Full wwPDB EM Validation Report (i)

Feb 25, 2024 – 08:06 AM EST

PDB ID	:	6X6K
EMDB ID	:	EMD-22076
Title	:	Cryo-EM Structure of the Helicobacter pylori dCag3 OMC
Authors	:	Sheedlo, M.J.; Chung, J.M.; Sawhney, N.; Durie, C.L.; Cover, T.L.; Ohi, M.D.; Lacy, D.B.
Deposited on	:	2020-05-28
Resolution	:	3.10 Å(reported)

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

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/EMValidationReportHelp
with specific help available everywhere you see the (i) 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 (i)) were used in the production of this report:

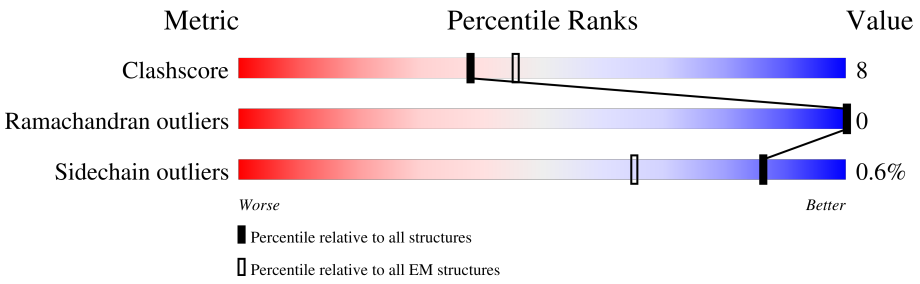
EMDB validation analysis	:	0.0.1.dev70
MolProbity	:	4.02b-467
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ	:	1.9.13
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 3.10 Å.

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



Metric	Whole archive (#Entries)	EM structures (#Entries)
Clashscore	158937	4297
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for >=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 <=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	AT	278	38% 53% 11% • 36%
1	BT	278	36% 52% 12% • 36%
1	CT	278	36% 51% 12% • 36%
1	DT	278	37% 52% 12% • 36%
1	ET	278	36% 53% 11% • 36%
1	FT	278	37% 51% 13% • 36%
1	GT	278	38% 53% 10% • 36%
1	HT	278	37% 52% 12% • 36%

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Mol	Chain	Length	Quality of chain
1	IT	278	37% 52% 12% • 36%
1	JT	278	37% 52% 12% • 36%
1	KT	278	37% 51% 13% • 36%
1	LT	278	36% 51% 13% • 36%
1	MT	278	37% 51% 12% • 36%
1	NT	278	37% 53% 10% • 36%
2	AX	522	7% 24% 6% 70%
2	BX	522	7% 24% 6% 70%
2	CX	522	7% 24% 5% 70%
2	DX	522	7% 24% 6% 70%
2	EX	522	7% 24% 6% 70%
2	FX	522	6% 23% 7% 70%
2	GX	522	7% 23% 7% 70%
2	HX	522	7% 23% 7% 70%
2	IX	522	7% 25% 5% 70%
2	JX	522	7% 24% 6% 70%
2	KX	522	6% 24% 6% 70%
2	LX	522	7% 24% 6% 70%
2	MX	522	6% 24% 6% 70%
2	NX	522	7% 23% 6% 70%
3	AY	1927	• 8% • 89%
3	BY	1927	• 8% • 89%
3	CY	1927	• 8% • 89%
3	DY	1927	• 8% • 89%
3	EY	1927	• 8% • 89%

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Mol	Chain	Length	Quality of chain
3	FY	1927	8% 89%
3	GY	1927	8% 89%
3	HY	1927	8% 89%
3	IY	1927	8% 89%
3	JY	1927	8% 89%
3	KY	1927	8% 89%
3	LY	1927	8% 89%
3	MY	1927	8% 89%
3	NY	1927	8% 89%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 60970 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 Cag pathogenicity island protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	AT	179	1502	957	261	282	2	0	0
1	BT	179	1502	957	261	282	2	0	0
1	CT	179	1502	957	261	282	2	0	0
1	DT	179	1502	957	261	282	2	0	0
1	ET	179	1502	957	261	282	2	0	0
1	FT	179	1502	957	261	282	2	0	0
1	GT	179	1502	957	261	282	2	0	0
1	HT	179	1502	957	261	282	2	0	0
1	IT	179	1502	957	261	282	2	0	0
1	JT	179	1502	957	261	282	2	0	0
1	KT	179	1502	957	261	282	2	0	0
1	LT	179	1502	957	261	282	2	0	0
1	MT	179	1502	957	261	282	2	0	0
1	NT	179	1502	957	261	282	2	0	0

- Molecule 2 is a protein called Type IV secretion system apparatus protein CagX.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
2	AX	155	1275	819	222	231	3	0	0

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Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
2	BX	155	1275	819	222	231	3	0	0
2	CX	155	1275	819	222	231	3	0	0
2	DX	155	1275	819	222	231	3	0	0
2	EX	155	1275	819	222	231	3	0	0
2	FX	155	1275	819	222	231	3	0	0
2	GX	155	1275	819	222	231	3	0	0
2	HX	155	1275	819	222	231	3	0	0
2	IX	155	1275	819	222	231	3	0	0
2	JX	155	1275	819	222	231	3	0	0
2	KX	155	1275	819	222	231	3	0	0
2	LX	155	1275	819	222	231	3	0	0
2	MX	155	1275	819	222	231	3	0	0
2	NX	155	1275	819	222	231	3	0	0

- Molecule 3 is a protein called Cag pathogenicity island protein (Cag7).

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
3	AY	209	1578	1014	255	300	9	0	0
3	BY	209	1578	1014	255	300	9	0	0
3	CY	209	1578	1014	255	300	9	0	0
3	DY	209	1578	1014	255	300	9	0	0
3	EY	209	1578	1014	255	300	9	0	0
3	FY	209	1578	1014	255	300	9	0	0

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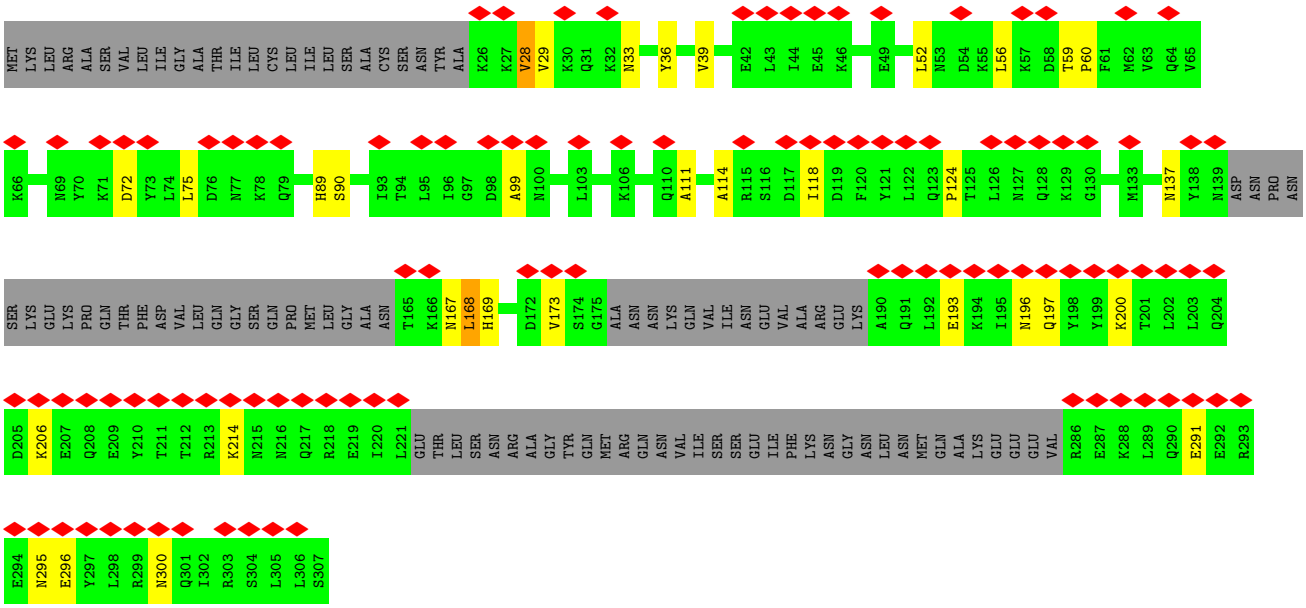
Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
3	GY	209	1578	1014	255	300	9	0	0
			Total	C	N	O	S		
3	HY	209	1578	1014	255	300	9	0	0
			Total	C	N	O	S		
3	IY	209	1578	1014	255	300	9	0	0
			Total	C	N	O	S		
3	JY	209	1578	1014	255	300	9	0	0
			Total	C	N	O	S		
3	KY	209	1578	1014	255	300	9	0	0
			Total	C	N	O	S		
3	LY	209	1578	1014	255	300	9	0	0
			Total	C	N	O	S		
3	MY	209	1578	1014	255	300	9	0	0
			Total	C	N	O	S		
3	NY	209	1578	1014	255	300	9	0	0

3 Residue-property plots (i)

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

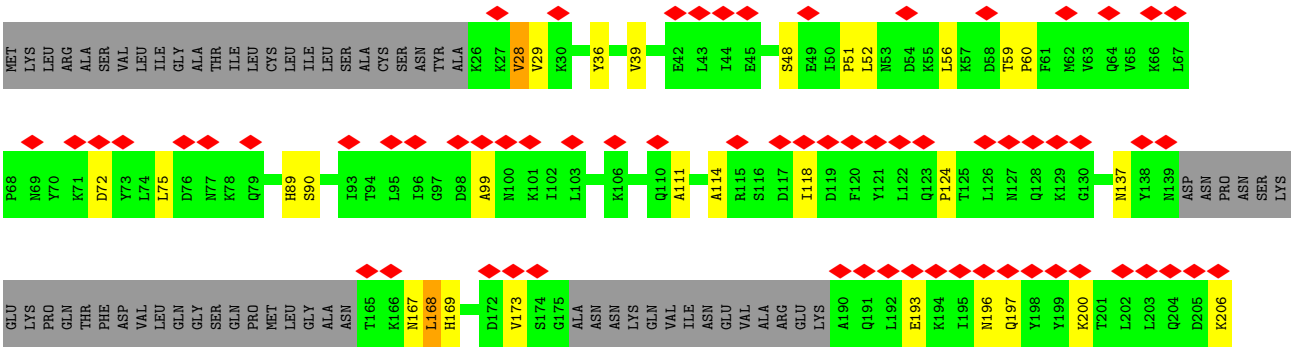
- Molecule 1: Cag pathogenicity island protein

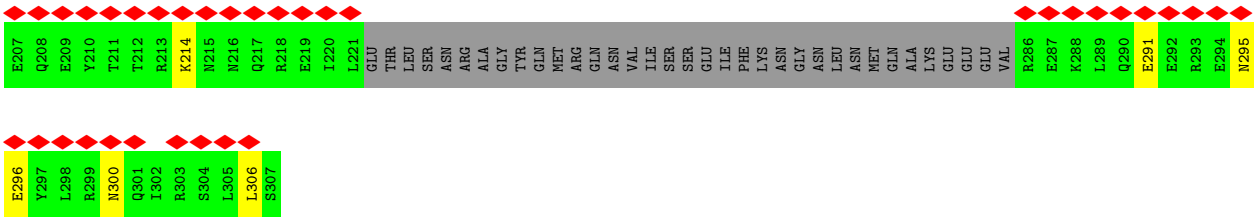
Chain AT:



- Molecule 1: Cag pathogenicity island protein

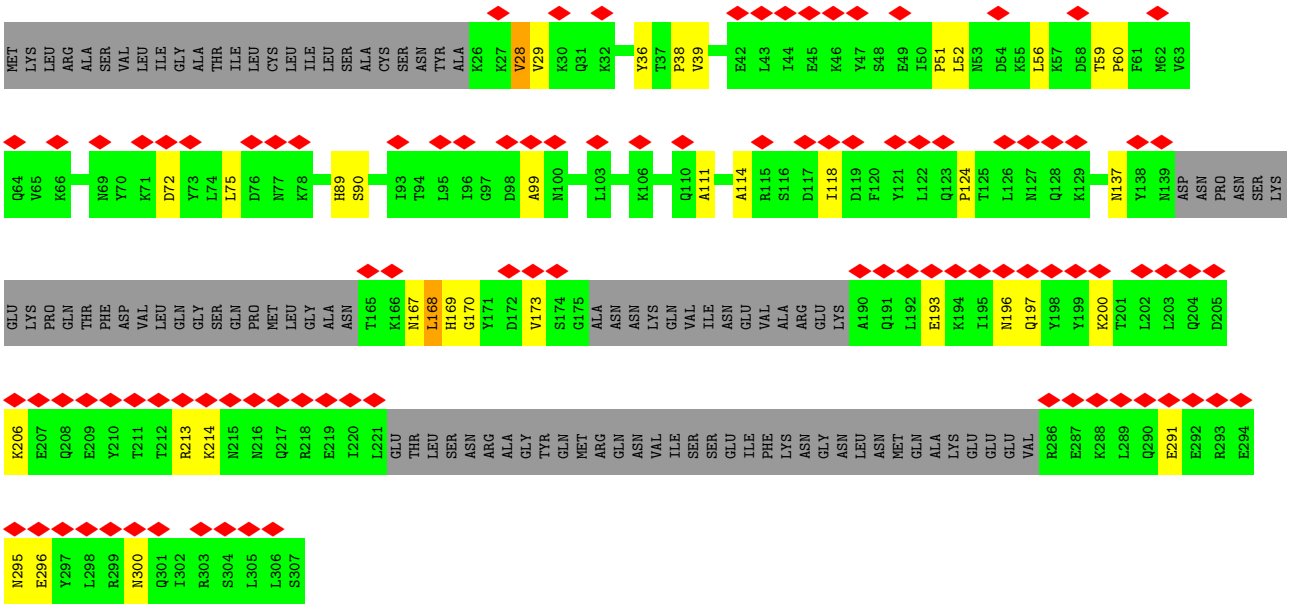
Chain BT:





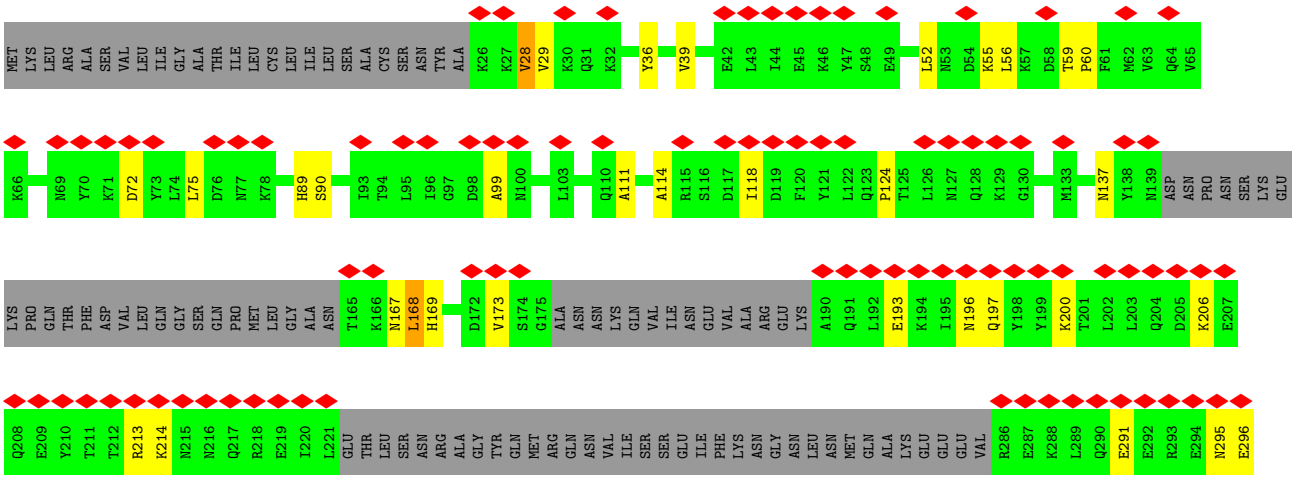
• Molecule 1: Cag pathogenicity island protein

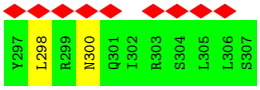
Chain CT:



