å)	561.0, 561.0, 561.0	wwPDB
Map dimensions	250, 250, 250	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	2.244, 2.244, 2.244	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	AG	0.15	0/1250	0.38	0/1699
1	Ag	0.12	0/278	0.23	0/377
1	BG	0.16	0/1250	0.40	0/1699
1	Bg	0.12	0/278	0.23	0/377
1	CG	0.16	0/1250	0.40	0/1699
1	Cg	0.11	0/278	0.24	0/377
1	DG	0.16	0/1250	0.40	0/1699
1	Dg	0.12	0/278	0.28	0/377
1	EG	0.16	0/1250	0.39	0/1699
1	Eg	0.12	0/278	0.25	0/377
1	FG	0.17	0/1250	0.43	0/1699
1	Fg	0.11	0/278	0.24	0/377
1	GG	0.17	0/1250	0.42	0/1699
1	Gg	0.11	0/278	0.26	0/377
1	HG	0.16	0/1250	0.40	0/1699
1	Hg	0.11	0/278	0.24	0/377
1	IG	0.16	0/1250	0.42	0/1699
1	Ig	0.11	0/278	0.26	0/377
1	JG	0.16	0/1250	0.40	0/1699
1	Jg	0.11	0/278	0.24	0/377
1	KG	0.16	0/1250	0.39	0/1699
1	Kg	0.11	0/278	0.27	0/377
1	LG	0.16	0/1250	0.42	0/1699
1	Lg	0.11	0/278	0.26	0/377
1	MG	0.17	0/1250	0.38	0/1699
1	Mg	0.11	0/278	0.26	0/377
1	NG	0.16	0/1250	0.39	0/1699
1	OG	0.16	0/1250	0.39	0/1699
1	PG	0.16	0/1250	0.39	0/1699
1	VG	0.12	0/278	0.32	0/377
1	WG	0.13	0/278	0.31	0/377
1	XG	0.15	0/278	0.36	0/377
1	YG	0.13	0/278	0.32	0/377
1	ZG	0.14	0/278	0.29	0/377

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
2	AH	0.15	0/2033	0.33	0/2775
2	BH	0.15	0/2033	0.32	0/2775
2	CH	0.14	0/2033	0.30	0/2775
2	DH	0.14	0/2033	0.30	0/2775
2	EH	0.15	0/2033	0.31	0/2775
2	FH	0.15	0/2033	0.31	0/2775
2	GH	0.14	0/2033	0.30	0/2775
2	HH	0.13	0/2033	0.27	0/2775
2	IH	0.14	0/2033	0.31	0/2775
2	JH	0.14	0/2033	0.29	0/2775
2	KH	0.14	0/2033	0.29	0/2775
2	LH	0.14	0/2033	0.33	0/2775
2	MH	0.14	0/2033	0.30	0/2775
2	VH	0.11	0/1921	0.29	0/2620
2	WH	0.11	0/1921	0.29	0/2620
2	XH	0.11	0/1921	0.29	0/2620
2	YH	0.11	0/1921	0.29	0/2620
2	ZH	0.12	0/1921	0.30	0/2620
3	AK	0.18	0/1195	0.31	0/1616
3	BK	0.18	0/1195	0.30	0/1616
3	CK	0.16	0/1195	0.32	0/1616
3	DK	0.16	0/1195	0.29	0/1616
3	EK	0.15	0/1195	0.29	0/1616
3	FK	0.16	0/1195	0.30	0/1616
3	GK	0.16	0/1195	0.32	0/1616
3	HK	0.16	0/1195	0.31	0/1616
3	IK	0.15	0/1195	0.30	0/1616
3	JK	0.16	0/1195	0.29	0/1616
3	KK	0.16	0/1195	0.30	0/1616
3	LK	0.16	0/1195	0.30	0/1616
3	MK	0.16	0/1195	0.30	0/1616
4	AL	0.14	0/1417	0.30	0/1912
4	BL	0.16	0/1417	0.34	0/1912
4	CL	0.15	0/1417	0.32	0/1912
4	DL	0.15	0/1417	0.32	0/1912
4	EL	0.15	0/1417	0.36	0/1912
4	FL	0.15	0/1417	0.32	0/1912
4	GL	0.15	0/1417	0.34	0/1912
4	HL	0.15	0/1417	0.34	0/1912
4	IL	0.15	0/1417	0.33	0/1912
4	JL	0.15	0/1417	0.33	0/1912
4	KL	0.15	0/1417	0.31	0/1912
4	LL	0.15	0/1417	0.31	0/1912

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
4	ML	0.15	0/1417	0.32	0/1912
5	AN	0.15	0/593	0.31	0/799
5	BN	0.13	0/593	0.27	0/799
5	CN	0.14	0/593	0.30	0/799
5	DN	0.15	0/593	0.29	0/799
5	EN	0.14	0/593	0.29	0/799
5	FN	0.15	0/593	0.27	0/799
5	GN	0.15	0/593	0.28	0/799
5	HN	0.13	0/593	0.26	0/799
5	IN	0.15	0/593	0.29	0/799
5	JN	0.13	0/593	0.27	0/799
5	KN	0.13	0/593	0.29	0/799
5	LN	0.14	0/593	0.29	0/799
5	MN	0.13	0/593	0.27	0/799
8	AD	0.16	0/1107	0.32	0/1502
8	Ad	0.15	0/1078	0.31	0/1463
8	BD	0.17	0/1107	0.33	0/1502
8	Bd	0.15	0/1078	0.27	0/1463
8	CD	0.15	0/1107	0.34	0/1502
8	Cd	0.14	0/1078	0.30	0/1463
8	DD	0.15	0/1107	0.32	0/1502
8	Dd	0.15	0/1078	0.30	0/1463
8	ED	0.16	0/1107	0.35	0/1502
8	Ed	0.14	0/1078	0.27	0/1463
8	FD	0.16	0/1107	0.36	0/1502
8	Fd	0.14	0/1078	0.28	0/1463
8	GD	0.16	0/1107	0.35	0/1502
8	Gd	0.16	0/1078	0.30	0/1463
8	HD	0.17	0/1107	0.34	0/1502
8	Hd	0.14	0/1078	0.27	0/1463
8	ID	0.16	0/1107	0.34	0/1502
8	Id	0.15	0/1078	0.28	0/1463
8	JD	0.15	0/1107	0.33	0/1502
8	Jd	0.14	0/1078	0.27	0/1463
8	KD	0.15	0/1107	0.34	0/1502
8	Kd	0.14	0/1078	0.27	0/1463
8	LD	0.16	0/1107	0.32	0/1502
8	Ld	0.15	0/1078	0.28	0/1463
8	MD	0.16	0/1107	0.33	0/1502
8	Md	0.14	0/1078	0.30	0/1463
9	AF	0.11	0/490	0.26	0/660
9	Af	0.11	0/456	0.31	0/615
9	BF	0.11	0/490	0.23	0/660

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
9	Bf	0.12	0/456	0.30	0/615
9	CF	0.11	0/490	0.30	0/660
9	Cf	0.11	0/456	0.24	0/615
9	DF	0.12	0/490	0.23	0/660
9	Df	0.13	0/456	0.28	0/615
9	EF	0.10	0/490	0.24	0/660
9	Ef	0.10	0/456	0.25	0/615
9	FF	0.10	0/490	0.25	0/660
9	Ff	0.12	0/456	0.30	0/615
9	GF	0.10	0/490	0.26	0/660
9	Gf	0.11	0/456	0.25	0/615
9	HF	0.11	0/490	0.28	0/660
9	Hf	0.12	0/456	0.32	0/615
9	IF	0.10	0/490	0.28	0/660
9	If	0.10	0/456	0.26	0/615
9	JF	0.10	0/490	0.27	0/660
9	Jf	0.11	0/456	0.26	0/615
9	KF	0.09	0/490	0.26	0/660
9	Kf	0.11	0/456	0.29	0/615
9	LF	0.10	0/490	0.30	0/660
9	Lf	0.12	0/456	0.25	0/615
9	MF	0.10	0/490	0.30	0/660
9	Mf	0.13	0/456	0.32	0/615
9	VF	0.12	0/490	0.26	0/660
9	WF	0.09	0/490	0.25	0/660
9	XF	0.11	0/490	0.26	0/660
9	YF	0.10	0/490	0.29	0/660
9	ZF	0.10	0/490	0.25	0/660
10	AC	0.15	0/1957	0.30	0/2651
10	BC	0.16	0/1957	0.31	0/2651
10	CC	0.17	0/1702	0.30	0/2315
10	DC	0.17	0/1702	0.32	0/2315
10	EC	0.17	0/1702	0.32	0/2315
10	FC	0.16	0/1957	0.30	0/2651
10	GC	0.16	0/1702	0.31	0/2315
10	HC	0.16	0/1702	0.30	0/2315
10	IC	0.16	0/1702	0.29	0/2315
10	JC	0.16	0/1957	0.32	0/2651
10	KC	0.16	0/1957	0.31	0/2651
10	LC	0.16	0/1702	0.30	0/2315
10	MC	0.17	0/1702	0.30	0/2315
11	AM	0.14	0/1678	0.32	0/2262
11	BM	0.13	0/1678	0.34	0/2262

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
11	CM	0.14	0/1678	0.32	0/2262
11	DM	0.14	0/1678	0.32	0/2262
11	EM	0.14	0/1678	0.32	0/2262
11	FM	0.13	0/1678	0.35	0/2262
11	GM	0.14	0/1678	0.33	0/2262
11	HM	0.14	0/1678	0.31	0/2262
11	IM	0.14	0/1678	0.33	0/2262
11	JM	0.14	0/1678	0.32	0/2262
11	KM	0.13	0/1678	0.31	0/2262
11	LM	0.14	0/1678	0.33	0/2262
11	MM	0.13	0/1678	0.32	0/2262
All	All	0.15	0/191071	0.32	0/258997