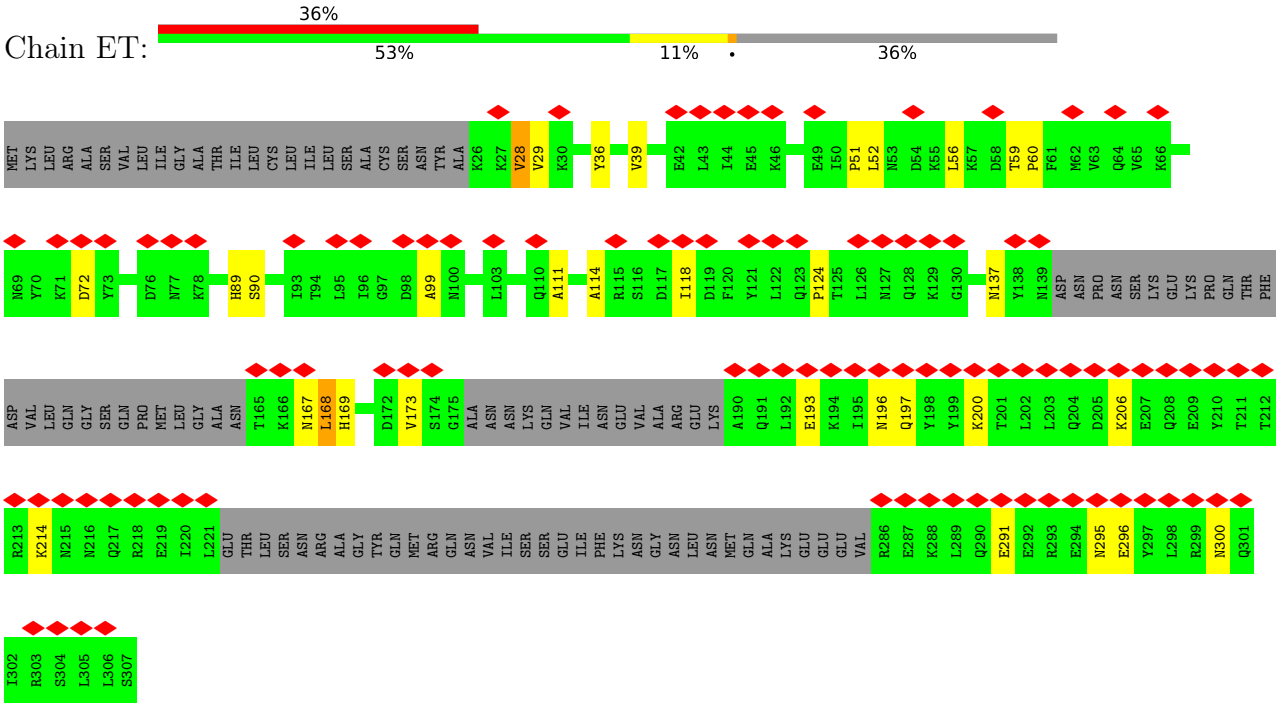
• Molecule 1: Cag pathogenicity island protein

Chain DT:

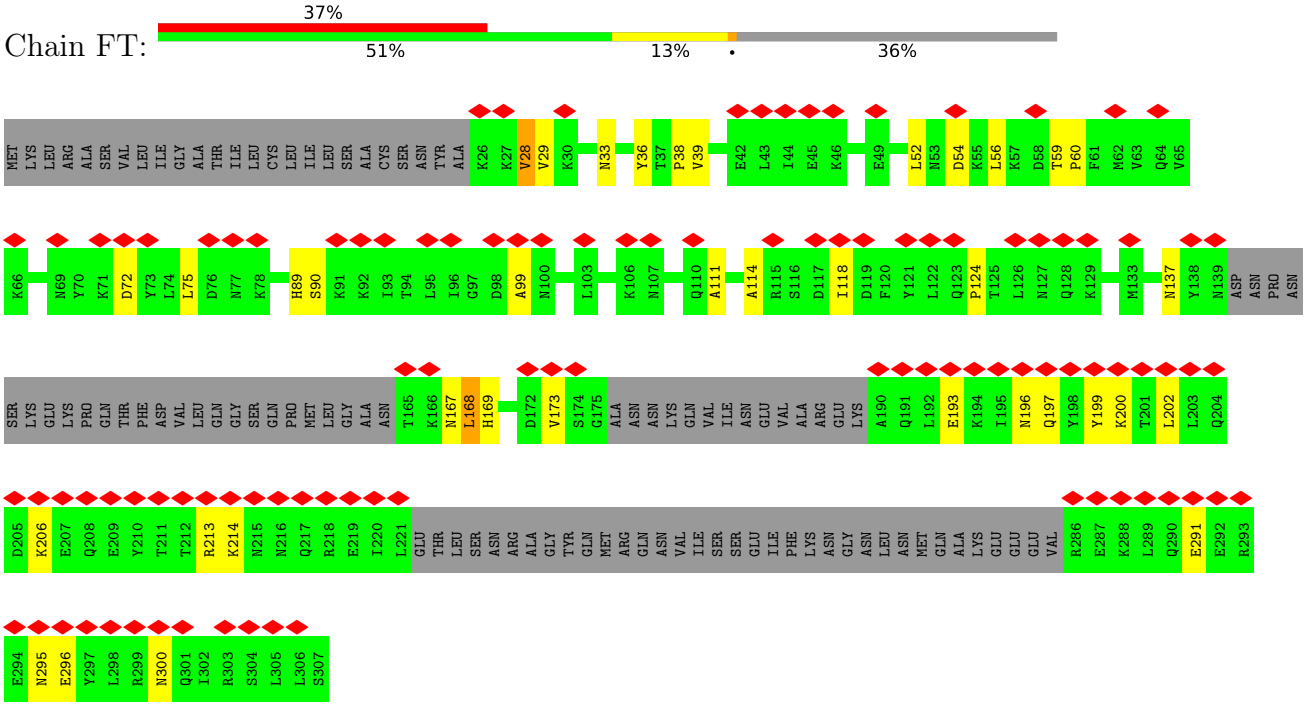




• Molecule 1: Cag pathogenicity island protein

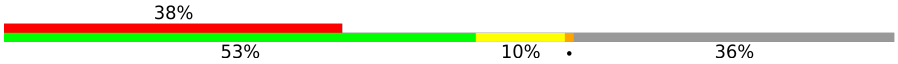


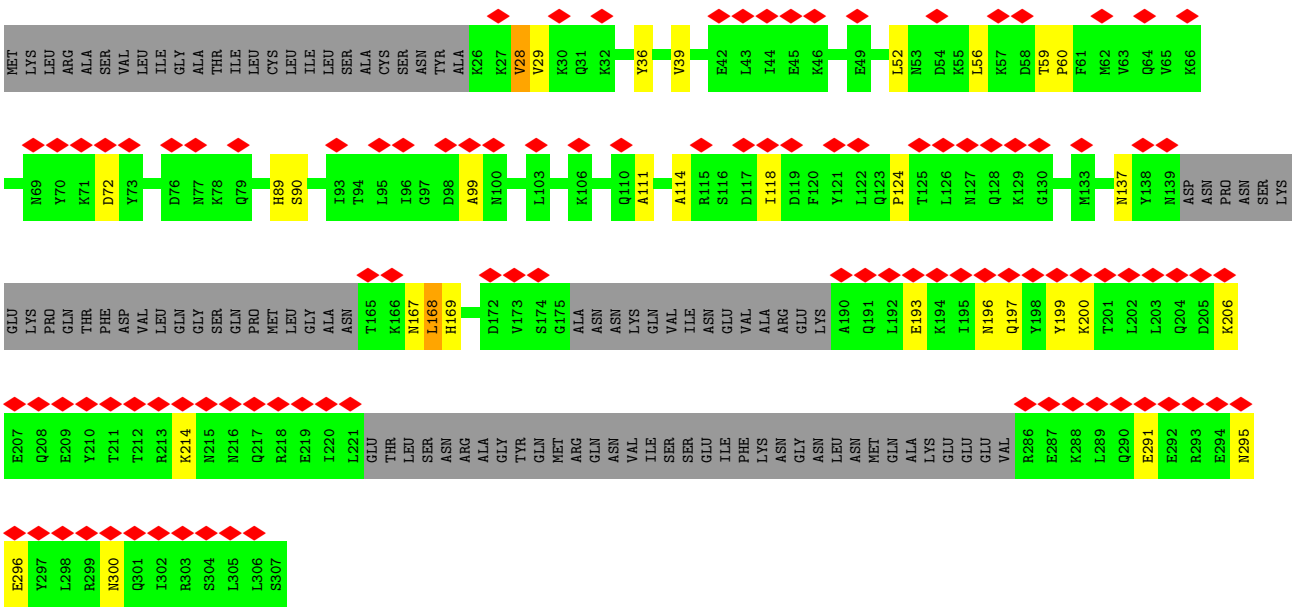
• Molecule 1: Cag pathogenicity island protein



• Molecule 1: Cag pathogenicity island protein

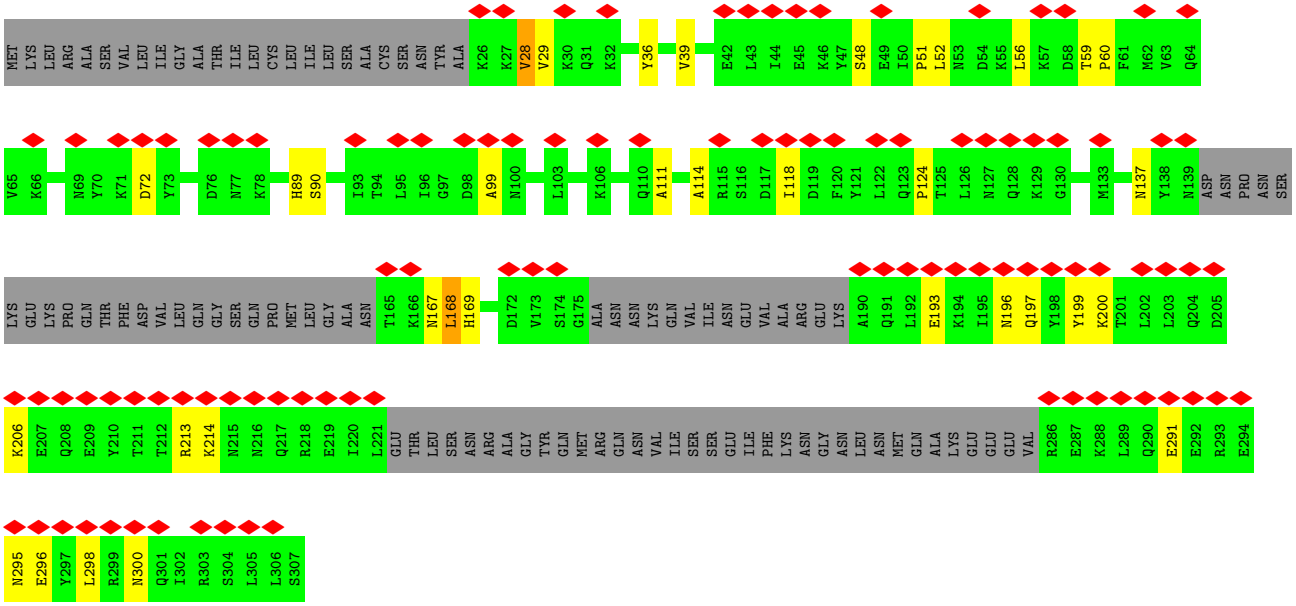
Chain GT:





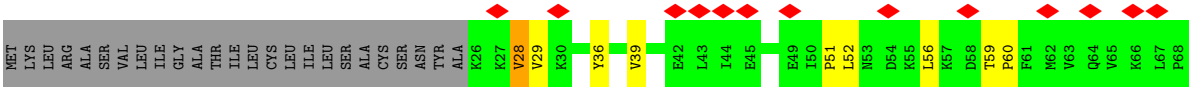
- Molecule 1: Cag pathogenicity island protein

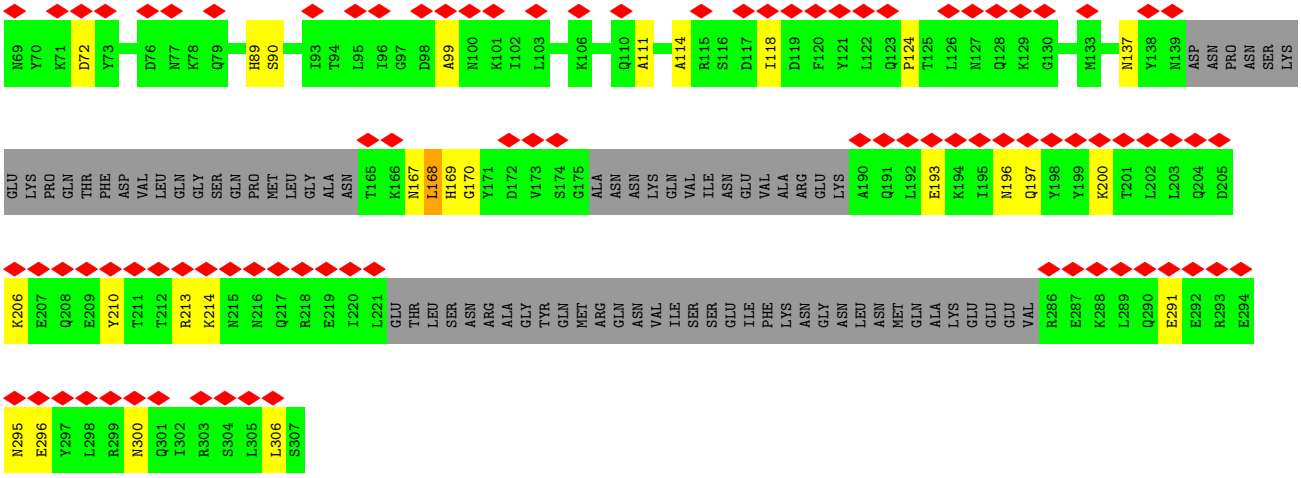
Chain HT:



- Molecule 1: Cag pathogenicity island protein

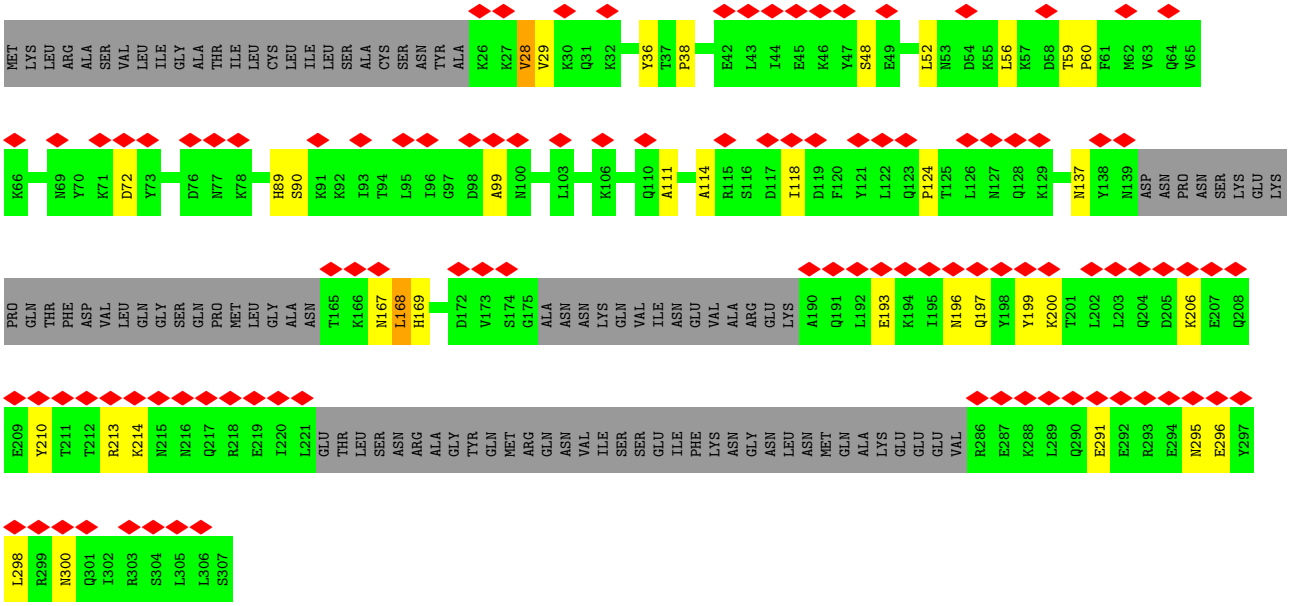
Chain IT:





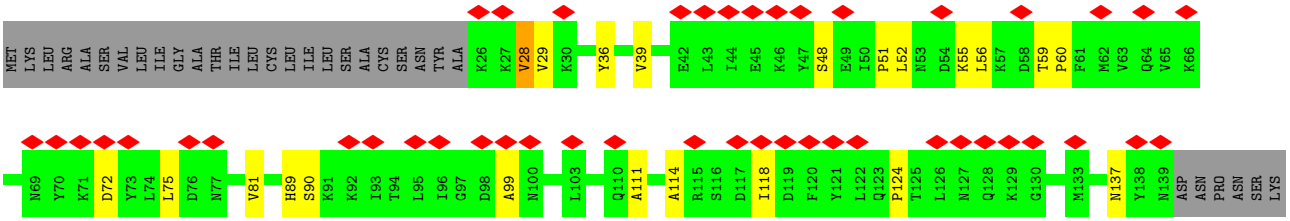
- Molecule 1: Cag pathogenicity island protein

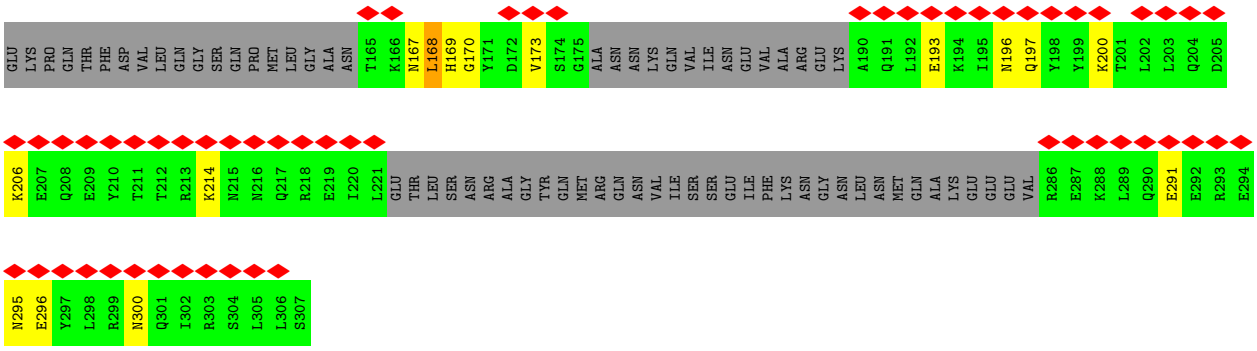
Chain JT:



- Molecule 1: Cag pathogenicity island protein

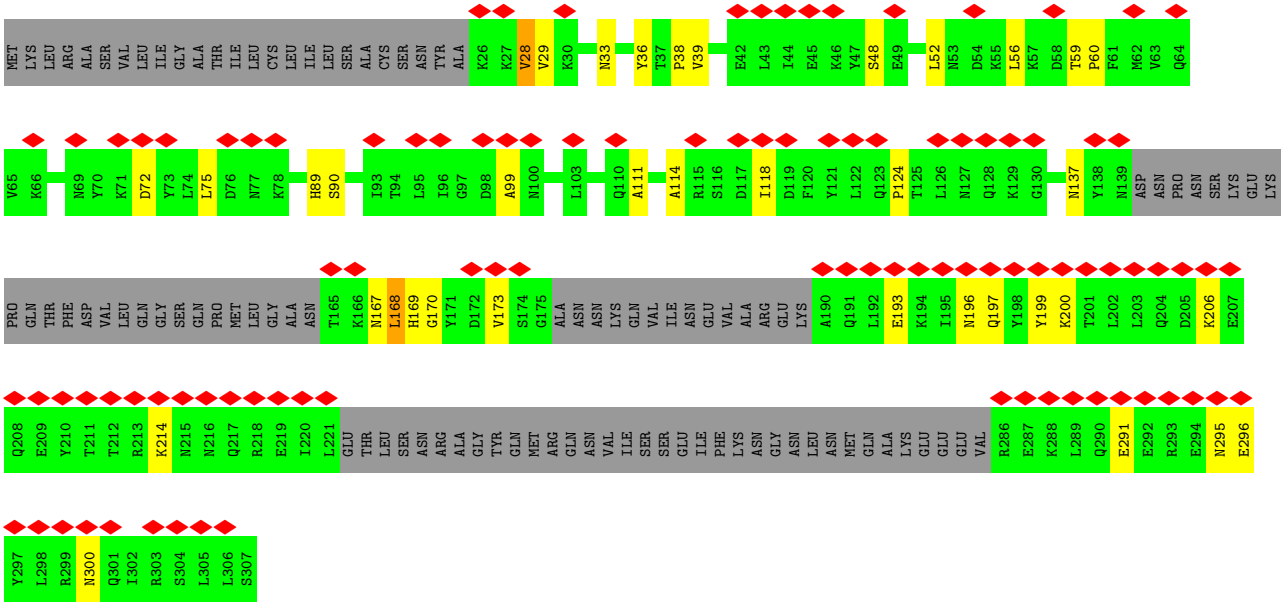
Chain KT:





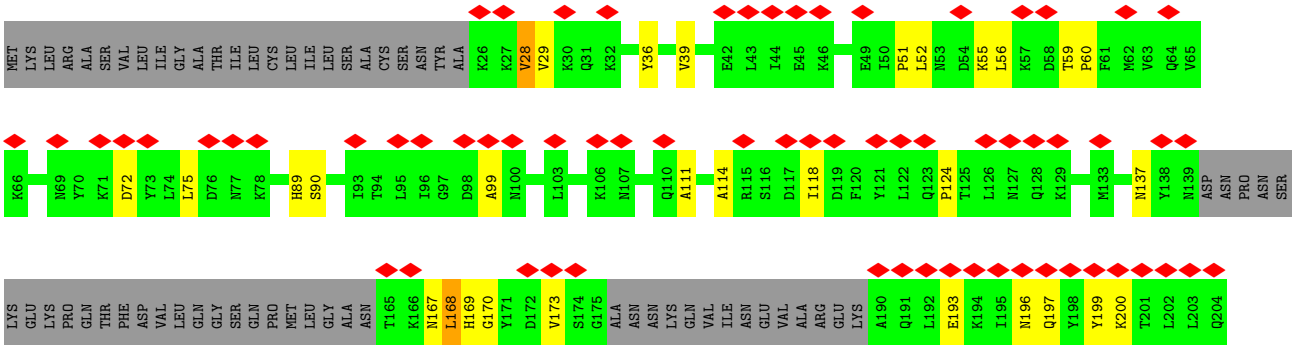
• Molecule 1: Cag pathogenicity island protein

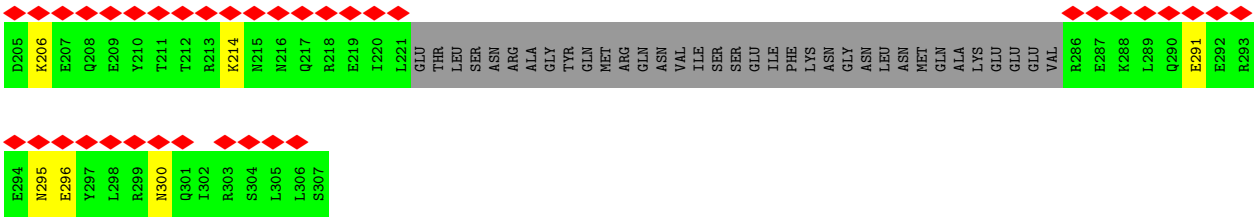
Chain LT:



• Molecule 1: Cag pathogenicity island protein

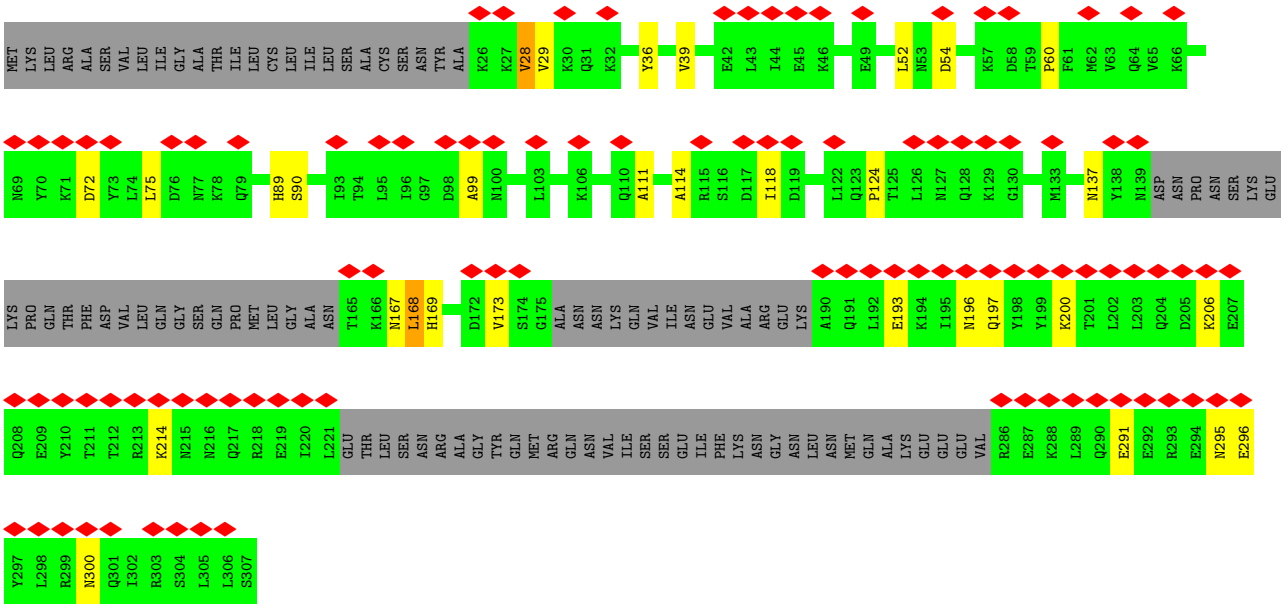
Chain MT:





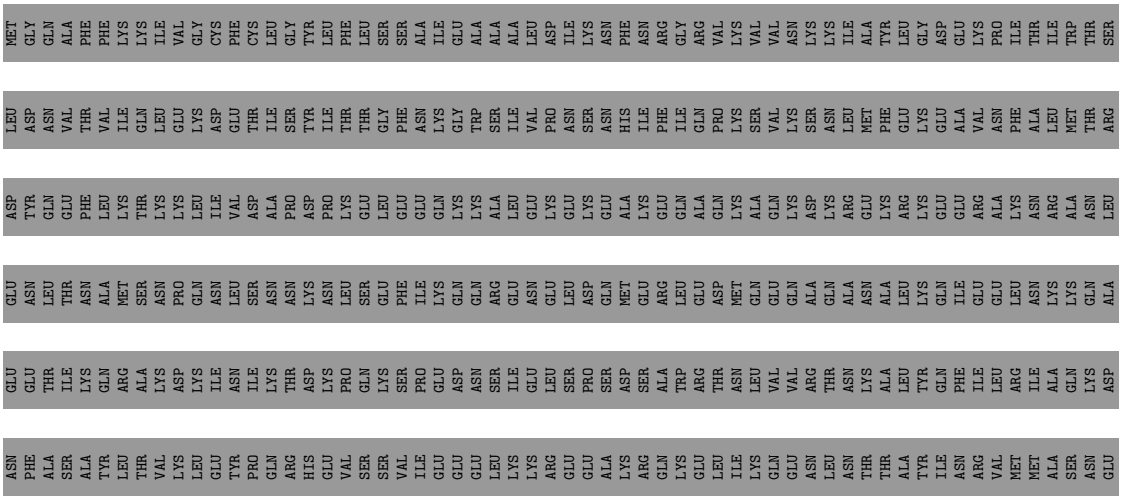
• Molecule 1: Cag pathogenicity island protein

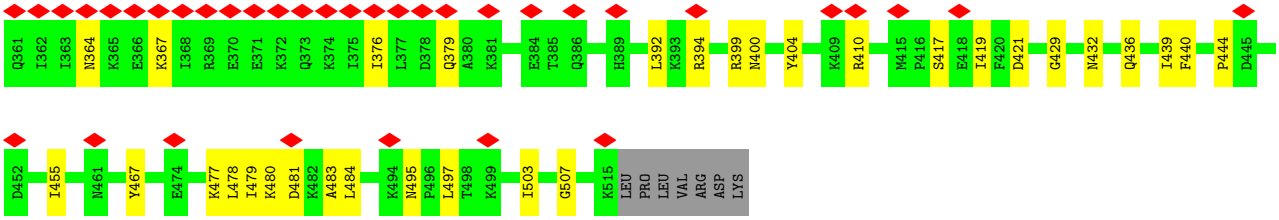
Chain NT:



• Molecule 2: Type IV secretion system apparatus protein CagX

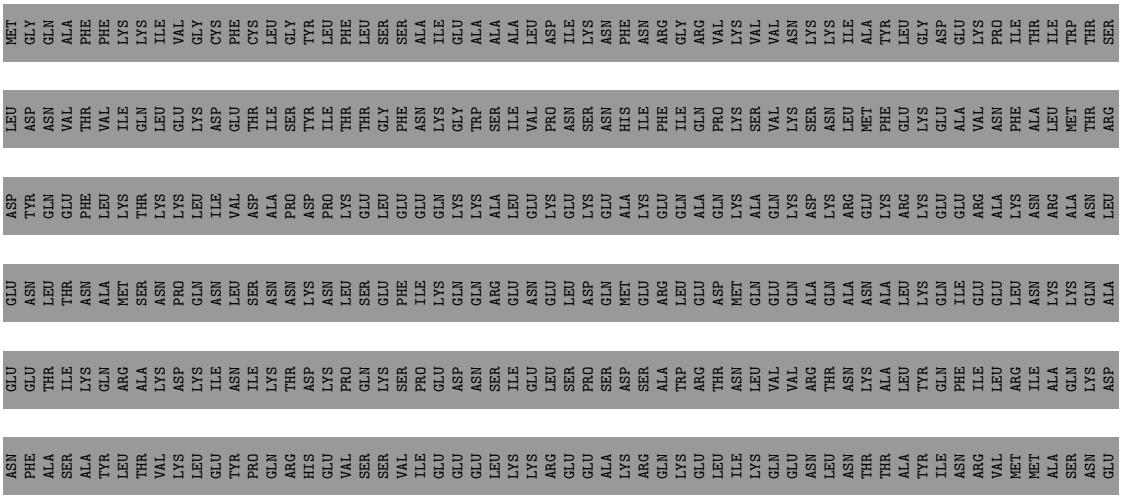
Chain AX:

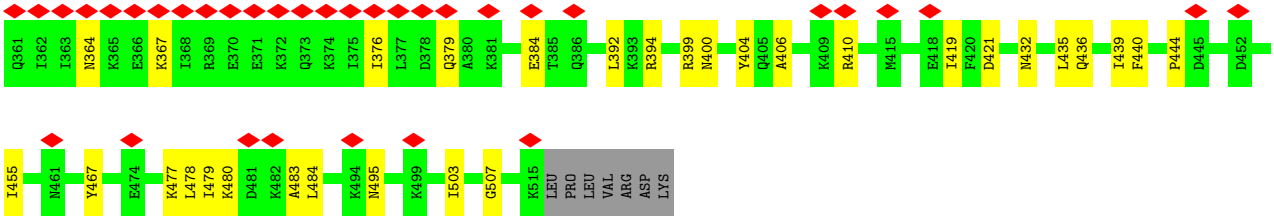




• Molecule 2: Type IV secretion system apparatus protein CagX

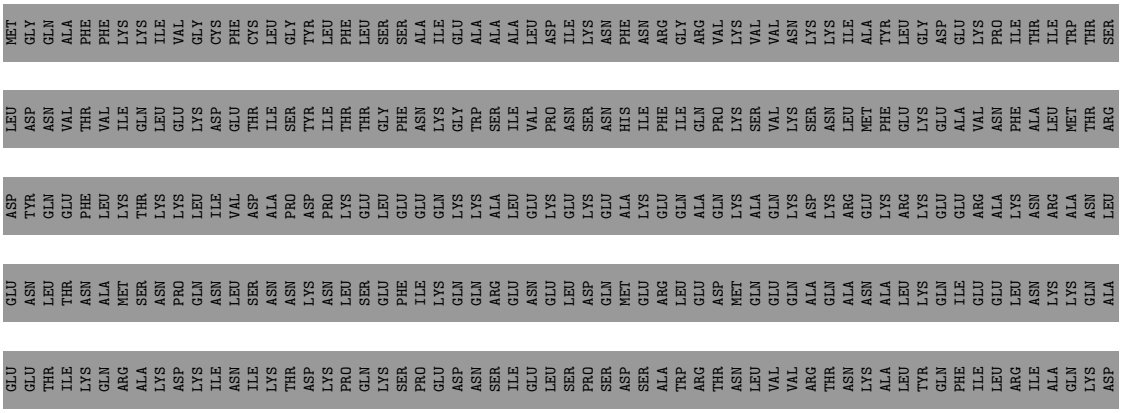
Chain BX:

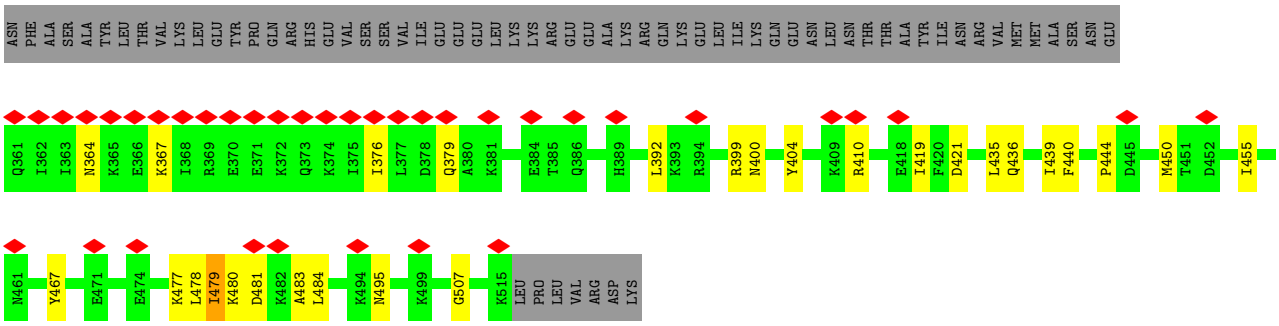




• Molecule 2: Type IV secretion system apparatus protein CagX

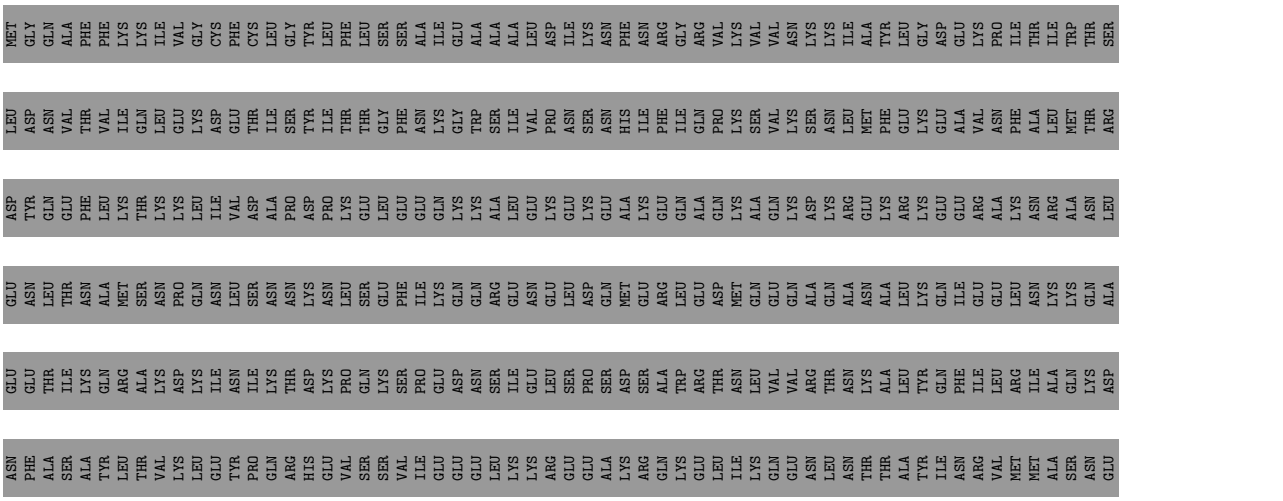
Chain CX:





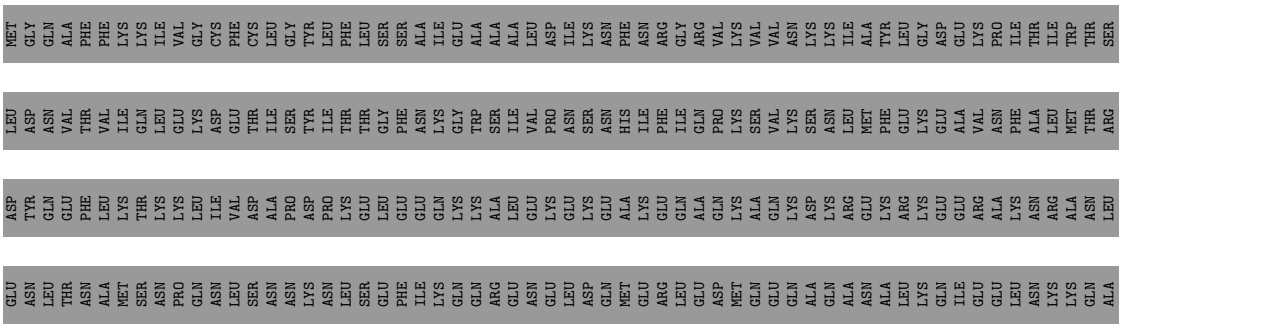
• Molecule 2: Type IV secretion system apparatus protein CagX

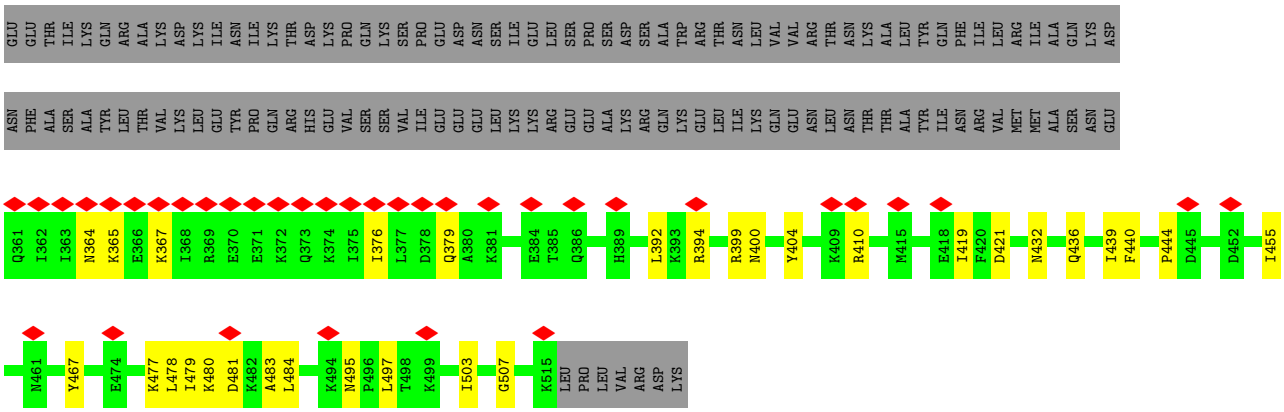
Chain DX: 7% 24% 6% 70%



• Molecule 2: Type IV secretion system apparatus protein CagX

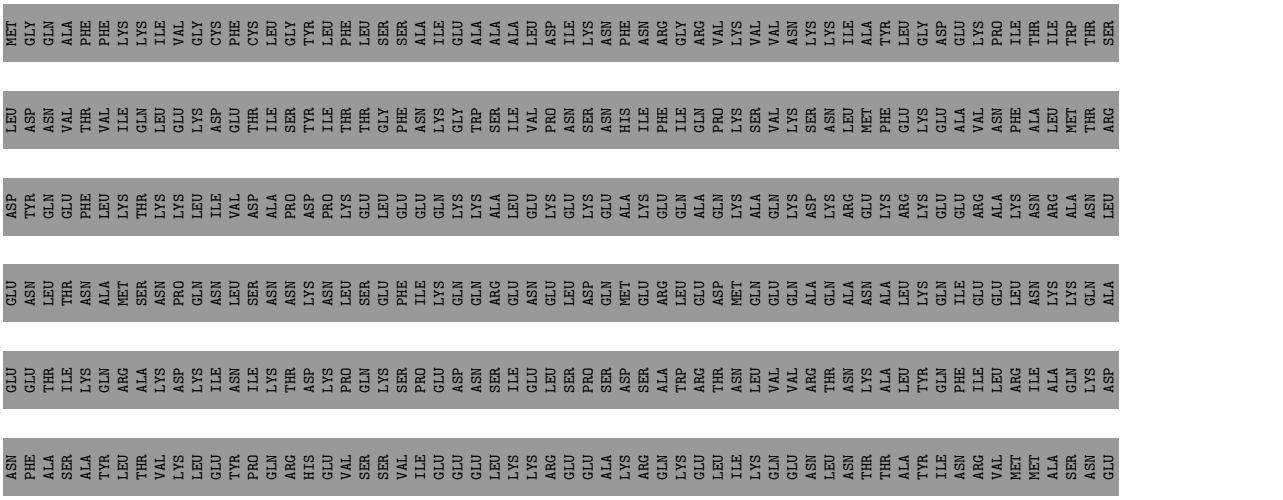
Chain EX: 7% 24% 6% 70%





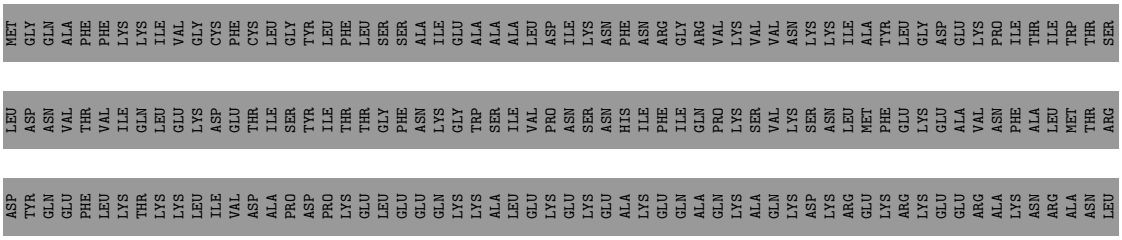
• Molecule 2: Type IV secretion system apparatus protein CagX

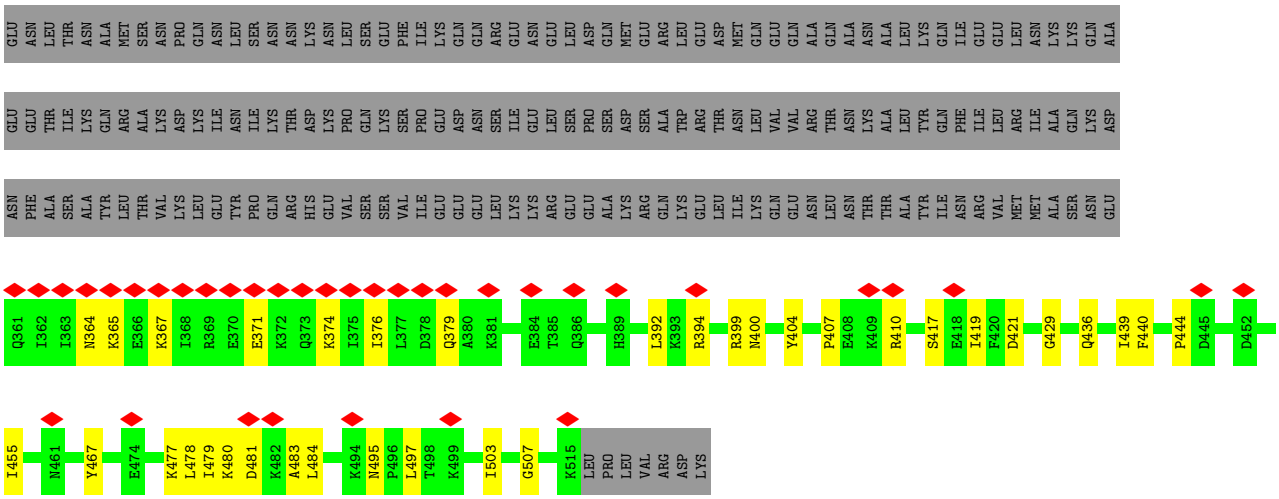
Chain FX:



• Molecule 2: Type IV secretion system apparatus protein CagX

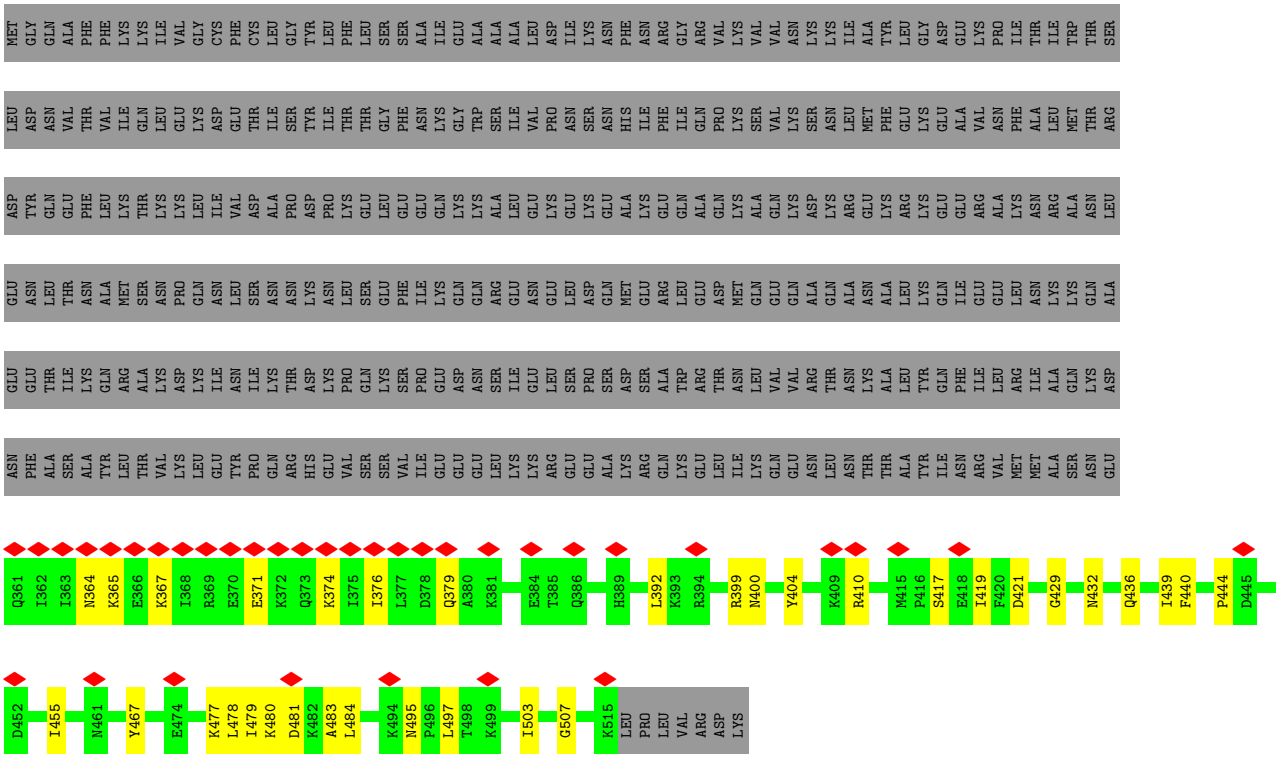
Chain GX:





• Molecule 2: Type IV secretion system apparatus protein CagX

Chain HX:



• Molecule 2: Type IV secretion system apparatus protein CagX

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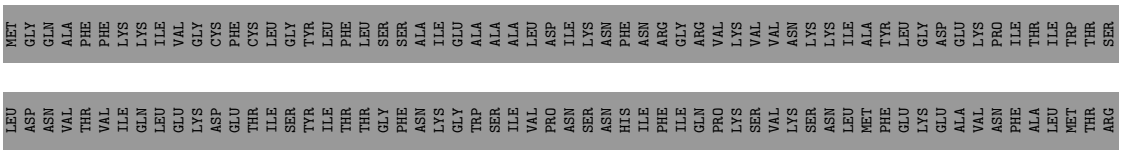
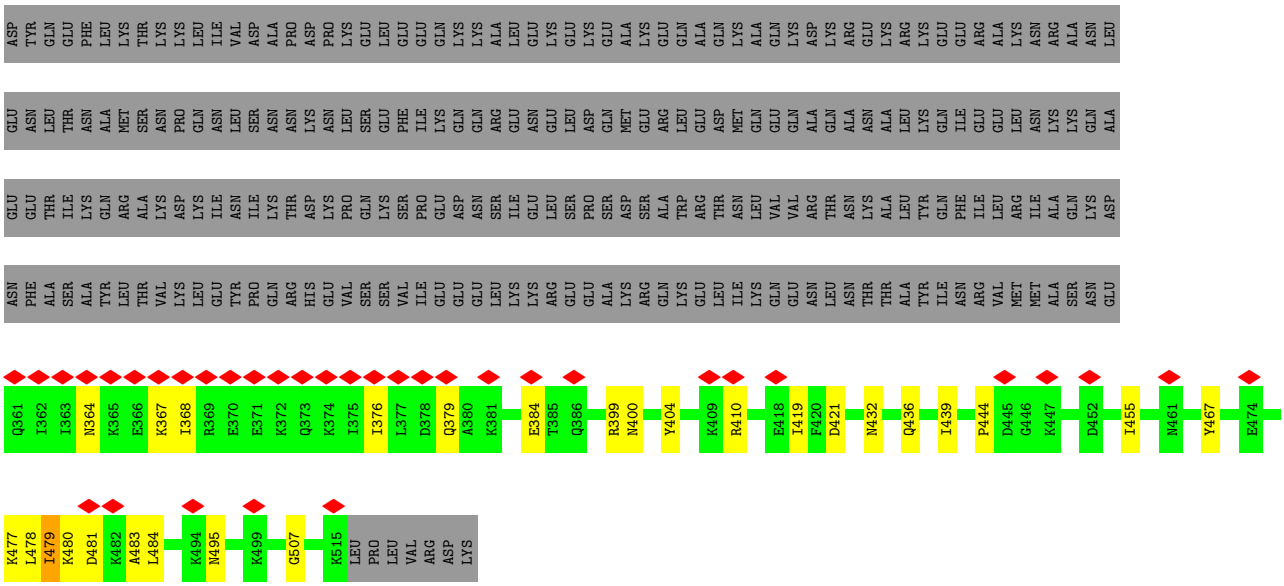
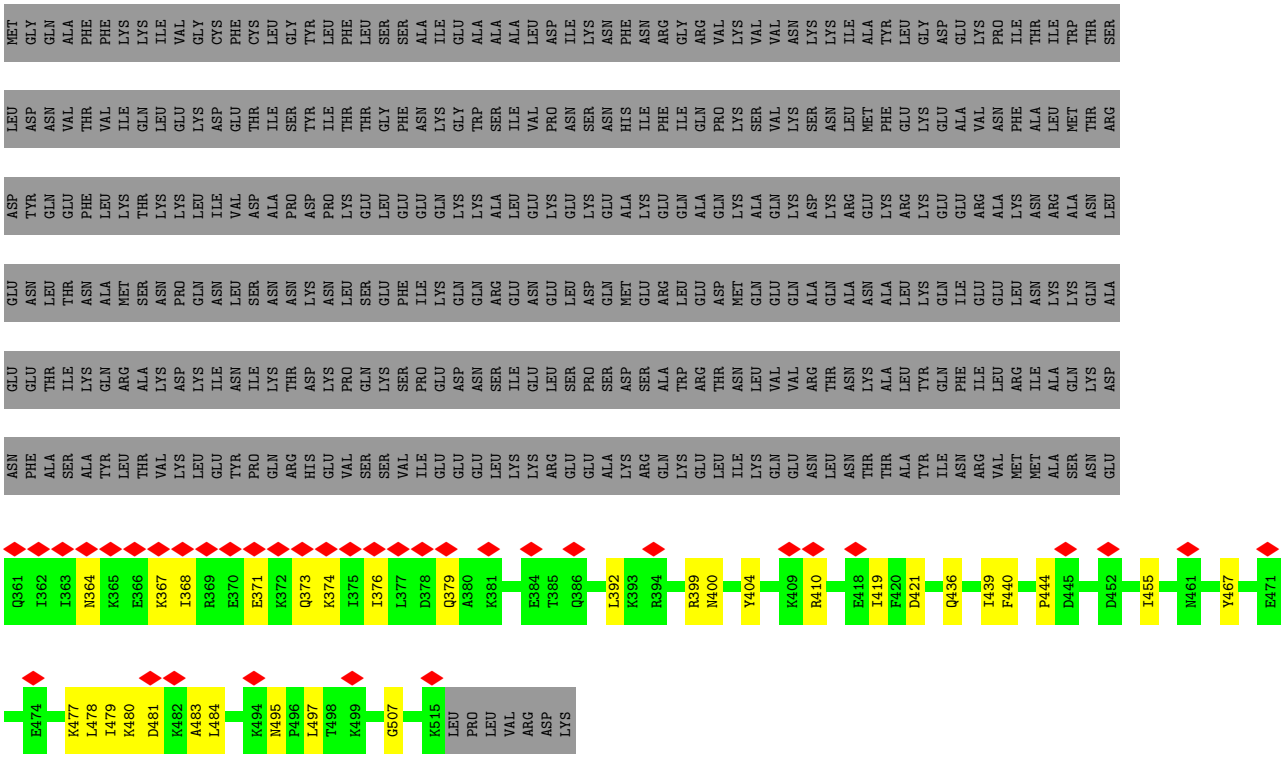




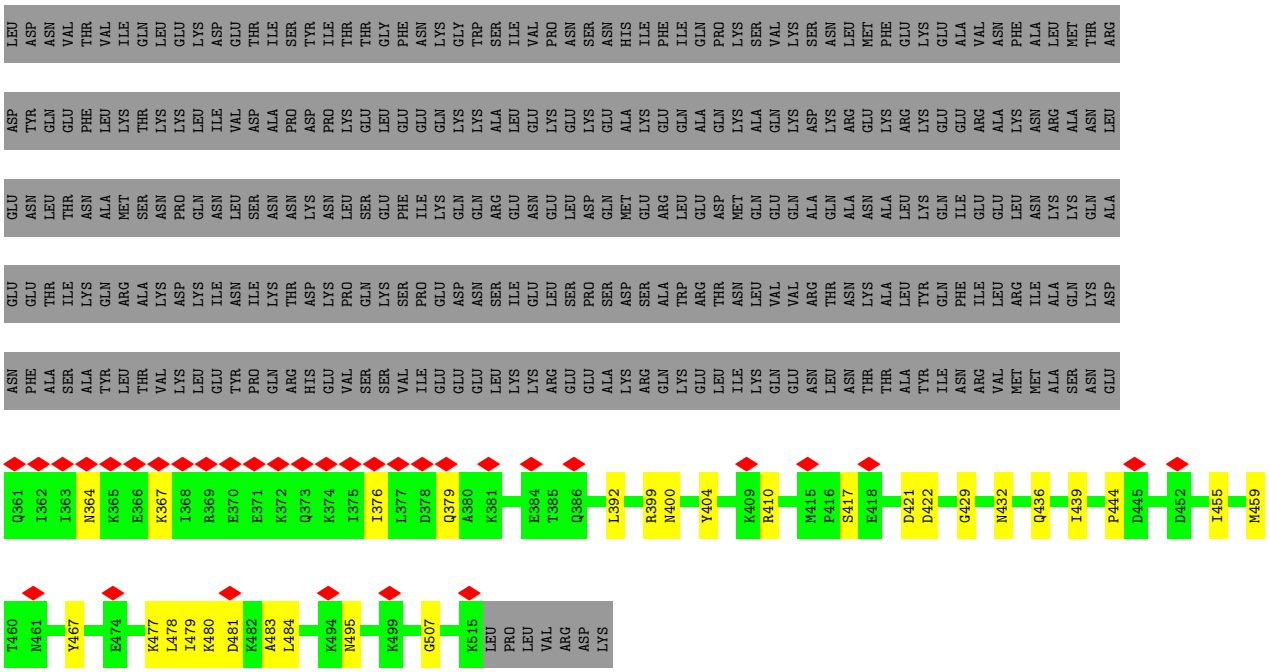
• Molecule 2: Type IV secretion system apparatus protein CagX

Chain JX:



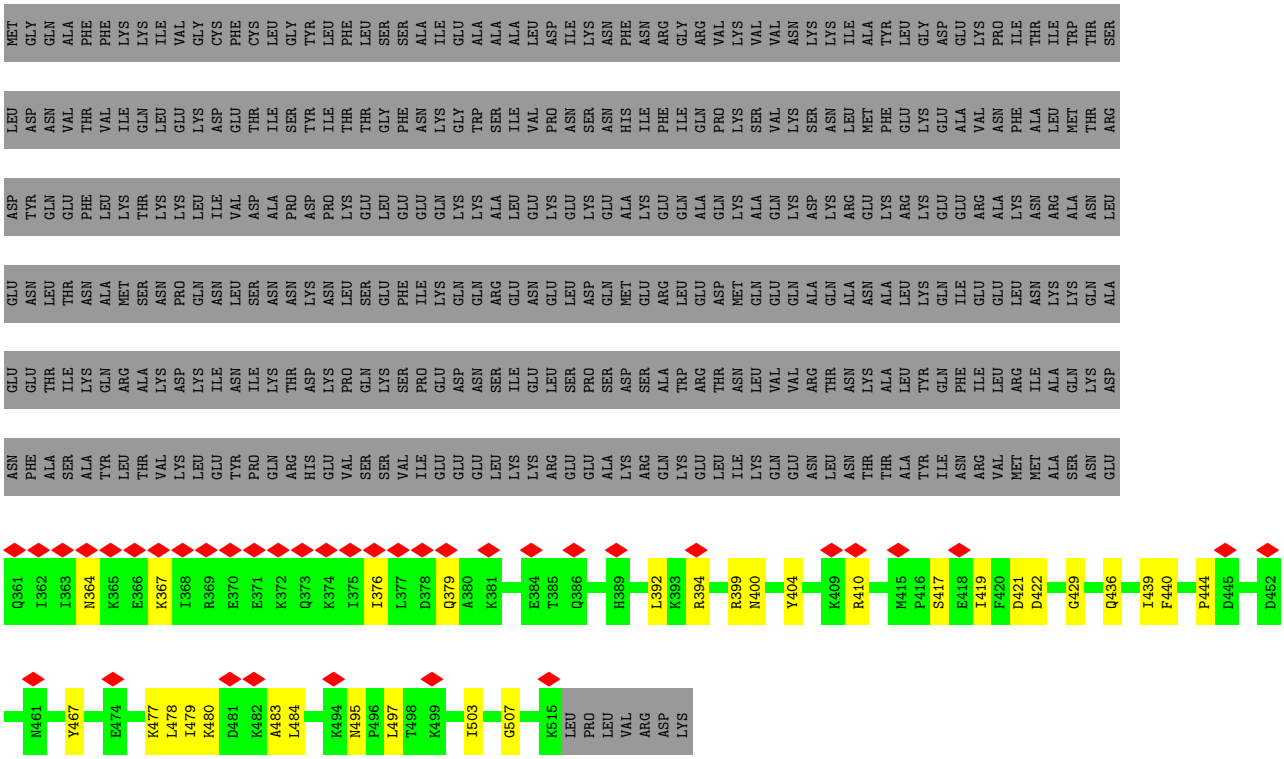
• Molecule 2: Type IV secretion system apparatus protein CagX

Chain KX:



• Molecule 2: Type IV secretion system apparatus protein CagX

Chain LX:



• Molecule 2: Type IV secretion system apparatus protein CagX

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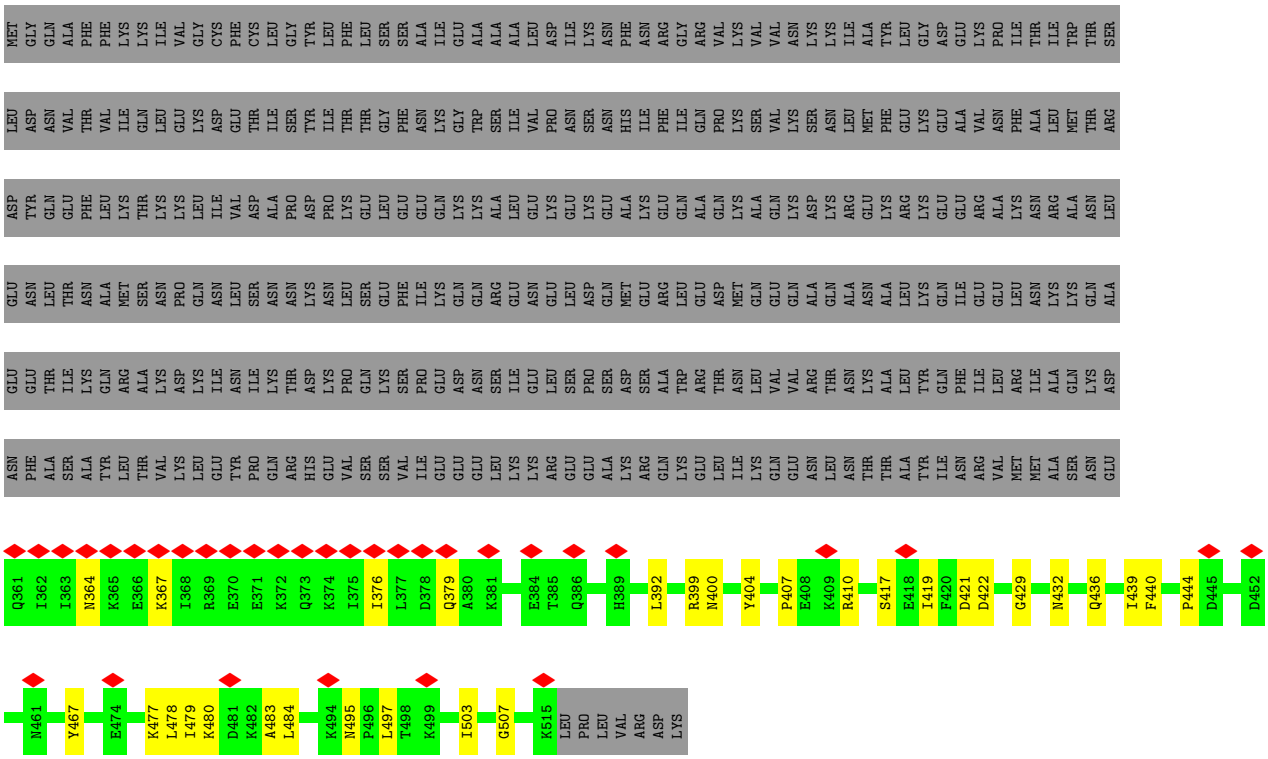

- Molecule 2: Type IV secretion system apparatus protein CagX

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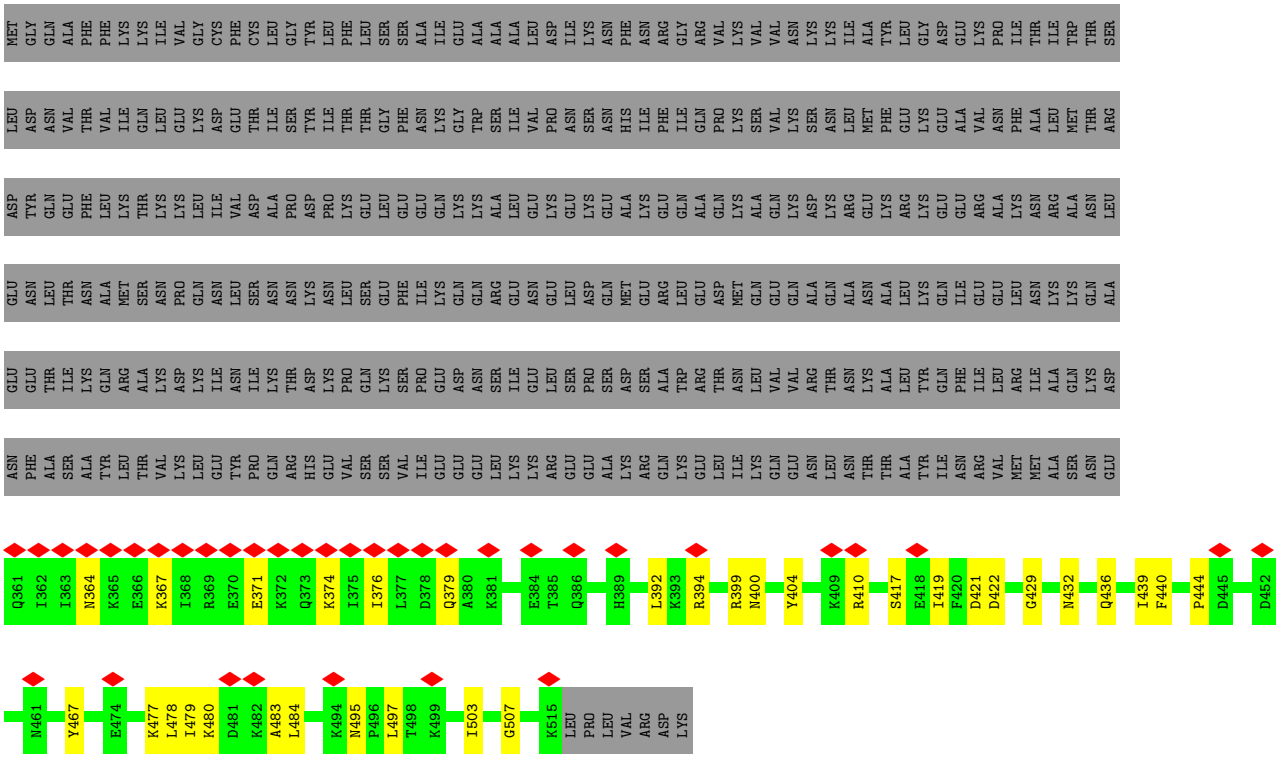


- Molecule 3: Cag pathogenicity island protein (Cag7)

Chain AY:

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

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

ASP GLN GLU ASN ASN GLU TYR GLN GLU GLU THR GLN THR ASP LEU ILE ASP ASP GLU THR SER LYS LYS THR GLN GLN HIS SER PRO GLN ASP LEU SER ASN GLU GLU ALA THR GLU ALA ASN HIS PHE GLU ASN LEU LEU LYS GLU SER LYS GLU SER SER ASP HIS HIS LEU ASP ASN

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

GLU GLU THR GLN GLU LEU ASN GLU GLU ASP ASP GLN GLU ASN ASN GLU TYR GLN GLU GLU THR GLN THR ASP LEU ILE ASP ASP GLU THR SER LYS LYS THR GLN GLN HIS SER PRO GLN ASP LEU SER ASN GLU GLU ALA THR GLU ALA ASN HIS PHE GLU ASN LEU LEU LYS GLU SER