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	AG	1229	0	1231	65	0
1	Ag	276	0	263	12	0
1	BG	1229	0	1231	59	0
1	Bg	276	0	263	18	0
1	CG	1229	0	1231	66	0
1	Cg	276	0	263	19	0
1	DG	1229	0	1231	65	0
1	Dg	276	0	263	18	0
1	EG	1229	0	1231	55	0
1	Eg	276	0	263	21	0
1	FG	1229	0	1231	55	0
1	Fg	276	0	263	15	0
1	GG	1229	0	1231	66	0
1	Gg	276	0	263	21	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	HG	1229	0	1231	64	0
1	Hg	276	0	263	12	0
1	IG	1229	0	1231	67	0
1	Ig	276	0	263	23	0
1	JG	1229	0	1231	60	0
1	Jg	276	0	263	17	0
1	KG	1229	0	1231	64	0
1	Kg	276	0	263	22	0
1	LG	1229	0	1231	50	0
1	Lg	276	0	263	18	0
1	MG	1229	0	1231	56	0
1	Mg	276	0	263	18	0
1	NG	1229	0	1231	72	0
1	OG	1229	0	1231	76	0
1	PG	1229	0	1231	63	0
1	VG	276	0	263	16	0
1	WG	276	0	263	17	0
1	XG	276	0	263	16	0
1	YG	276	0	263	16	0
1	ZG	276	0	263	17	0
2	AH	1983	0	2009	78	0
2	BH	1983	0	2009	82	0
2	CH	1983	0	2009	65	0
2	DH	1983	0	2009	80	0
2	EH	1983	0	2009	85	0
2	FH	1983	0	2009	74	0
2	GH	1983	0	2009	56	0
2	HH	1983	0	2009	72	0
2	IH	1983	0	2009	70	0
2	JH	1983	0	2009	65	0
2	KH	1983	0	2009	69	0
2	LH	1983	0	2009	73	0
2	MH	1983	0	2009	66	0
2	VH	1875	0	1900	70	0
2	WH	1875	0	1900	51	0
2	XH	1875	0	1900	68	0
2	YH	1875	0	1900	72	0
2	ZH	1875	0	1900	63	0
3	AK	1175	0	1221	29	0
3	BK	1175	0	1221	35	0
3	CK	1175	0	1221	27	0
3	DK	1175	0	1221	36	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	EK	1175	0	1221	34	0
3	FK	1175	0	1221	47	0
3	GK	1175	0	1221	40	0
3	HK	1175	0	1221	38	0
3	IK	1175	0	1221	39	0
3	JK	1175	0	1221	35	0
3	KK	1175	0	1221	43	0
3	LK	1175	0	1221	39	0
3	MK	1175	0	1221	42	0
4	AL	1388	0	1378	41	0
4	BL	1388	0	1378	48	0
4	CL	1388	0	1378	33	0
4	DL	1388	0	1378	51	0
4	EL	1388	0	1378	47	0
4	FL	1388	0	1378	40	0
4	GL	1388	0	1378	56	0
4	HL	1388	0	1378	52	0
4	IL	1388	0	1378	46	0
4	JL	1388	0	1378	48	0
4	KL	1388	0	1378	57	0
4	LL	1388	0	1378	49	0
4	ML	1388	0	1378	39	0
5	AN	582	0	530	21	0
5	BN	582	0	530	27	0
5	CN	582	0	530	19	0
5	DN	582	0	530	20	0
5	EN	582	0	530	19	0
5	FN	582	0	530	17	0
5	GN	582	0	530	17	0
5	HN	582	0	530	19	0
5	IN	582	0	530	24	0
5	JN	582	0	530	25	0
5	KN	582	0	530	16	0
5	LN	582	0	530	25	0
5	MN	582	0	530	20	0
6	AU	45	0	12	1	0
6	BU	45	0	12	0	0
6	CU	45	0	12	0	0
6	DU	45	0	12	1	0
6	EU	45	0	12	1	0
6	FU	45	0	12	1	0
6	GU	45	0	12	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	HU	45	0	12	1	0
6	IU	45	0	12	1	0
6	JU	45	0	13	1	0
6	KU	45	0	13	1	0
6	LU	45	0	12	1	0
6	MU	45	0	12	0	0
7	AX	240	0	56	2	0
7	BX	240	0	57	5	0
7	CX	240	0	57	2	0
7	DX	240	0	56	2	0
7	EX	240	0	55	0	0
7	FX	240	0	54	0	0
7	GX	240	0	54	0	0
7	HX	240	0	55	1	0
7	IX	240	0	55	1	0
7	JX	240	0	57	2	0
7	KX	240	0	55	1	0
7	LX	240	0	55	0	0
7	MX	240	0	58	3	0
7	VX	240	0	54	1	0
7	WX	240	0	52	1	0
7	XX	240	0	54	1	0
7	YX	240	0	55	4	0
7	ZX	240	0	54	1	0
8	AD	1086	0	1121	44	0
8	Ad	1058	0	1092	36	0
8	BD	1086	0	1121	37	0
8	Bd	1058	0	1092	36	0
8	CD	1086	0	1121	34	0
8	Cd	1058	0	1092	40	0
8	DD	1086	0	1121	42	0
8	Dd	1058	0	1092	23	0
8	ED	1086	0	1121	38	0
8	Ed	1058	0	1092	38	0
8	FD	1086	0	1121	44	0
8	Fd	1058	0	1092	46	0
8	GD	1086	0	1121	41	0
8	Gd	1058	0	1092	31	0
8	HD	1086	0	1121	37	0
8	Hd	1058	0	1092	33	0
8	ID	1086	0	1121	34	0
8	Id	1058	0	1092	28	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
8	JD	1086	0	1121	32	0
8	Jd	1058	0	1092	30	0
8	KD	1086	0	1121	36	0
8	Kd	1058	0	1092	28	0
8	LD	1086	0	1121	33	0
8	Ld	1058	0	1092	34	0
8	MD	1086	0	1121	25	0
8	Md	1058	0	1092	28	0
9	AF	483	0	502	23	0
9	Af	449	0	474	11	0
9	BF	483	0	502	15	0
9	Bf	449	0	474	15	0
9	CF	483	0	502	23	0
9	Cf	449	0	474	14	0
9	DF	483	0	502	22	0
9	Df	449	0	474	18	0
9	EF	483	0	502	17	0
9	Ef	449	0	474	17	0
9	FF	483	0	502	17	0
9	Ff	449	0	474	19	0
9	GF	483	0	502	20	0
9	Gf	449	0	474	12	0
9	HF	483	0	502	18	0
9	Hf	449	0	474	12	0
9	IF	483	0	502	15	0
9	If	449	0	474	26	0
9	JF	483	0	502	16	0
9	Jf	449	0	474	11	0
9	KF	483	0	502	18	0
9	Kf	449	0	474	19	0
9	LF	483	0	502	13	0
9	Lf	449	0	474	15	0
9	MF	483	0	502	18	0
9	Mf	449	0	474	11	0
9	VF	483	0	502	22	0
9	WF	483	0	502	20	0
9	XF	483	0	502	22	0
9	YF	483	0	502	21	0
9	ZF	483	0	502	10	0
10	AC	1921	0	1952	84	0
10	BC	1921	0	1952	63	0
10	CC	1667	0	1682	68	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
10	DC	1667	0	1682	70	0
10	EC	1667	0	1682	56	0
10	FC	1921	0	1952	65	0
10	GC	1667	0	1682	74	0
10	HC	1667	0	1682	53	0
10	IC	1667	0	1682	74	0
10	JC	1921	0	1952	65	0
10	KC	1921	0	1952	77	0
10	LC	1667	0	1682	67	0
10	MC	1667	0	1682	66	0
11	AM	1650	0	1658	65	0
11	BM	1650	0	1658	72	0
11	CM	1650	0	1658	57	0
11	DM	1650	0	1658	46	0
11	EM	1650	0	1658	68	0
11	FM	1650	0	1658	73	0
11	GM	1650	0	1658	67	0
11	HM	1650	0	1658	59	0
11	IM	1650	0	1658	51	0
11	JM	1650	0	1658	72	0
11	KM	1650	0	1658	64	0
11	LM	1650	0	1658	45	0
11	MM	1650	0	1658	54	0
All	All	192370	0	190612	6122	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 16.

The worst 5 of 6122 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)
11:EM:198:GLY:O	11:EM:219:ALA:HB3	1.66	0.96
1:ZG:798:THR:HA	2:BH:115:MET:HE1	1.50	0.92
11:JM:275:ASP:HB3	11:JM:294:PHE:HB2	1.50	0.91
5:LN:78:LEU:H	2:MH:354:MET:HE1	1.43	0.83
4:DL:166:ALA:HA	4:DL:206:GLU:O	1.77	0.83

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	AG	161/1048 (15%)	149 (92%)	12 (8%)	0	100	100
1	Ag	32/1048 (3%)	30 (94%)	2 (6%)	0	100	100
1	BG	161/1048 (15%)	154 (96%)	7 (4%)	0	100	100
1	Bg	32/1048 (3%)	30 (94%)	2 (6%)	0	100	100
1	CG	161/1048 (15%)	155 (96%)	6 (4%)	0	100	100
1	Cg	32/1048 (3%)	30 (94%)	2 (6%)	0	100	100
1	DG	161/1048 (15%)	153 (95%)	8 (5%)	0	100	100
1	Dg	32/1048 (3%)	31 (97%)	1 (3%)	0	100	100
1	EG	161/1048 (15%)	153 (95%)	8 (5%)	0	100	100
1	Eg	32/1048 (3%)	31 (97%)	1 (3%)	0	100	100
1	FG	161/1048 (15%)	151 (94%)	10 (6%)	0	100	100
1	Fg	32/1048 (3%)	31 (97%)	1 (3%)	0	100	100
1	GG	161/1048 (15%)	153 (95%)	8 (5%)	0	100	100
1	Gg	32/1048 (3%)	32 (100%)	0	0	100	100
1	HG	161/1048 (15%)	156 (97%)	5 (3%)	0	100	100
1	Hg	32/1048 (3%)	32 (100%)	0	0	100	100
1	IG	161/1048 (15%)	154 (96%)	7 (4%)	0	100	100
1	Ig	32/1048 (3%)	30 (94%)	2 (6%)	0	100	100
1	JG	161/1048 (15%)	154 (96%)	7 (4%)	0	100	100
1	Jg	32/1048 (3%)	31 (97%)	1 (3%)	0	100	100
1	KG	161/1048 (15%)	150 (93%)	11 (7%)	0	100	100
1	Kg	32/1048 (3%)	30 (94%)	2 (6%)	0	100	100
1	LG	161/1048 (15%)	151 (94%)	10 (6%)	0	100	100
1	Lg	32/1048 (3%)	31 (97%)	1 (3%)	0	100	100
1	MG	161/1048 (15%)	150 (93%)	11 (7%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	Mg	32/1048 (3%)	31 (97%)	1 (3%)	0	100	100
1	NG	161/1048 (15%)	153 (95%)	8 (5%)	0	100	100
1	OG	161/1048 (15%)	154 (96%)	7 (4%)	0	100	100
1	PG	161/1048 (15%)	155 (96%)	6 (4%)	0	100	100
1	VG	32/1048 (3%)	32 (100%)	0	0	100	100
1	WG	32/1048 (3%)	32 (100%)	0	0	100	100
1	XG	32/1048 (3%)	31 (97%)	1 (3%)	0	100	100
1	YG	32/1048 (3%)	31 (97%)	1 (3%)	0	100	100
1	ZG	32/1048 (3%)	32 (100%)	0	0	100	100
2	AH	256/361 (71%)	241 (94%)	15 (6%)	0	100	100
2	BH	256/361 (71%)	242 (94%)	14 (6%)	0	100	100
2	CH	256/361 (71%)	244 (95%)	12 (5%)	0	100	100
2	DH	256/361 (71%)	241 (94%)	15 (6%)	0	100	100
2	EH	256/361 (71%)	239 (93%)	17 (7%)	0	100	100
2	FH	256/361 (71%)	245 (96%)	11 (4%)	0	100	100
2	GH	256/361 (71%)	238 (93%)	18 (7%)	0	100	100
2	HH	256/361 (71%)	246 (96%)	10 (4%)	0	100	100
2	IH	256/361 (71%)	240 (94%)	16 (6%)	0	100	100
2	JH	256/361 (71%)	245 (96%)	11 (4%)	0	100	100
2	KH	256/361 (71%)	242 (94%)	14 (6%)	0	100	100
2	LH	256/361 (71%)	243 (95%)	13 (5%)	0	100	100
2	MH	256/361 (71%)	240 (94%)	16 (6%)	0	100	100
2	VH	239/361 (66%)	227 (95%)	12 (5%)	0	100	100
2	WH	239/361 (66%)	235 (98%)	4 (2%)	0	100	100
2	XH	239/361 (66%)	234 (98%)	5 (2%)	0	100	100
2	YH	239/361 (66%)	229 (96%)	10 (4%)	0	100	100
2	ZH	239/361 (66%)	230 (96%)	9 (4%)	0	100	100
3	AK	149/189 (79%)	143 (96%)	6 (4%)	0	100	100
3	BK	149/189 (79%)	143 (96%)	6 (4%)	0	100	100
3	CK	149/189 (79%)	143 (96%)	6 (4%)	0	100	100
3	DK	149/189 (79%)	142 (95%)	7 (5%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	EK	149/189 (79%)	144 (97%)	5 (3%)	0	100	100
3	FK	149/189 (79%)	143 (96%)	6 (4%)	0	100	100
3	GK	149/189 (79%)	143 (96%)	6 (4%)	0	100	100
3	HK	149/189 (79%)	140 (94%)	9 (6%)	0	100	100
3	IK	149/189 (79%)	143 (96%)	6 (4%)	0	100	100
3	JK	149/189 (79%)	144 (97%)	5 (3%)	0	100	100
3	KK	149/189 (79%)	144 (97%)	5 (3%)	0	100	100
3	LK	149/189 (79%)	141 (95%)	8 (5%)	0	100	100
3	MK	149/189 (79%)	143 (96%)	6 (4%)	0	100	100
4	AL	169/249 (68%)	162 (96%)	7 (4%)	0	100	100
4	BL	169/249 (68%)	165 (98%)	4 (2%)	0	100	100
4	CL	169/249 (68%)	162 (96%)	7 (4%)	0	100	100
4	DL	169/249 (68%)	161 (95%)	8 (5%)	0	100	100
4	EL	169/249 (68%)	161 (95%)	8 (5%)	0	100	100
4	FL	169/249 (68%)	163 (96%)	6 (4%)	0	100	100
4	GL	169/249 (68%)	162 (96%)	7 (4%)	0	100	100
4	HL	169/249 (68%)	162 (96%)	7 (4%)	0	100	100
4	IL	169/249 (68%)	162 (96%)	7 (4%)	0	100	100
4	JL	169/249 (68%)	163 (96%)	6 (4%)	0	100	100
4	KL	169/249 (68%)	165 (98%)	4 (2%)	0	100	100
4	LL	169/249 (68%)	162 (96%)	7 (4%)	0	100	100
4	ML	169/249 (68%)	161 (95%)	8 (5%)	0	100	100
5	AN	76/124 (61%)	72 (95%)	4 (5%)	0	100	100
5	BN	76/124 (61%)	70 (92%)	6 (8%)	0	100	100
5	CN	76/124 (61%)	69 (91%)	7 (9%)	0	100	100
5	DN	76/124 (61%)	69 (91%)	7 (9%)	0	100	100
5	EN	76/124 (61%)	67 (88%)	9 (12%)	0	100	100
5	FN	76/124 (61%)	70 (92%)	6 (8%)	0	100	100
5	GN	76/124 (61%)	72 (95%)	4 (5%)	0	100	100
5	HN	76/124 (61%)	73 (96%)	3 (4%)	0	100	100
5	IN	76/124 (61%)	72 (95%)	4 (5%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
5	JN	76/124 (61%)	70 (92%)	6 (8%)	0	100	100
5	KN	76/124 (61%)	70 (92%)	6 (8%)	0	100	100
5	LN	76/124 (61%)	72 (95%)	4 (5%)	0	100	100
5	MN	76/124 (61%)	71 (93%)	5 (7%)	0	100	100
8	AD	138/163 (85%)	132 (96%)	6 (4%)	0	100	100
8	Ad	135/163 (83%)	128 (95%)	7 (5%)	0	100	100
8	BD	138/163 (85%)	132 (96%)	6 (4%)	0	100	100
8	Bd	135/163 (83%)	128 (95%)	7 (5%)	0	100	100
8	CD	138/163 (85%)	130 (94%)	8 (6%)	0	100	100
8	Cd	135/163 (83%)	129 (96%)	6 (4%)	0	100	100
8	DD	138/163 (85%)	132 (96%)	6 (4%)	0	100	100
8	Dd	135/163 (83%)	131 (97%)	4 (3%)	0	100	100
8	ED	138/163 (85%)	132 (96%)	6 (4%)	0	100	100
8	Ed	135/163 (83%)	130 (96%)	5 (4%)	0	100	100
8	FD	138/163 (85%)	133 (96%)	5 (4%)	0	100	100
8	Fd	135/163 (83%)	129 (96%)	6 (4%)	0	100	100
8	GD	138/163 (85%)	132 (96%)	6 (4%)	0	100	100
8	Gd	135/163 (83%)	125 (93%)	10 (7%)	0	100	100
8	HD	138/163 (85%)	131 (95%)	7 (5%)	0	100	100
8	Hd	135/163 (83%)	127 (94%)	8 (6%)	0	100	100
8	ID	138/163 (85%)	131 (95%)	7 (5%)	0	100	100
8	Id	135/163 (83%)	131 (97%)	4 (3%)	0	100	100
8	JD	138/163 (85%)	128 (93%)	10 (7%)	0	100	100
8	Jd	135/163 (83%)	128 (95%)	7 (5%)	0	100	100
8	KD	138/163 (85%)	131 (95%)	7 (5%)	0	100	100
8	Kd	135/163 (83%)	125 (93%)	10 (7%)	0	100	100
8	LD	138/163 (85%)	127 (92%)	11 (8%)	0	100	100
8	Ld	135/163 (83%)	129 (96%)	6 (4%)	0	100	100
8	MD	138/163 (85%)	132 (96%)	6 (4%)	0	100	100
8	Md	135/163 (83%)	129 (96%)	6 (4%)	0	100	100
9	AF	61/269 (23%)	59 (97%)	2 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
9	Af	57/269 (21%)	53 (93%)	4 (7%)	0	100	100
9	BF	61/269 (23%)	57 (93%)	4 (7%)	0	100	100
9	Bf	57/269 (21%)	54 (95%)	3 (5%)	0	100	100
9	CF	61/269 (23%)	60 (98%)	1 (2%)	0	100	100
9	Cf	57/269 (21%)	55 (96%)	2 (4%)	0	100	100
9	DF	61/269 (23%)	61 (100%)	0	0	100	100
9	Df	57/269 (21%)	54 (95%)	3 (5%)	0	100	100
9	EF	61/269 (23%)	59 (97%)	2 (3%)	0	100	100
9	Ef	57/269 (21%)	53 (93%)	4 (7%)	0	100	100
9	FF	61/269 (23%)	58 (95%)	3 (5%)	0	100	100
9	Ff	57/269 (21%)	53 (93%)	4 (7%)	0	100	100
9	GF	61/269 (23%)	59 (97%)	2 (3%)	0	100	100
9	Gf	57/269 (21%)	53 (93%)	4 (7%)	0	100	100
9	HF	61/269 (23%)	57 (93%)	4 (7%)	0	100	100
9	Hf	57/269 (21%)	55 (96%)	2 (4%)	0	100	100
9	IF	61/269 (23%)	53 (87%)	8 (13%)	0	100	100
9	If	57/269 (21%)	54 (95%)	3 (5%)	0	100	100
9	JF	61/269 (23%)	59 (97%)	2 (3%)	0	100	100
9	Jf	57/269 (21%)	55 (96%)	2 (4%)	0	100	100
9	KF	61/269 (23%)	55 (90%)	6 (10%)	0	100	100
9	Kf	57/269 (21%)	51 (90%)	6 (10%)	0	100	100
9	LF	61/269 (23%)	56 (92%)	5 (8%)	0	100	100
9	Lf	57/269 (21%)	53 (93%)	4 (7%)	0	100	100
9	MF	61/269 (23%)	57 (93%)	4 (7%)	0	100	100
9	Mf	57/269 (21%)	56 (98%)	1 (2%)	0	100	100
9	VF	61/269 (23%)	56 (92%)	5 (8%)	0	100	100
9	WF	61/269 (23%)	57 (93%)	4 (7%)	0	100	100
9	XF	61/269 (23%)	58 (95%)	3 (5%)	0	100	100
9	YF	61/269 (23%)	60 (98%)	1 (2%)	0	100	100
9	ZF	61/269 (23%)	58 (95%)	3 (5%)	0	100	100
10	AC	241/303 (80%)	225 (93%)	16 (7%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
10	BC	241/303 (80%)	229 (95%)	12 (5%)	0	100	100
10	CC	207/303 (68%)	191 (92%)	16 (8%)	0	100	100
10	DC	207/303 (68%)	188 (91%)	19 (9%)	0	100	100
10	EC	207/303 (68%)	192 (93%)	15 (7%)	0	100	100
10	FC	241/303 (80%)	231 (96%)	10 (4%)	0	100	100
10	GC	207/303 (68%)	194 (94%)	13 (6%)	0	100	100
10	HC	207/303 (68%)	198 (96%)	9 (4%)	0	100	100
10	IC	207/303 (68%)	188 (91%)	19 (9%)	0	100	100
10	JC	241/303 (80%)	227 (94%)	14 (6%)	0	100	100
10	KC	241/303 (80%)	226 (94%)	15 (6%)	0	100	100
10	LC	207/303 (68%)	193 (93%)	14 (7%)	0	100	100
10	MC	207/303 (68%)	193 (93%)	14 (7%)	0	100	100
11	AM	206/320 (64%)	188 (91%)	18 (9%)	0	100	100
11	BM	206/320 (64%)	190 (92%)	16 (8%)	0	100	100
11	CM	206/320 (64%)	194 (94%)	12 (6%)	0	100	100
11	DM	206/320 (64%)	192 (93%)	14 (7%)	0	100	100
11	EM	206/320 (64%)	187 (91%)	19 (9%)	0	100	100
11	FM	206/320 (64%)	192 (93%)	14 (7%)	0	100	100
11	GM	206/320 (64%)	187 (91%)	19 (9%)	0	100	100
11	HM	206/320 (64%)	185 (90%)	21 (10%)	0	100	100
11	IM	206/320 (64%)	188 (91%)	18 (9%)	0	100	100
11	JM	206/320 (64%)	192 (93%)	14 (7%)	0	100	100
11	KM	206/320 (64%)	193 (94%)	13 (6%)	0	100	100
11	LM	206/320 (64%)	193 (94%)	13 (6%)	0	100	100
11	MM	206/320 (64%)	194 (94%)	12 (6%)	0	100	100
All	All	23724/70112 (34%)	22448 (95%)	1276 (5%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