LYS GLU SER SER ASP HIS HIS LEU ASP ASN PRO THR GLU THR GLN THR ASN PHE ASP GLY ASP LYS SER GLU GLU ILE THR ASP ASP SER ASN ASP GLN GLU ILE ILE LYS GLY SER LYS LYS LYS TYR ILE ILE GLY GLY ILE VAL VAL ALA VAL LEU ILE VAL ILE ILE LEU PHE SER

ARG SER ILE PHE HIS TYR PHE MET PRO LEU GLU ASP LYS SER SER ARG PHE SER LYS ASP ARG ASN LEU TYR VAL ASN ASP GLU ILE GLN ILE ARG GLN GLU TYR ASN ARG LEU LEU LYS GLU ARG ASN GLU LYS GLY ASN MET ILE ASP LYS ASN LEU PHE PHE ASN ASP ASP PRO ASN

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

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

PHE ARG LYS GLU LEU GLU LEU GLN LYS GLU LEU GLN GLU TYR LYS ASP CYS ILE LYS ASN ALA LYS THR GLU ALA GLU LYS ASN LYS CYS LEU LYS GLY LEU SER LYS GLU ALA ILE GLU ARG LEU LYS GLN GLN ALA LEU ASP CYS LEU LYS ASN ALA LYS THR ASP GLU GLU ARG ASN

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

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

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

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

CYS VAL SER GLN ALA LYS THR GLU ALA GLU LYS LYS GLU CYS GLU LYS LEU LEU THR PRO GLU ALA ARG LYS LEU LEU GLU GLU ALA LYS LYS SER VAL LYS ALA TYR LEU ASP CYS VAL SER GLN ALA LYS THR GLU ALA GLU LYS LYS GLU CYS GLU LYS LEU LEU THR PRO GLU ALA

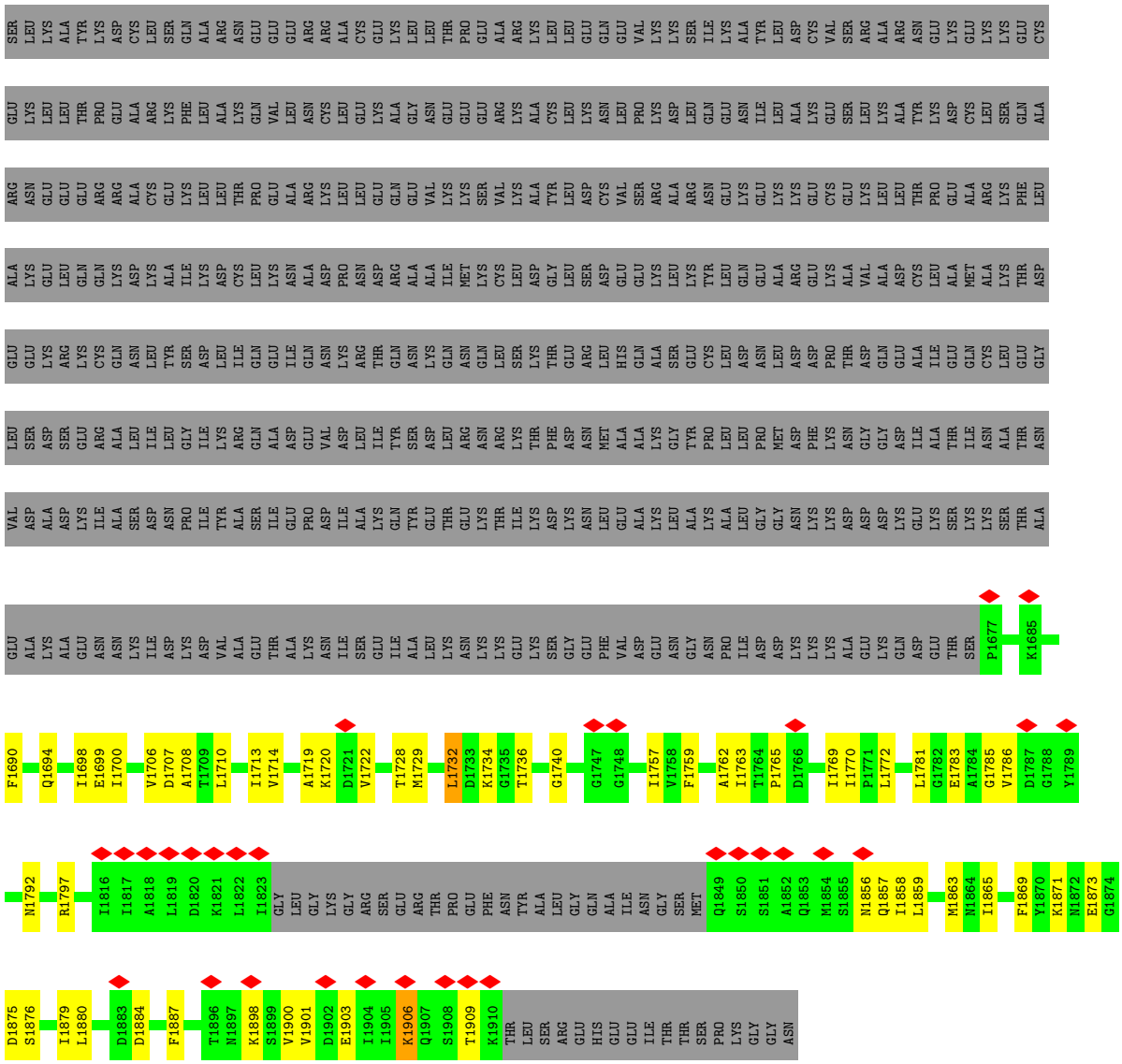
ARG LYS LEU LEU GLU GLU UNK ALA LYS GLU SER VAL LYS ALA TYR LEU ASP CYS VAL SER GLN ALA LYS ASN GLU ALA GLU LYS LYS GLU CYS GLU LYS LEU LEU THR LEU GLU SER LYS LYS LEU GLU GLU ALA LYS LYS SER VAL LYS ALA TYR LEU ASP CYS VAL SER GLN ALA

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

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

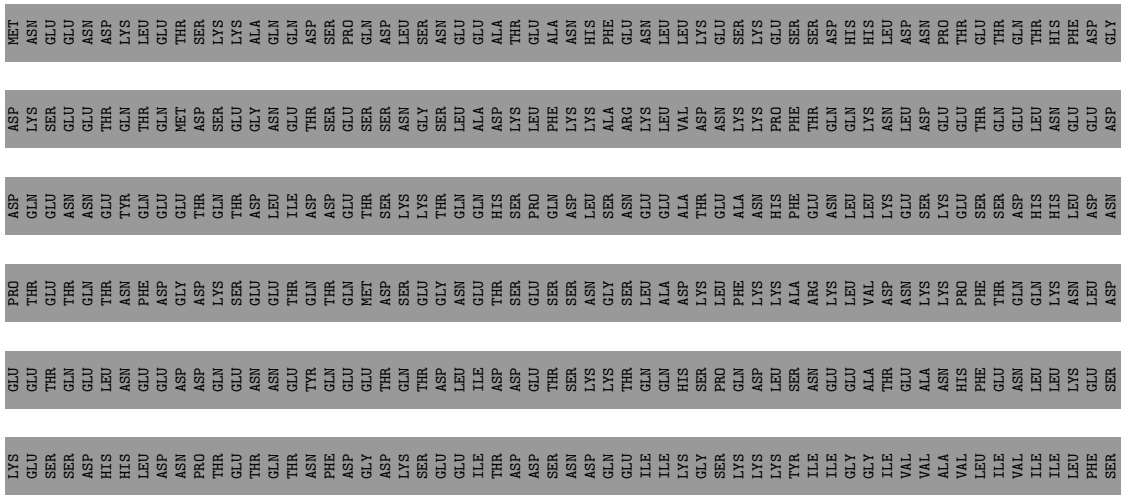
THR PRO GLU ALA ARG LYS LEU LEU GLU GLU ALA LYS GLU SER VAL LYS ALA TYR LYS ASP CYS VAL SER LYS ALA ARG ASN GLU LYS GLU LYS LYS GLU CYS GLU LYS LEU LEU THR PRO GLU ALA LYS LYS LEU LEU GLU GLN GLN VAL LEU ASP CYS LEU LYS ASN ALA LYS THR GLU

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



• Molecule 3: Cag pathogenicity island protein (Cag7)

Chain BY: 8% 89%



ARG SER ILE PHE HIS TYR PHE MET PRO LEU GLU ASP LYS SER SER ARG PHE SER LYS ASP ARG ASN LEU TYR VAL ASN ASP GLU ILE GLN ILE ARG GLN GLU TYR ASN ARG LEU LEU LYS GLU ARG ASN GLU LYS GLY ASN MET ILE ASP LYS ASN LEU PHE PHE ASN ASP ASP PRO ASN

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

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

PHE ARG LYS GLU LEU GLU LEU GLN LYS GLU LEU GLN GLU TYR LYS ASP CYS ILE LYS ASN ALA LYS THR GLU ALA GLU LYS ASN LYS CYS LEU LYS GLY LEU SER LYS GLU ALA ILE GLU ARG LEU LYS GLN GLN ALA LEU ASP CYS LEU LYS ASN ALA LYS THR ASP GLU GLU ARG ASN

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

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

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

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

CYS VAL SER GLN ALA LYS THR GLU ALA GLU LYS LYS GLU CYS GLU LYS LEU LEU THR PRO GLU ALA ARG LYS LEU LEU GLU GLU ALA LYS LYS SER VAL LYS ALA TYR LEU ASP CYS VAL SER GLN ALA LYS THR GLU ALA GLU LYS LYS GLU CYS GLU LYS LEU LEU THR PRO GLU ALA

ARG LYS LEU LEU GLU GLU UNK ALA LYS GLU SER VAL LYS ALA TYR LEU ASP CYS VAL SER GLN ALA LYS ASN GLU ALA GLU LYS LYS GLU CYS GLU LYS LEU LEU THR LEU GLU SER LYS LYS LYS LEU GLU GLU ALA LYS LYS SER VAL LYS ALA TYR LEU ASP CYS VAL SER GLN ALA

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

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

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

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

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

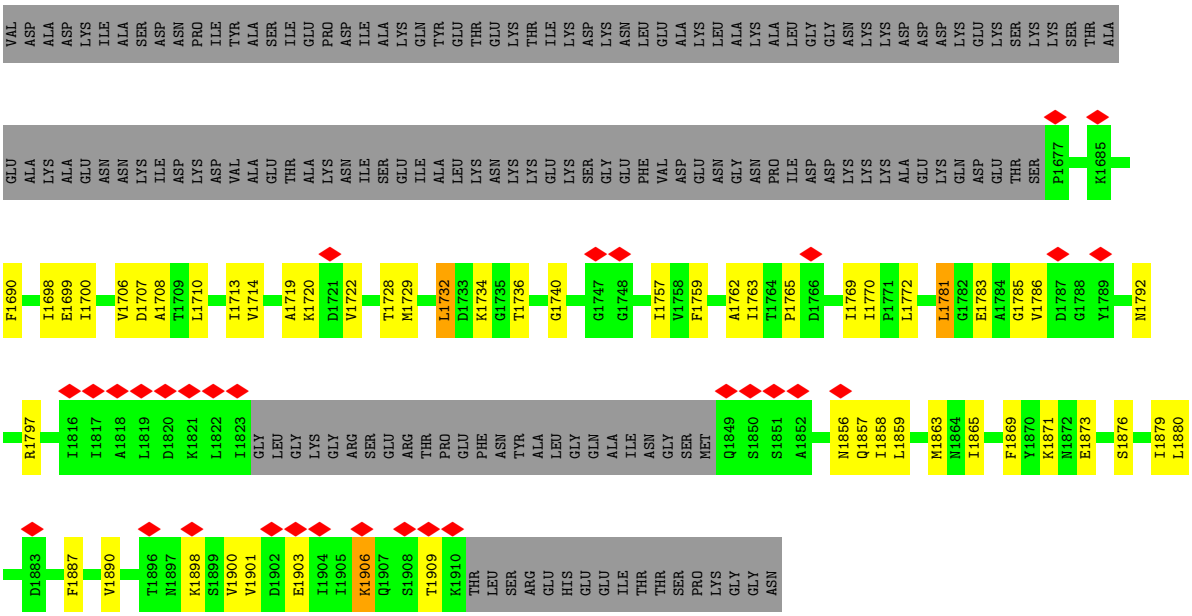
GLU LYS LEU LEU THR PRO GLU ALA ARG LYS PHE LEU ALA LYS GLN VAL LEU ASN CYS LEU GLU LYS ALA GLY ASN GLU GLU ARG LYS ALA CYS LEU LYS ASN LEU PRO LYS ASP LEU GLN GLU ASN ILE LEU ALA LYS GLU SER LEU LYS ALA TYR LYS ASP CYS LEU SER GLN ALA

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

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

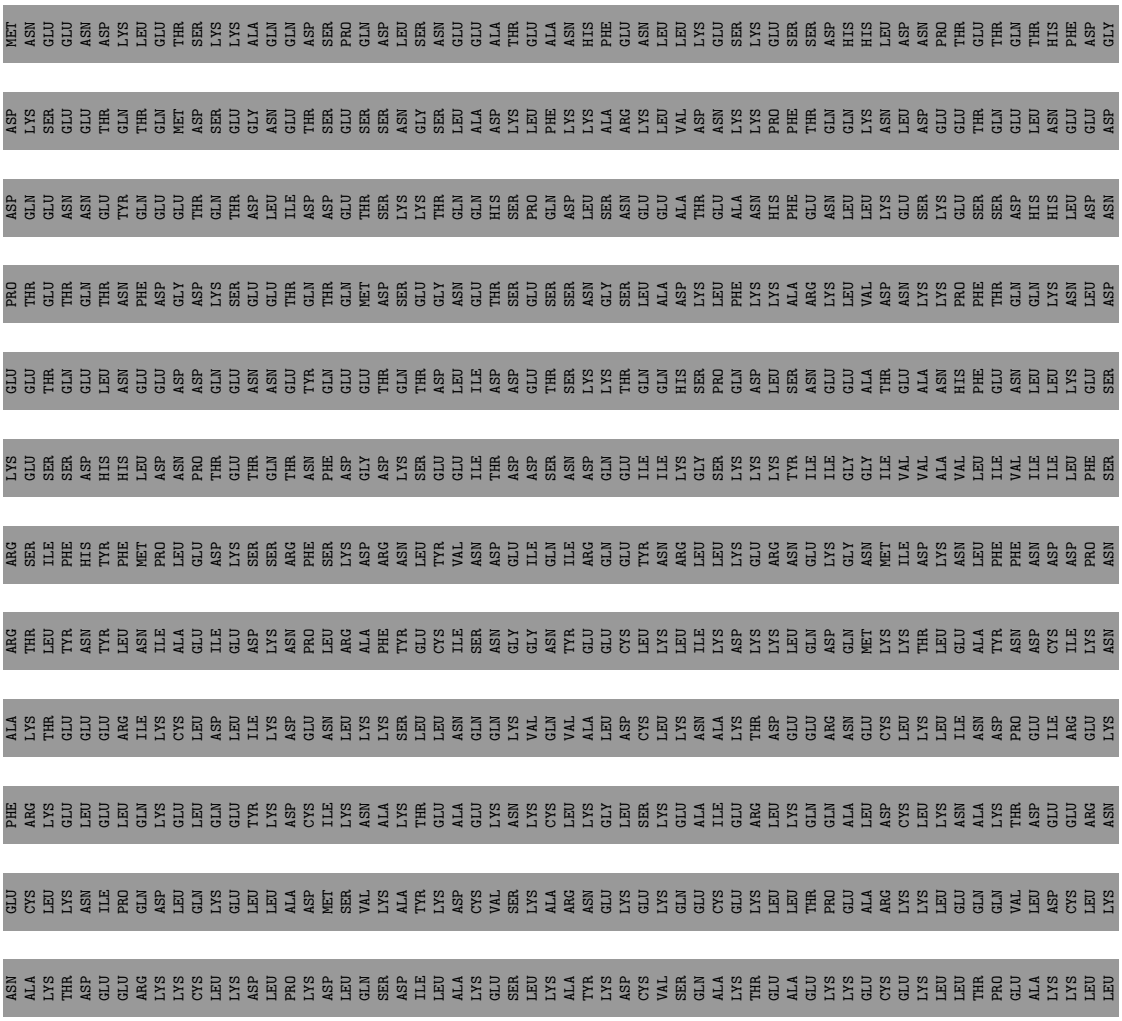
GLU GLU LYS ARG LYS CYS GLN ASN LEU TYR SER ASP LEU ILE GLN GLU ILE GLN ASN LYS ARG THR GLN ASN LYS GLN ASN GLN LEU SER LYS THR GLU ARG LEU HIS GLN ALA SER GLU CYS LEU ASP ASN LEU ASP ASP PRO THR ASP GLN GLU ALA ILE GLU GLN CYS LEU GLN GLY