entries.

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	AG	135/765 (18%)	125 (93%)	10 (7%)	13	33
1	Ag	31/765 (4%)	26 (84%)	5 (16%)	2	12
1	BG	135/765 (18%)	125 (93%)	10 (7%)	13	33
1	Bg	31/765 (4%)	27 (87%)	4 (13%)	4	17
1	CG	135/765 (18%)	124 (92%)	11 (8%)	11	31
1	Cg	31/765 (4%)	30 (97%)	1 (3%)	34	55
1	DG	135/765 (18%)	120 (89%)	15 (11%)	6	20
1	Dg	31/765 (4%)	27 (87%)	4 (13%)	4	17
1	EG	135/765 (18%)	123 (91%)	12 (9%)	9	28
1	Eg	31/765 (4%)	29 (94%)	2 (6%)	15	37
1	FG	135/765 (18%)	119 (88%)	16 (12%)	5	19
1	Fg	31/765 (4%)	30 (97%)	1 (3%)	34	55
1	GG	135/765 (18%)	123 (91%)	12 (9%)	9	28
1	Gg	31/765 (4%)	25 (81%)	6 (19%)	1	8
1	HG	135/765 (18%)	122 (90%)	13 (10%)	8	25
1	Hg	31/765 (4%)	26 (84%)	5 (16%)	2	12
1	IG	135/765 (18%)	122 (90%)	13 (10%)	8	25
1	Ig	31/765 (4%)	27 (87%)	4 (13%)	4	17
1	JG	135/765 (18%)	121 (90%)	14 (10%)	7	22
1	Jg	31/765 (4%)	28 (90%)	3 (10%)	8	25
1	KG	135/765 (18%)	123 (91%)	12 (9%)	9	28
1	Kg	31/765 (4%)	27 (87%)	4 (13%)	4	17
1	LG	135/765 (18%)	124 (92%)	11 (8%)	11	31
1	Lg	31/765 (4%)	26 (84%)	5 (16%)	2	12
1	MG	135/765 (18%)	125 (93%)	10 (7%)	13	33
1	Mg	31/765 (4%)	27 (87%)	4 (13%)	4	17
1	NG	135/765 (18%)	125 (93%)	10 (7%)	13	33
1	OG	135/765 (18%)	122 (90%)	13 (10%)	8	25
1	PG	135/765 (18%)	124 (92%)	11 (8%)	11	31

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	VG	31/765 (4%)	29 (94%)	2 (6%)	15	37
1	WG	31/765 (4%)	26 (84%)	5 (16%)	2	12
1	XG	31/765 (4%)	28 (90%)	3 (10%)	8	25
1	YG	31/765 (4%)	25 (81%)	6 (19%)	1	8
1	ZG	31/765 (4%)	25 (81%)	6 (19%)	1	8
2	AH	220/300 (73%)	196 (89%)	24 (11%)	6	21
2	BH	220/300 (73%)	203 (92%)	17 (8%)	12	32
2	CH	220/300 (73%)	204 (93%)	16 (7%)	13	34
2	DH	220/300 (73%)	205 (93%)	15 (7%)	14	36
2	EH	220/300 (73%)	200 (91%)	20 (9%)	9	28
2	FH	220/300 (73%)	206 (94%)	14 (6%)	16	37
2	GH	220/300 (73%)	206 (94%)	14 (6%)	16	37
2	HH	220/300 (73%)	209 (95%)	11 (5%)	22	43
2	IH	220/300 (73%)	212 (96%)	8 (4%)	31	52
2	JH	220/300 (73%)	202 (92%)	18 (8%)	10	31
2	KH	220/300 (73%)	201 (91%)	19 (9%)	10	29
2	LH	220/300 (73%)	203 (92%)	17 (8%)	12	32
2	MH	220/300 (73%)	200 (91%)	20 (9%)	9	28
2	VH	207/300 (69%)	190 (92%)	17 (8%)	10	31
2	WH	207/300 (69%)	184 (89%)	23 (11%)	6	20
2	XH	207/300 (69%)	186 (90%)	21 (10%)	7	23
2	YH	207/300 (69%)	184 (89%)	23 (11%)	6	20
2	ZH	207/300 (69%)	190 (92%)	17 (8%)	10	31
3	AK	129/163 (79%)	119 (92%)	10 (8%)	11	32
3	BK	129/163 (79%)	120 (93%)	9 (7%)	14	36
3	CK	129/163 (79%)	121 (94%)	8 (6%)	16	38
3	DK	129/163 (79%)	118 (92%)	11 (8%)	10	29
3	EK	129/163 (79%)	122 (95%)	7 (5%)	20	42
3	FK	129/163 (79%)	122 (95%)	7 (5%)	20	42
3	GK	129/163 (79%)	115 (89%)	14 (11%)	6	21
3	HK	129/163 (79%)	122 (95%)	7 (5%)	20	42

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	IK	129/163 (79%)	116 (90%)	13 (10%)	7	23
3	JK	129/163 (79%)	116 (90%)	13 (10%)	7	23
3	KK	129/163 (79%)	120 (93%)	9 (7%)	14	36
3	LK	129/163 (79%)	116 (90%)	13 (10%)	7	23
3	MK	129/163 (79%)	111 (86%)	18 (14%)	3	15
4	AL	148/203 (73%)	140 (95%)	8 (5%)	20	42
4	BL	148/203 (73%)	136 (92%)	12 (8%)	11	31
4	CL	148/203 (73%)	139 (94%)	9 (6%)	17	39
4	DL	148/203 (73%)	138 (93%)	10 (7%)	14	36
4	EL	148/203 (73%)	139 (94%)	9 (6%)	17	39
4	FL	148/203 (73%)	134 (90%)	14 (10%)	8	25
4	GL	148/203 (73%)	141 (95%)	7 (5%)	23	45
4	HL	148/203 (73%)	139 (94%)	9 (6%)	17	39
4	IL	148/203 (73%)	137 (93%)	11 (7%)	13	33
4	JL	148/203 (73%)	139 (94%)	9 (6%)	17	39
4	KL	148/203 (73%)	132 (89%)	16 (11%)	6	22
4	LL	148/203 (73%)	141 (95%)	7 (5%)	23	45
4	ML	148/203 (73%)	135 (91%)	13 (9%)	9	29
5	AN	66/107 (62%)	61 (92%)	5 (8%)	12	33
5	BN	66/107 (62%)	60 (91%)	6 (9%)	9	28
5	CN	66/107 (62%)	59 (89%)	7 (11%)	6	22
5	DN	66/107 (62%)	61 (92%)	5 (8%)	12	33
5	EN	66/107 (62%)	62 (94%)	4 (6%)	17	39
5	FN	66/107 (62%)	62 (94%)	4 (6%)	17	39
5	GN	66/107 (62%)	60 (91%)	6 (9%)	9	28
5	HN	66/107 (62%)	63 (96%)	3 (4%)	24	46
5	IN	66/107 (62%)	61 (92%)	5 (8%)	12	33
5	JN	66/107 (62%)	61 (92%)	5 (8%)	12	33
5	KN	66/107 (62%)	63 (96%)	3 (4%)	24	46
5	LN	66/107 (62%)	63 (96%)	3 (4%)	24	46
5	MN	66/107 (62%)	56 (85%)	10 (15%)	3	13

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
8	AD	121/139 (87%)	114 (94%)	7 (6%)	18	40
8	Ad	118/139 (85%)	108 (92%)	10 (8%)	10	29
8	BD	121/139 (87%)	114 (94%)	7 (6%)	18	40
8	Bd	118/139 (85%)	109 (92%)	9 (8%)	12	33
8	CD	121/139 (87%)	118 (98%)	3 (2%)	42	62
8	Cd	118/139 (85%)	111 (94%)	7 (6%)	18	40
8	DD	121/139 (87%)	114 (94%)	7 (6%)	18	40
8	Dd	118/139 (85%)	110 (93%)	8 (7%)	14	36
8	ED	121/139 (87%)	114 (94%)	7 (6%)	18	40
8	Ed	118/139 (85%)	104 (88%)	14 (12%)	5	19
8	FD	121/139 (87%)	112 (93%)	9 (7%)	13	33
8	Fd	118/139 (85%)	109 (92%)	9 (8%)	12	33
8	GD	121/139 (87%)	110 (91%)	11 (9%)	9	28
8	Gd	118/139 (85%)	111 (94%)	7 (6%)	18	40
8	HD	121/139 (87%)	112 (93%)	9 (7%)	13	33
8	Hd	118/139 (85%)	108 (92%)	10 (8%)	10	29
8	ID	121/139 (87%)	113 (93%)	8 (7%)	15	37
8	Id	118/139 (85%)	114 (97%)	4 (3%)	32	54
8	JD	121/139 (87%)	112 (93%)	9 (7%)	13	33
8	Jd	118/139 (85%)	109 (92%)	9 (8%)	12	33
8	KD	121/139 (87%)	112 (93%)	9 (7%)	13	33
8	Kd	118/139 (85%)	112 (95%)	6 (5%)	21	43
8	LD	121/139 (87%)	106 (88%)	15 (12%)	4	18
8	Ld	118/139 (85%)	110 (93%)	8 (7%)	14	36
8	MD	121/139 (87%)	113 (93%)	8 (7%)	15	37
8	Md	118/139 (85%)	107 (91%)	11 (9%)	8	26
9	AF	53/237 (22%)	49 (92%)	4 (8%)	12	33
9	Af	49/237 (21%)	44 (90%)	5 (10%)	7	23
9	BF	53/237 (22%)	50 (94%)	3 (6%)	18	40
9	Bf	49/237 (21%)	45 (92%)	4 (8%)	10	31
9	CF	53/237 (22%)	48 (91%)	5 (9%)	8	26