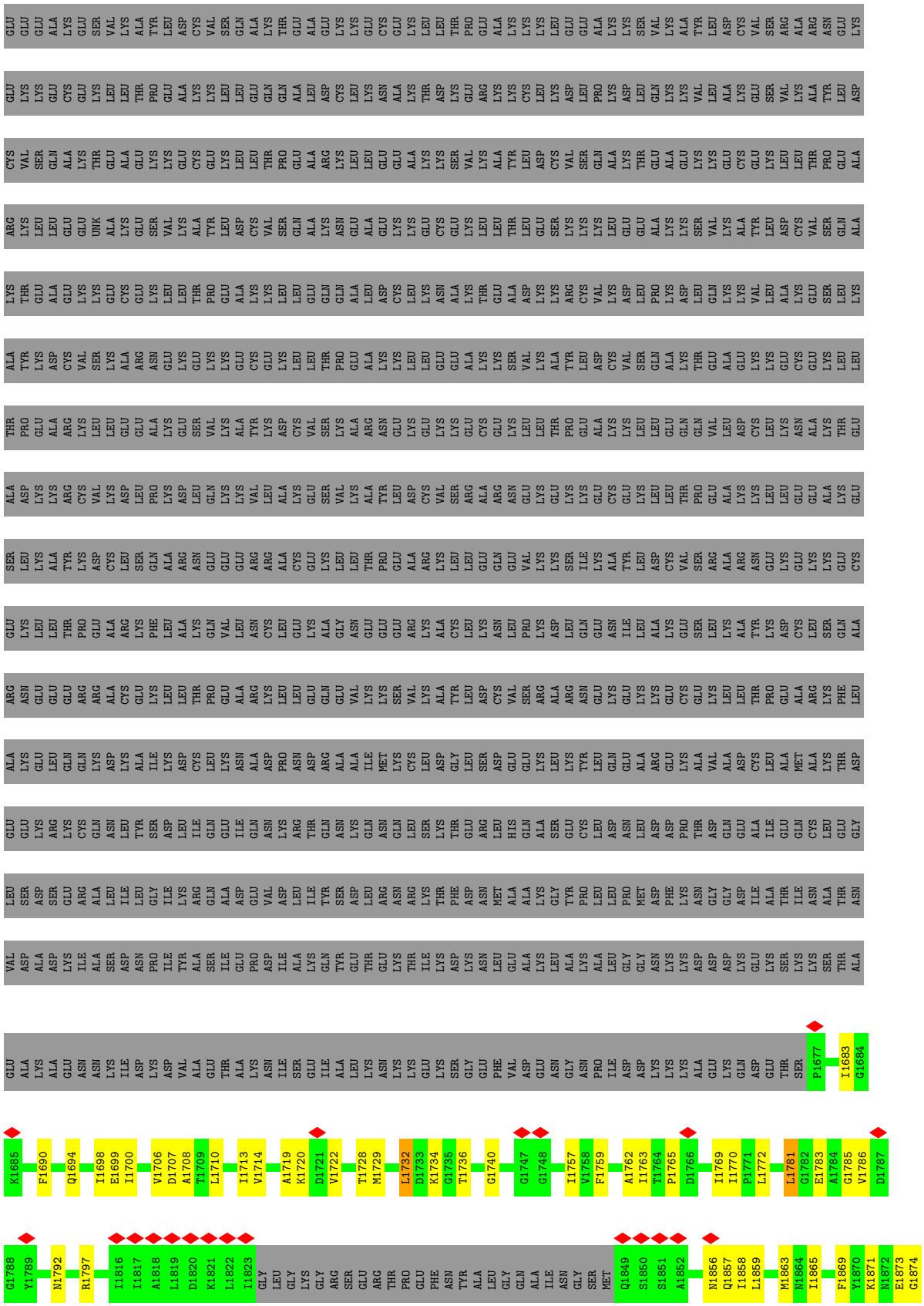
LEU SER ASP SER GLU ARG ALA LEU ILE LEU GLY ILE LYS ARG GLN ALA ASP GLU VAL ASP LEU ILE TYR SER ASP LEU ARG ASN ARG LYS THR PHE ASP ASN MET ALA ALA LYS GLY TYR PRO LEU LEU PRO MET ASP PHE LYS ASN GLY GLY ASP ILE ALA THR ILE ASN ALA THR ASN



• Molecule 3: Cag pathogenicity island protein (Cag7)

Chain CY: 8% 89%

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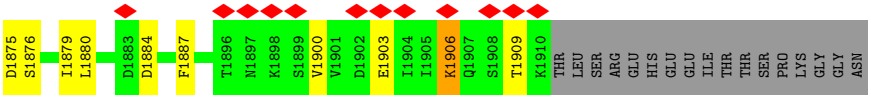
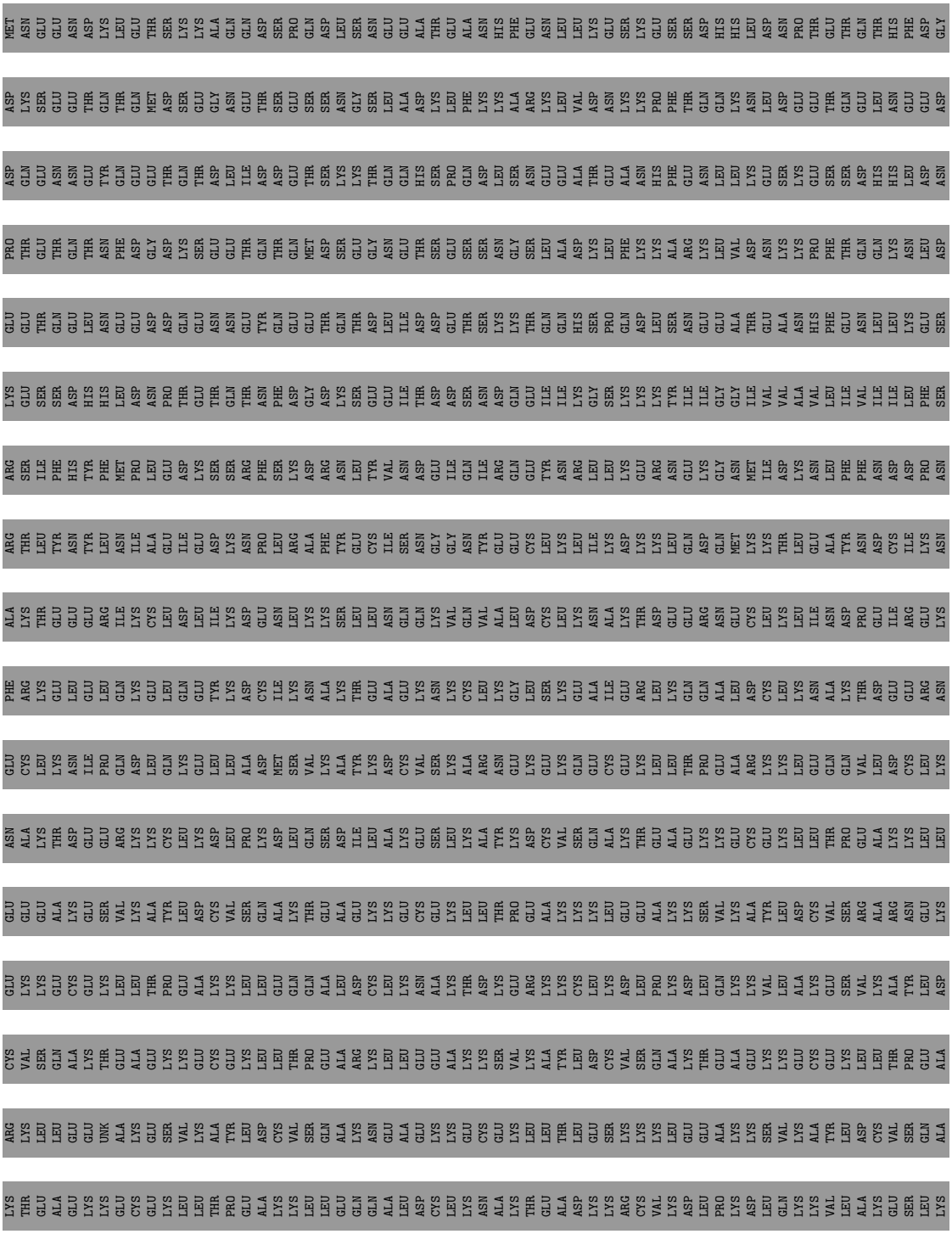
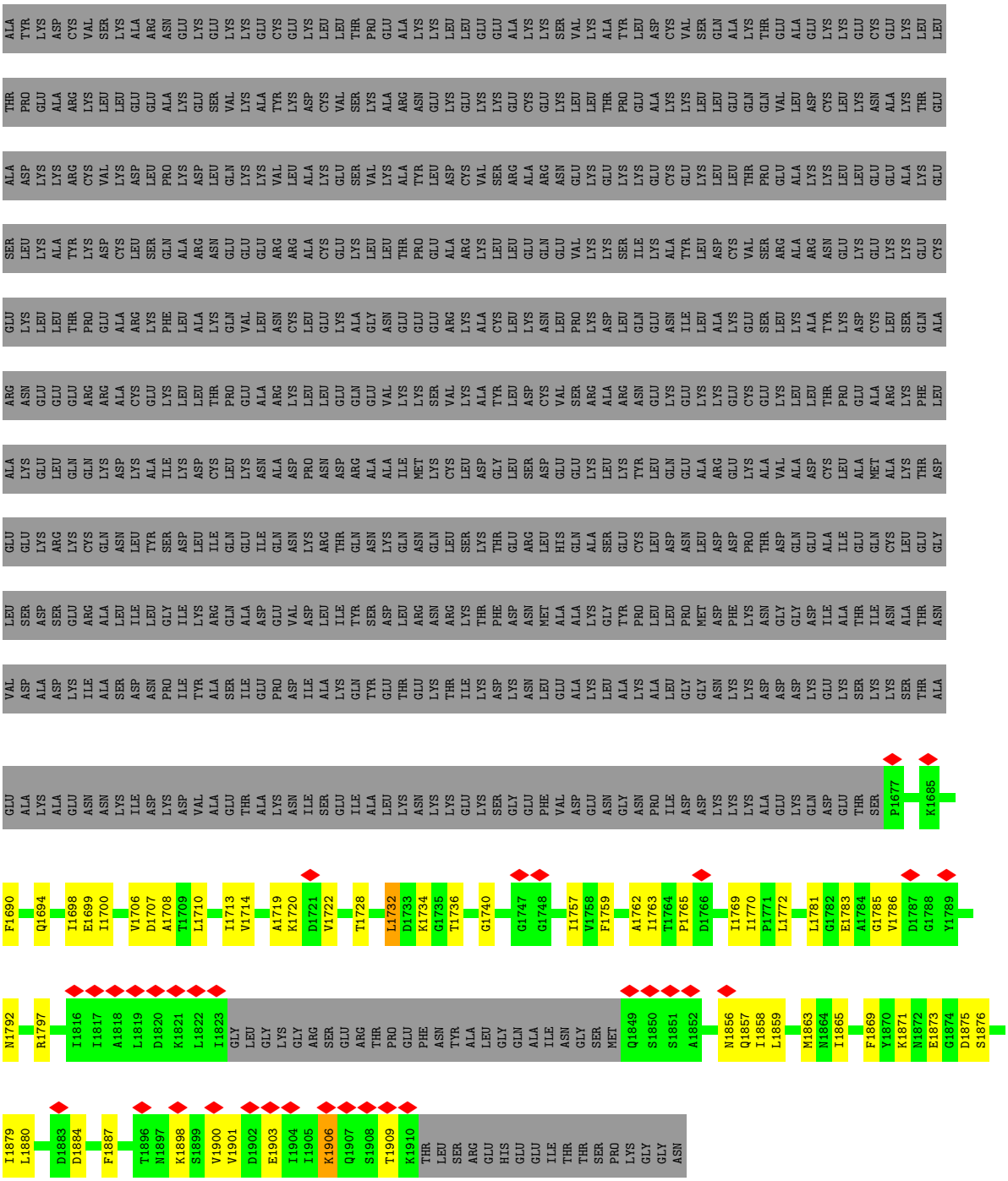


• Molecule 3: Cag pathogenicity island protein (Cag7)

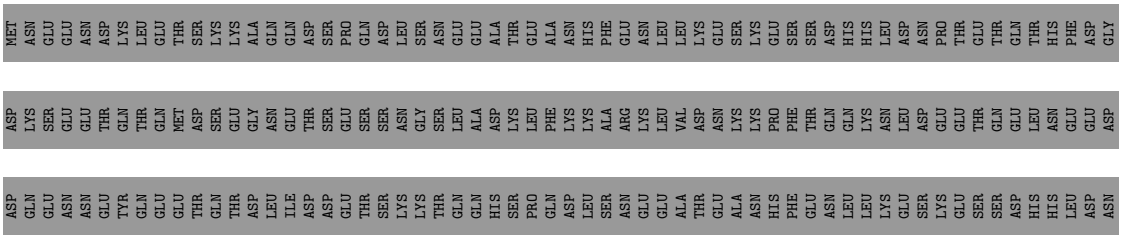
Chain DY: 8% 89%





• Molecule 3: Cag pathogenicity island protein (Cag7)

Chain EY:



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

GLU GLU THR GLN GLU LEU ASN GLU GLU ASP ASP GLN GLU ASN ASN GLU TYR GLN GLU GLU THR GLN THR ASP LEU ILE ASP ASP GLU THR SER LYS LYS THR GLN GLN HIS SER PRO GLN ASP LEU SER ASN GLU GLU ALA THR GLU ALA ASN HIS PHE GLU ASN LEU LEU LYS GLU SER

LYS GLU SER SER ASP HIS HIS LEU ASP ASN PRO THR GLU THR GLN THR ASN PHE ASP GLY ASP LYS SER GLU GLU ILE THR ASP ASP SER ASN ASP GLN GLU ILE ILE LYS GLY SER LYS LYS LYS TYR ILE ILE GLY GLY ILE VAL VAL ALA VAL LEU ILE VAL ILE ILE LEU PHE SER

ARG SER ILE PHE HIS TYR PHE MET PRO LEU GLU ASP LYS SER SER ARG PHE SER LYS ASP ARG ASN LEU TYR VAL ASN ASP GLU ILE GLN ILE ARG GLN GLU TYR ASN ARG LEU LEU LYS GLU ARG ASN GLU LYS GLY ASN MET ILE ASP LYS ASN LEU PHE PHE ASN ASP ASP PRO ASN

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

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

PHE ARG LYS GLU LEU GLU LEU GLN LYS GLU LEU GLN GLU TYR LYS ASP CYS ILE LYS ASN ALA LYS THR GLU ALA GLU LYS ASN LYS CYS LEU LYS GLY LEU SER LYS GLU ALA ILE GLU ARG LEU LYS GLN GLN ALA LEU ASP CYS LEU LYS ASN ALA LYS THR ASP GLU GLU ARG ASN

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

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

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

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

CYS VAL SER GLN ALA LYS THR GLU ALA GLU LYS LYS GLU CYS GLU LYS LEU LEU THR PRO GLU ALA ARG LYS LEU LEU GLU GLU ALA LYS LYS SER VAL LYS ALA TYR LEU ASP CYS VAL SER GLN ALA LYS THR GLU ALA GLU LYS LYS GLU CYS GLU LYS LEU LEU THR PRO GLU ALA

ARG LYS LEU LEU GLU GLU UNK ALA LYS GLU SER VAL LYS ALA TYR LEU ASP CYS VAL SER GLN ALA LYS ASN GLU ALA GLU LYS LYS GLU CYS GLU LYS LEU LEU THR LEU GLU SER LYS LYS LYS LEU GLU GLU ALA LYS LYS SER VAL LYS ALA TYR LEU ASP CYS VAL SER GLN ALA