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
9	Cf	49/237 (21%)	44 (90%)	5 (10%)	7	23
9	DF	53/237 (22%)	50 (94%)	3 (6%)	18	40
9	Df	49/237 (21%)	44 (90%)	5 (10%)	7	23
9	EF	53/237 (22%)	49 (92%)	4 (8%)	12	33
9	Ef	49/237 (21%)	45 (92%)	4 (8%)	10	31
9	FF	53/237 (22%)	49 (92%)	4 (8%)	12	33
9	Ff	49/237 (21%)	45 (92%)	4 (8%)	10	31
9	GF	53/237 (22%)	49 (92%)	4 (8%)	12	33
9	Gf	49/237 (21%)	47 (96%)	2 (4%)	27	49
9	HF	53/237 (22%)	48 (91%)	5 (9%)	8	26
9	Hf	49/237 (21%)	46 (94%)	3 (6%)	17	39
9	IF	53/237 (22%)	50 (94%)	3 (6%)	18	40
9	If	49/237 (21%)	44 (90%)	5 (10%)	7	23
9	JF	53/237 (22%)	46 (87%)	7 (13%)	4	16
9	Jf	49/237 (21%)	46 (94%)	3 (6%)	17	39
9	KF	53/237 (22%)	45 (85%)	8 (15%)	3	13
9	Kf	49/237 (21%)	46 (94%)	3 (6%)	17	39
9	LF	53/237 (22%)	49 (92%)	4 (8%)	12	33
9	Lf	49/237 (21%)	44 (90%)	5 (10%)	7	23
9	MF	53/237 (22%)	49 (92%)	4 (8%)	12	33
9	Mf	49/237 (21%)	46 (94%)	3 (6%)	17	39
9	VF	53/237 (22%)	52 (98%)	1 (2%)	50	66
9	WF	53/237 (22%)	50 (94%)	3 (6%)	18	40
9	XF	53/237 (22%)	51 (96%)	2 (4%)	29	51
9	YF	53/237 (22%)	49 (92%)	4 (8%)	12	33
9	ZF	53/237 (22%)	48 (91%)	5 (9%)	8	26
10	AC	203/257 (79%)	188 (93%)	15 (7%)	13	33
10	BC	203/257 (79%)	176 (87%)	27 (13%)	4	16
10	CC	178/257 (69%)	159 (89%)	19 (11%)	6	22
10	DC	178/257 (69%)	161 (90%)	17 (10%)	8	25
10	EC	178/257 (69%)	155 (87%)	23 (13%)	4	17

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
10	FC	203/257 (79%)	189 (93%)	14 (7%)	14	36
10	GC	178/257 (69%)	161 (90%)	17 (10%)	8	25
10	HC	178/257 (69%)	158 (89%)	20 (11%)	6	20
10	IC	178/257 (69%)	161 (90%)	17 (10%)	8	25
10	JC	203/257 (79%)	189 (93%)	14 (7%)	14	36
10	KC	203/257 (79%)	185 (91%)	18 (9%)	9	28
10	LC	178/257 (69%)	157 (88%)	21 (12%)	5	19
10	MC	178/257 (69%)	156 (88%)	22 (12%)	4	18
11	AM	175/276 (63%)	163 (93%)	12 (7%)	14	36
11	BM	175/276 (63%)	162 (93%)	13 (7%)	13	33
11	CM	175/276 (63%)	162 (93%)	13 (7%)	13	33
11	DM	175/276 (63%)	162 (93%)	13 (7%)	13	33
11	EM	175/276 (63%)	160 (91%)	15 (9%)	10	29
11	FM	175/276 (63%)	157 (90%)	18 (10%)	7	23
11	GM	175/276 (63%)	161 (92%)	14 (8%)	11	31
11	HM	175/276 (63%)	168 (96%)	7 (4%)	28	49
11	IM	175/276 (63%)	156 (89%)	19 (11%)	6	21
11	JM	175/276 (63%)	155 (89%)	20 (11%)	5	20
11	KM	175/276 (63%)	161 (92%)	14 (8%)	11	31
11	LM	175/276 (63%)	158 (90%)	17 (10%)	8	25
11	MM	175/276 (63%)	162 (93%)	13 (7%)	13	33
All	All	20484/55449 (37%)	18791 (92%)	1693 (8%)	12	30

5 of 1693 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	XH	233	LEU
8	Ad	97	ARG
1	MG	930	ILE
1	YG	815	VAL
2	XH	220	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 206 such sidechains are listed below:

Mol	Chain	Res	Type
1	ZG	818	GLN
4	BL	225	GLN
1	PG	1017	GLN
2	ZH	205	GLN
11	FM	174	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 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
7	EX	1
7	GX	1
7	YX	1
7	VX	1
7	CX	1
7	AX	1
7	FX	1
7	JX	1

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Mol	Chain	Number of breaks
7	KX	1
7	MX	1
7	WX	1
7	XX	1
7	LX	1
7	BX	1
7	HX	1
7	ZX	1
7	IX	1
7	DX	1

The worst 5 of 18 chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	EX	38:UNK	C	70:UNK	N	24.84
1	GX	38:UNK	C	70:UNK	N	24.65
1	YX	38:UNK	C	70:UNK	N	24.61
1	VX	38:UNK	C	70:UNK	N	24.52
1	CX	38:UNK	C	70:UNK	N	24.52

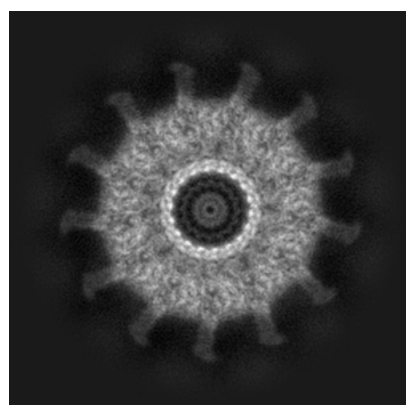
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-24023. These allow visual inspection of the internal detail of the map and identification of artifacts.

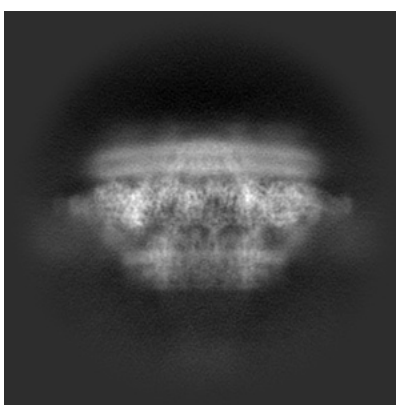
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

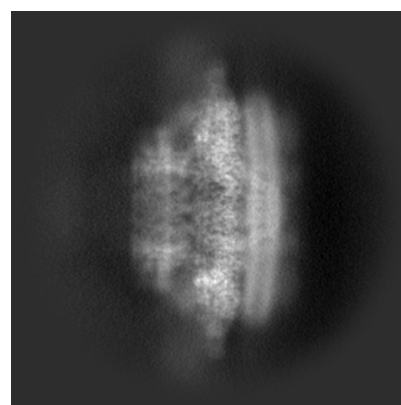
6.1.1 Primary map



X



Y

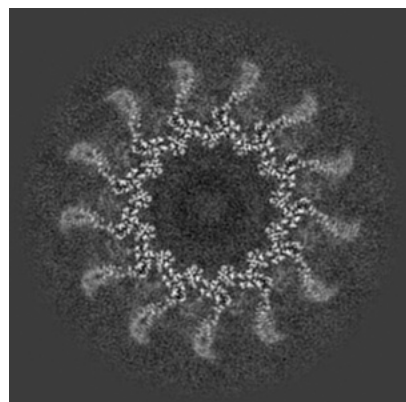


Z

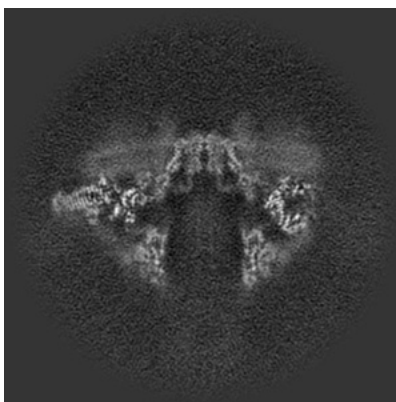
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

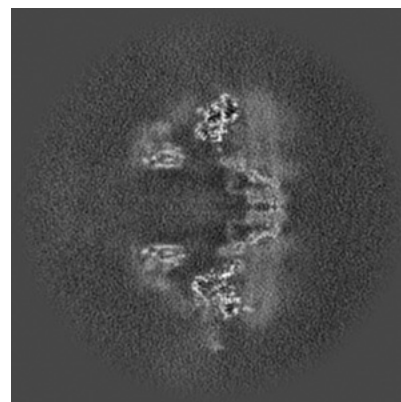
6.2.1 Primary map



X Index: 125



Y Index: 125

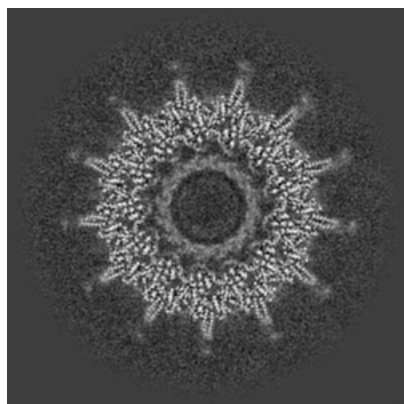


Z Index: 125

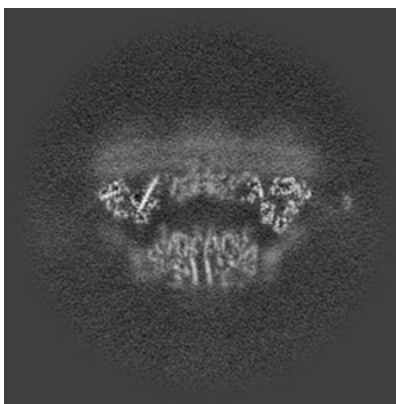
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

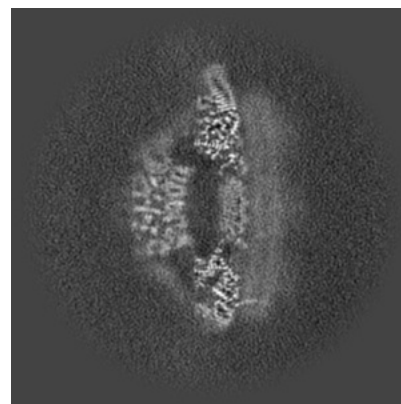
6.3.1 Primary map



X Index: 134



Y Index: 101

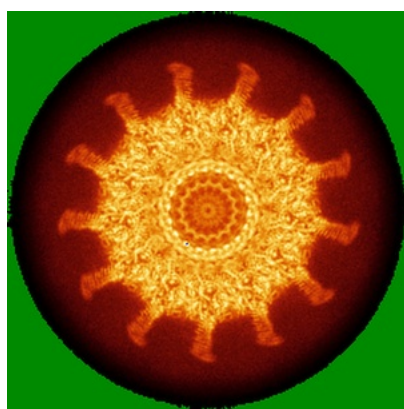


Z Index: 151

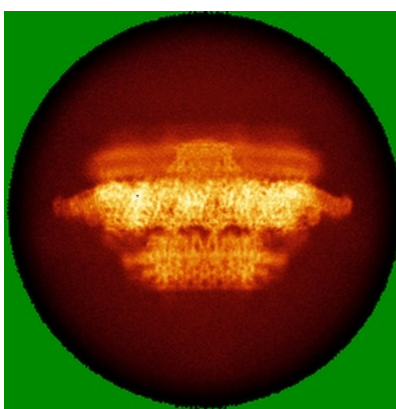
The images above show the largest variance slices of the map in three orthogonal directions.

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

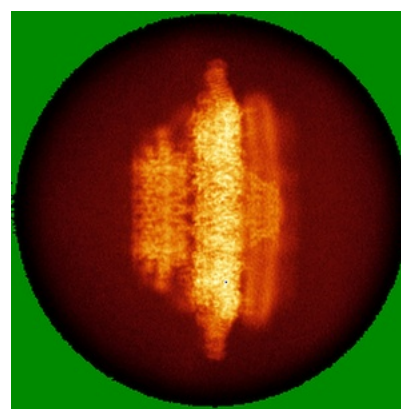
6.4.1 Primary map



X



Y

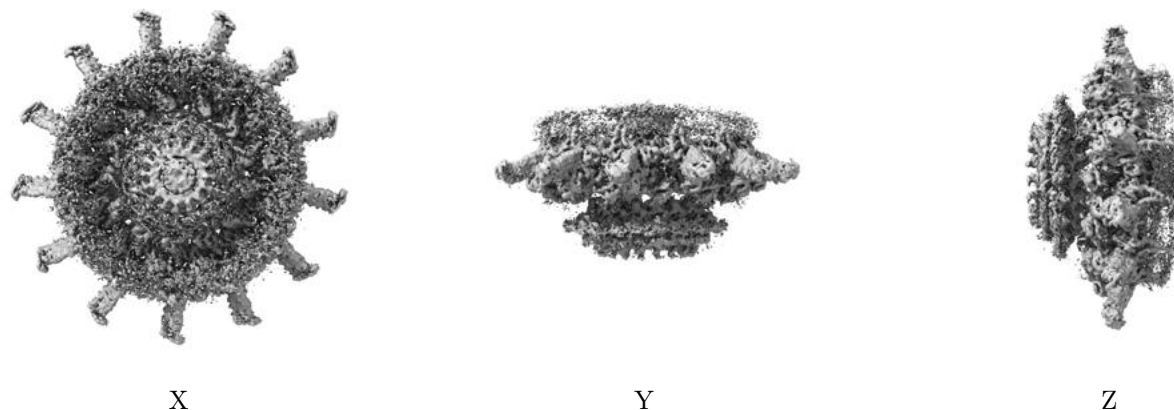


Z

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 2.25. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

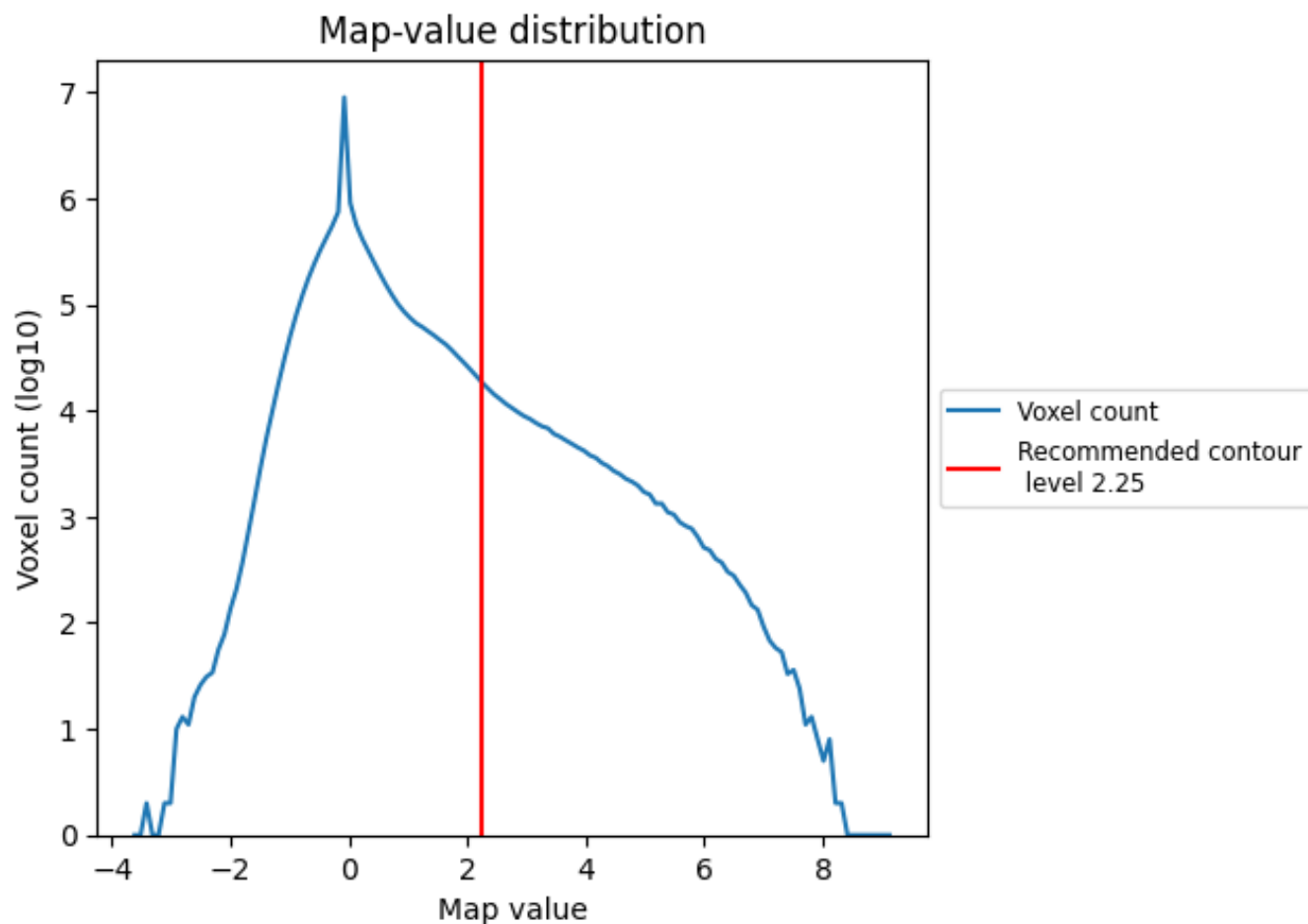
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

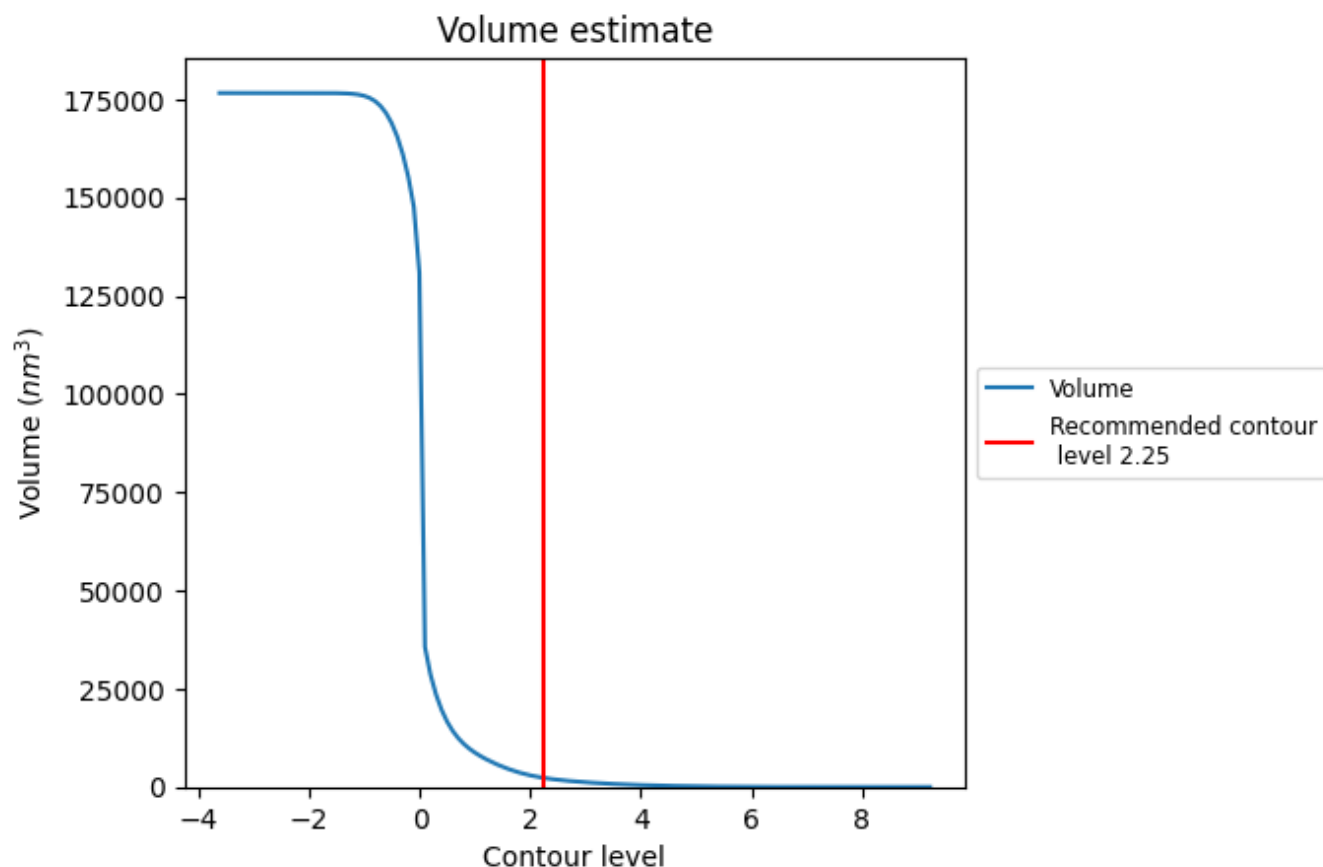
This section contains the results of statistical analysis of the map.

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.

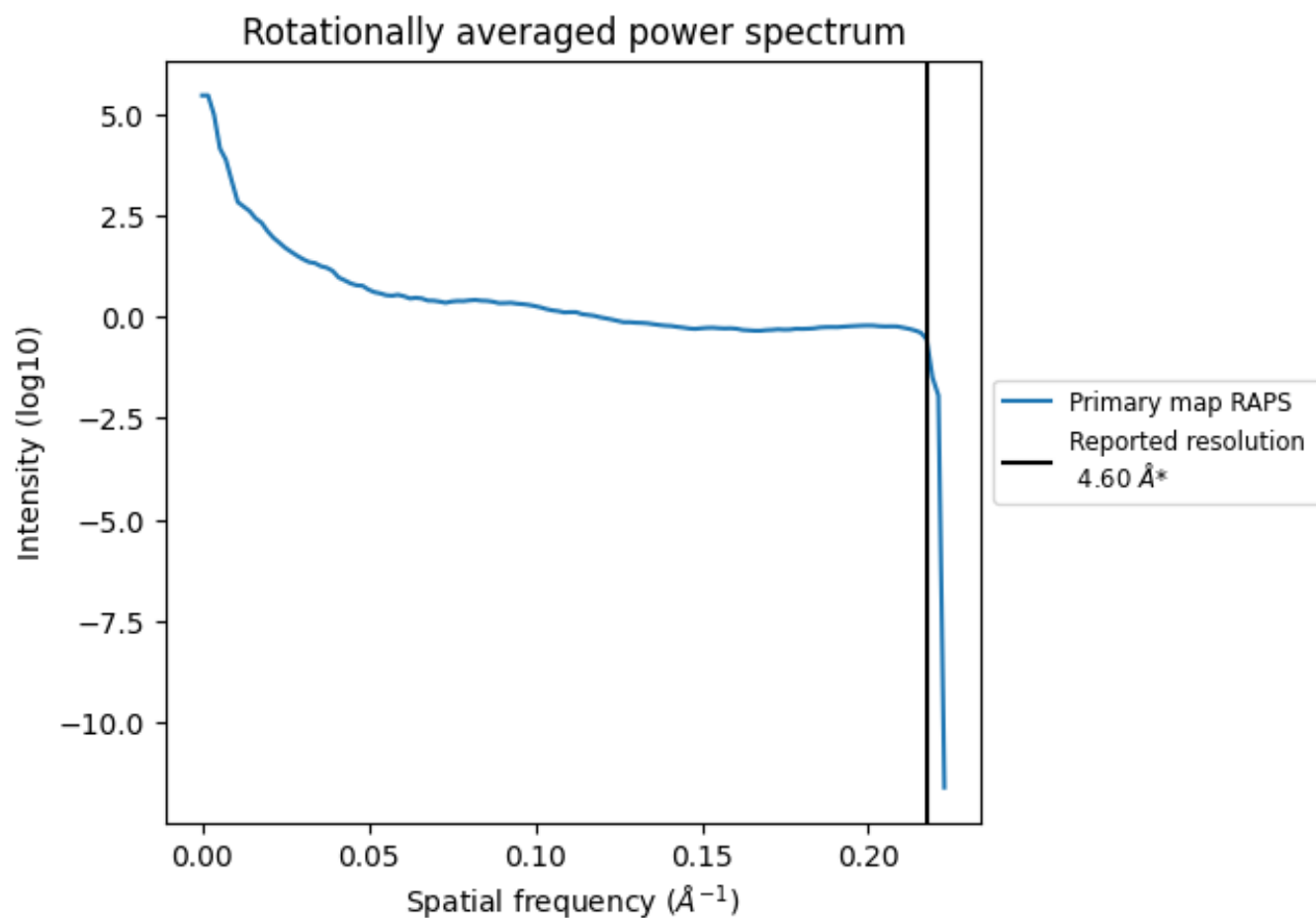
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 2290 nm³; this corresponds to an approximate mass of 2069 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 0.217 $\text{\AA}^{-1}$

8 Fourier-Shell correlation

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

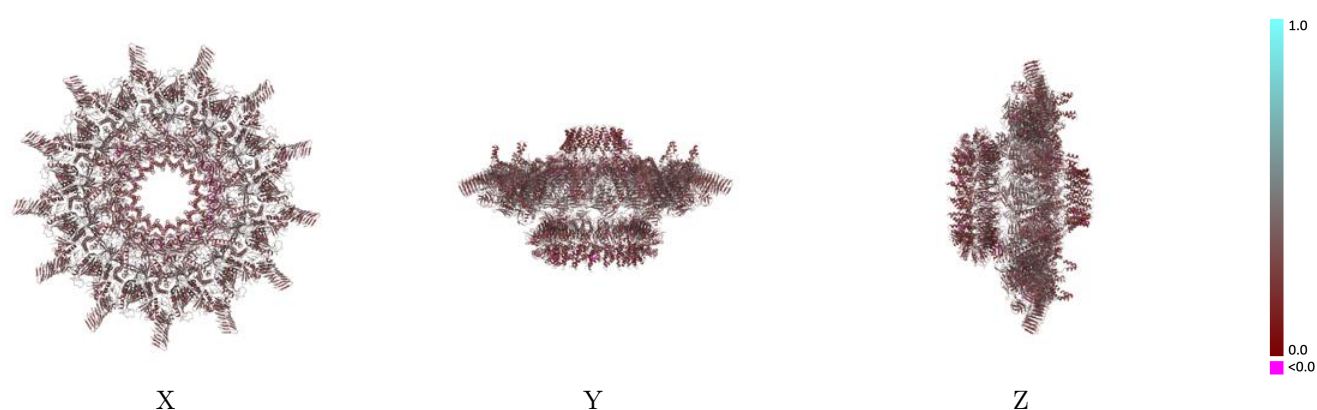
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-24023 and PDB model 7MUV. Per-residue inclusion information can be found in section [3](#) on page [23](#).

9.1 Map-model overlay [i](#)

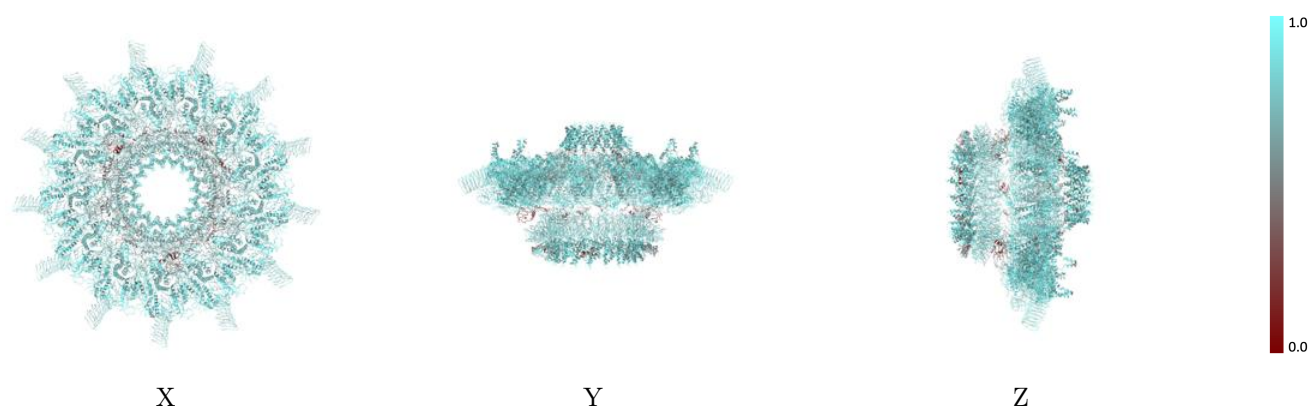
This section was not generated.

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



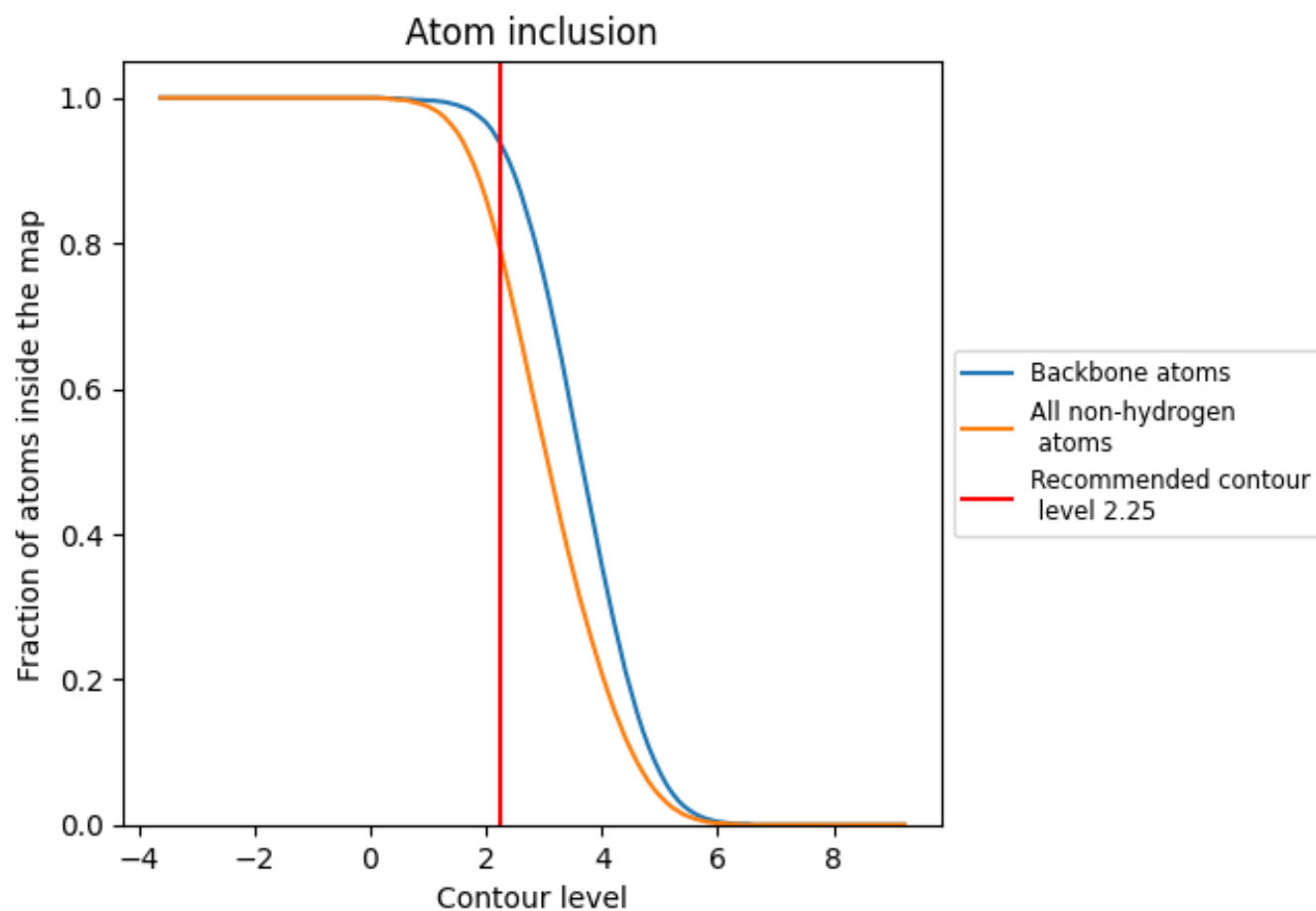
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (2.25).

9.4 Atom inclusion [i](#)



At the recommended contour level, 94% of all backbone atoms, 79% 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 (2.25) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.7900	 0.3160
AC	 0.7910	 0.3300
AD	 0.8170	 0.3540
AF	 0.7810	 0.2420
AG	 0.6540	 0.2680
AH	 0.8060	 0.3150
AK	 0.8390	 0.3520
AL	 0.8400	 0.3280
AM	 0.8120	 0.3270
AN	 0.7920	 0.3720
AU	 0.9780	 0.4790
AX	 0.6210	 0.2160
Ad	 0.8370	 0.3450
Af	 0.7950	 0.3120
Ag	 0.8310	 0.2930
BC	 0.7890	 0.3280
BD	 0.8170	 0.3640
BF	 0.7880	 0.2480
BG	 0.6470	 0.2540
BH	 0.7950	 0.3180
BK	 0.8320	 0.3460
BL	 0.8350	 0.3320
BM	 0.8190	 0.3220
BN	 0.7670	 0.3530
BU	 1.0000	 0.4510
BX	 0.5710	 0.2670
Bd	 0.8310	 0.3400
Bf	 0.8180	 0.3170
Bg	 0.8310	 0.2720
CC	 0.8250	 0.3270
CD	 0.8240	 0.3540
CF	 0.7220	 0.2560
CG	 0.6560	 0.2470
CH	 0.8150	 0.3260
CK	 0.8350	 0.3520



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Chain	Atom inclusion	Q-score
CL	 0.8490	 0.3310
CM	 0.8060	 0.3270
CN	 0.7710	 0.3610
CU	 1.0000	 0.4840
CX	 0.6330	 0.3280
Cd	 0.8370	 0.3330
Cf	 0.7860	 0.3050
Cg	 0.8050	 0.2830
DC	 0.8530	 0.3310
DD	 0.8330	 0.3500
DF	 0.7900	 0.2560
DG	 0.6650	 0.2520
DH	 0.8020	 0.3090
DK	 0.8520	 0.3550
DL	 0.8430	 0.3320
DM	 0.8000	 0.3230
DN	 0.8090	 0.3700
DU	 0.9560	 0.4600
DX	 0.6000	 0.2440
Dd	 0.8240	 0.3380
Df	 0.7900	 0.3130
Dg	 0.8490	 0.2700
EC	 0.8360	 0.3380
ED	 0.8120	 0.3470
EF	 0.7920	 0.2660
EG	 0.6560	 0.2530
EH	 0.8160	 0.3140
EK	 0.8390	 0.3450
EL	 0.8470	 0.3380
EM	 0.8230	 0.3300
EN	 0.8040	 0.3600
EU	 1.0000	 0.4750
EX	 0.5330	 0.2410
Ed	 0.8380	 0.3450
Ef	 0.7930	 0.3120
Eg	 0.8710	 0.2920
FC	 0.7930	 0.3300
FD	 0.8320	 0.3460
FF	 0.7620	 0.2750
FG	 0.6910	 0.2440
FH	 0.7920	 0.3080
FK	 0.8280	 0.3490

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Chain	Atom inclusion	Q-score
FL	 0.8580	 0.3350
FM	 0.8190	 0.3170
FN	 0.7850	 0.3610
FU	 0.9330	 0.4430
FX	 0.6620	 0.2630
Fd	 0.8400	 0.3330
Ff	 0.7810	 0.3070
Fg	 0.8640	 0.2760
GC	 0.8310	 0.3350
GD	 0.8190	 0.3590
GF	 0.7660	 0.2370
GG	 0.6690	 0.2290
GH	 0.8110	 0.3180
GK	 0.8410	 0.3520
GL	 0.8440	 0.3350
GM	 0.8420	 0.3320
GN	 0.7710	 0.3580
GU	 1.0000	 0.4310
GX	 0.6210	 0.2760
Gd	 0.8340	 0.3420
Gf	 0.7490	 0.3190
Gg	 0.8270	 0.2770
HC	 0.8490	 0.3370
HD	 0.8240	 0.3560
HF	 0.7920	 0.2750
HG	 0.6470	 0.2210
HH	 0.8060	 0.3160
HK	 0.8490	 0.3450
HL	 0.8490	 0.3290
HM	 0.8400	 0.3300
HN	 0.8250	 0.3720
HU	 0.9560	 0.4740
HX	 0.5330	 0.2320
Hd	 0.8470	 0.3440
Hf	 0.8430	 0.3040
Hg	 0.8090	 0.2740
IC	 0.8410	 0.3450
ID	 0.8390	 0.3520
IF	 0.8020	 0.2490
IG	 0.6130	 0.2190
IH	 0.8020	 0.3150
IK	 0.8500	 0.3530

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Chain	Atom inclusion	Q-score
IL	 0.8360	 0.3350
IM	 0.8140	 0.3230
IN	 0.8250	 0.3640
IU	 0.9780	 0.4440
IX	 0.5460	 0.2540
Id	 0.8420	 0.3400
If	 0.8130	 0.2930
Ig	 0.8460	 0.2780
JC	 0.7810	 0.3290
JD	 0.8290	 0.3430
JF	 0.7750	 0.2470
JG	 0.6050	 0.2030
JH	 0.8210	 0.3110
JK	 0.8500	 0.3450
JL	 0.8510	 0.3270
JM	 0.8310	 0.3370
JN	 0.7970	 0.3470
JU	 1.0000	 0.4940
JX	 0.6540	 0.2890
Jd	 0.8310	 0.3310
Jf	 0.7930	 0.3150
Jg	 0.8380	 0.2510
KC	 0.7800	 0.3320
KD	 0.8290	 0.3440
KF	 0.7620	 0.2430
KG	 0.6030	 0.1930
KH	 0.8140	 0.3070
KK	 0.8530	 0.3620
KL	 0.8500	 0.3270
KM	 0.8350	 0.3340
KN	 0.7760	 0.3630
KU	 0.9560	 0.4680
KX	 0.5670	 0.2750
Kd	 0.8520	 0.3390
Kf	 0.8020	 0.3060
Kg	 0.8350	 0.2850
LC	 0.8370	 0.3450
LD	 0.8400	 0.3450
LF	 0.7640	 0.2590
LG	 0.6220	 0.1990
LH	 0.8190	 0.3190
LK	 0.8510	 0.3500

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Chain	Atom inclusion	Q-score
LL	 0.8410	 0.3350
LM	 0.8200	 0.3280
LN	 0.7880	 0.3650
LU	 0.9780	 0.4900
LX	 0.6710	 0.2650
Ld	 0.8210	 0.3420
Lf	 0.8060	 0.3300
Lg	 0.8600	 0.2730
MC	 0.8440	 0.3440
MD	 0.8250	 0.3510
MF	 0.7600	 0.2660
MG	 0.6780	 0.2680
MH	 0.8180	 0.3190
MK	 0.8430	 0.3510
ML	 0.8410	 0.3290
MM	 0.8130	 0.3180
MN	 0.8510	 0.3560
MU	 1.0000	 0.4940
MX	 0.6250	 0.2650
Md	 0.8440	 0.3390
Mf	 0.8130	 0.3070
Mg	 0.8310	 0.2970
NG	 0.6620	 0.2630
OG	 0.7310	 0.2650
PG	 0.6860	 0.2770
VF	 0.8130	 0.2690
VG	 0.8420	 0.2760
VH	 0.6680	 0.2990
VX	 0.5210	 0.2560
WF	 0.7640	 0.2630
WG	 0.8160	 0.2840
WH	 0.5420	 0.2340
WX	 0.5880	 0.2760
XF	 0.7880	 0.2710
XG	 0.7900	 0.2520
XH	 0.6960	 0.2910
XX	 0.6420	 0.2460
YF	 0.7750	 0.2460
YG	 0.8460	 0.2770
YH	 0.5510	 0.2760
YX	 0.6370	 0.2940
ZF	 0.7960	 0.2710

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Chain	Atom inclusion	Q-score
ZG	 0.8120	 0.2820
ZH	 0.6650	 0.3070
ZX	 0.5500	 0.2810