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

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

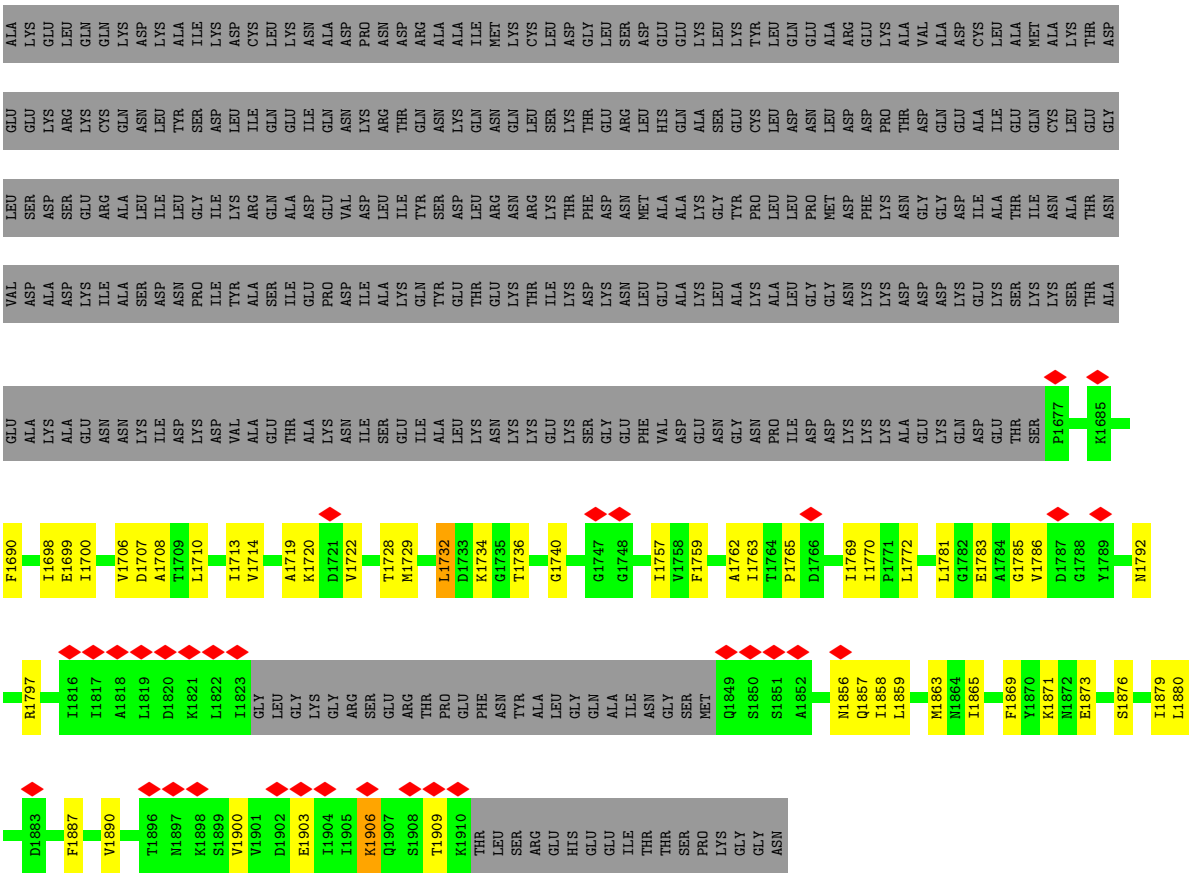
THR PRO GLU ALA ARG LYS LEU LEU GLU GLU ALA LYS GLU SER VAL LYS ALA TYR LYS ASP CYS VAL SER LYS ALA ARG ASN GLU LYS GLU LYS LYS GLU CYS GLU LYS LEU LEU THR PRO GLU ALA LYS LYS LEU LEU GLU GLN GLN VAL LEU ASP CYS LEU LYS ASN ALA LYS THR GLU

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

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

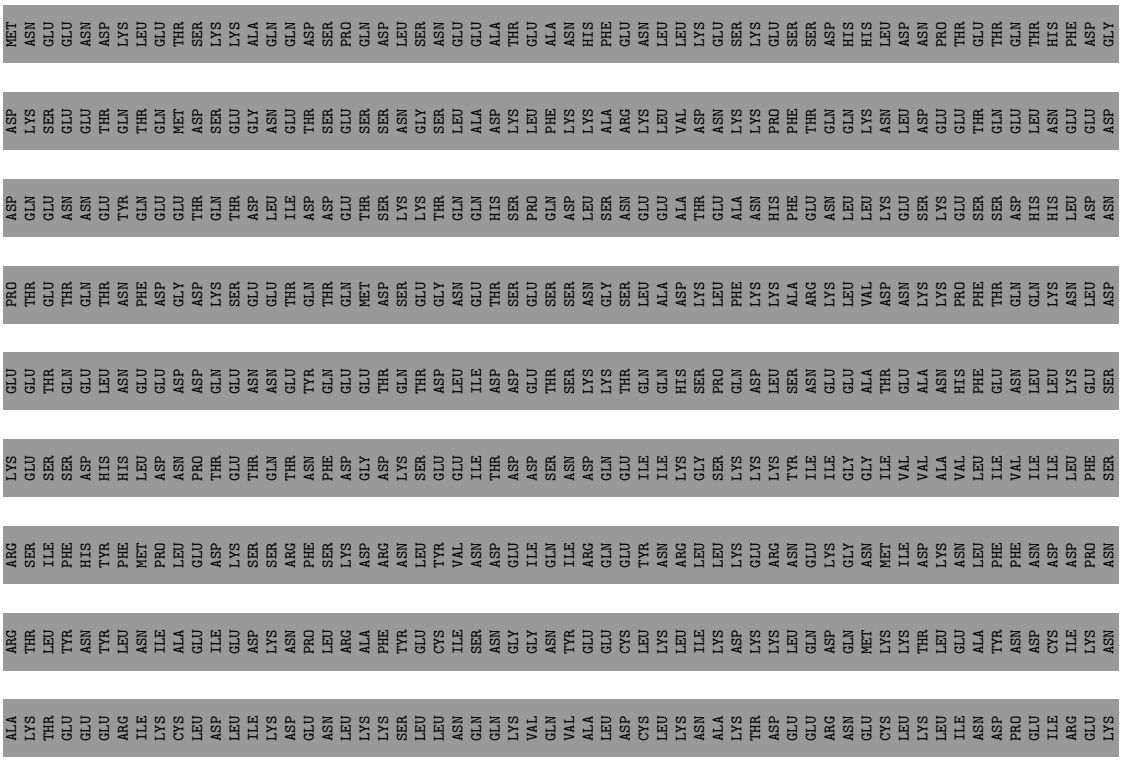
GLU LYS LEU LEU THR PRO GLU ALA ARG LYS PHE LEU ALA LYS GLN VAL LEU ASN CYS LEU GLU LYS ALA GLY ASN GLU GLU GLU ARG LYS ALA CYS LEU LYS ASN LEU PRO LYS ASP LEU GLN GLU ASN ILE LEU ALA LYS GLU SER LEU LYS ALA TYR LYS ASP CYS LEU SER GLN ALA

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



• Molecule 3: Cag pathogenicity island protein (Cag7)

Chain FY: 8% 89%



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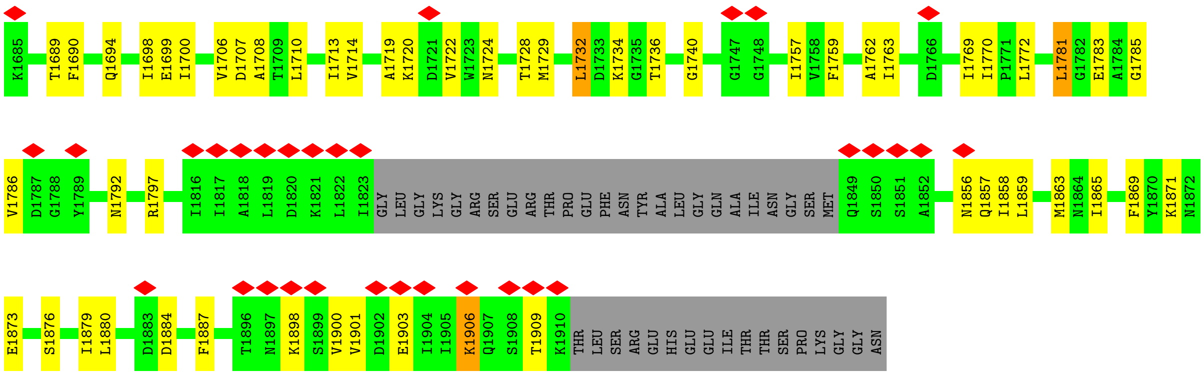
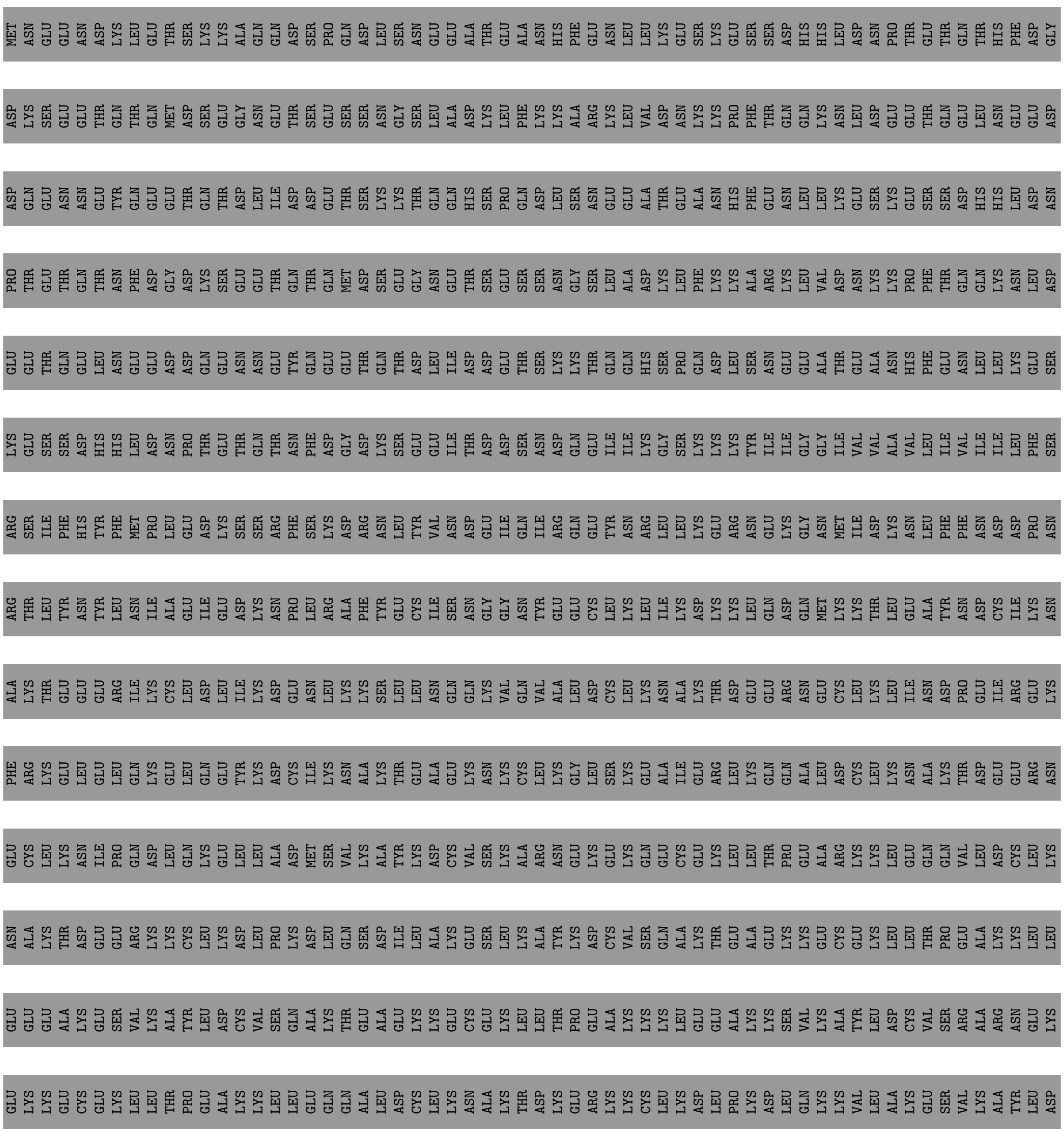
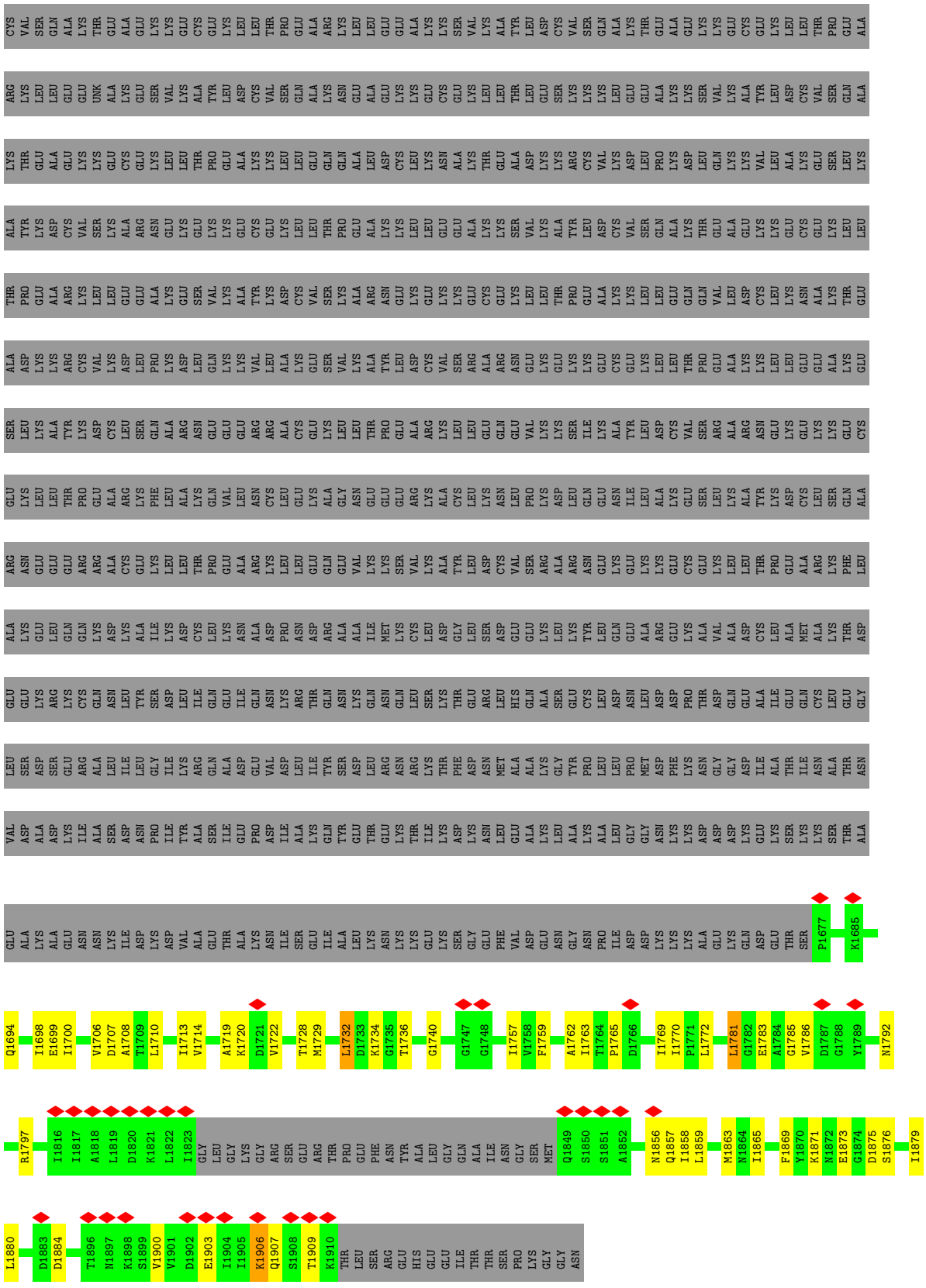


- Molecule 3: Cag pathogenicity island protein (Cag7)

Chain GY:





- Molecule 3: Cag pathogenicity island protein (Cag7)

Chain HY:

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

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

ASP GLN GLU ASN ASN GLU TYR GLN GLU GLU THR GLN THR ASP LEU ILE ASP ASP GLU THR SER LYS LYS THR GLN GLN HIS SER PRO GLN ASP LEU SER ASN GLU GLU ALA THR GLU ALA ASN HIS PHE GLU ASN LEU LEU LYS GLU SER LYS GLU SER SER ASP HIS HIS LEU ASP ASN

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

GLU GLU THR GLN GLU LEU ASN GLU GLU ASP ASP GLN GLU ASN ASN GLU TYR GLN GLU GLU THR GLN THR ASP LEU ILE ASP ASP GLU THR SER LYS LYS THR GLN GLN HIS SER PRO GLN ASP LEU SER ASN GLU GLU ALA THR GLU ALA ASN HIS PHE GLU ASN LEU LEU LYS GLU SER

LYS GLU SER SER ASP HIS HIS LEU ASP ASN PRO THR GLU THR GLN THR ASN PHE ASP GLY ASP LYS SER GLU GLU ILE THR ASP ASP SER ASN ASP GLN GLU ILE ILE LYS GLY SER LYS LYS LYS TYR ILE ILE GLY GLY ILE VAL VAL ALA VAL LEU ILE VAL ILE ILE LEU PHE SER

ARG SER ILE PHE HIS TYR PHE MET PRO LEU GLU ASP LYS SER SER ARG PHE SER LYS ASP ARG ASN LEU TYR VAL ASN ASP GLU ILE GLN ILE ARG GLN GLU TYR ASN ARG LEU LEU LYS GLU ARG ASN GLU LYS GLY ASN MET ILE ASP LYS ASN LEU PHE PHE ASN ASP ASP PRO ASN

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

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

PHE ARG LYS GLU LEU GLU LEU GLN LYS GLU LEU GLN GLU TYR LYS ASP CYS ILE LYS ASN ALA LYS THR GLU ALA GLU LYS ASN LYS CYS LEU LYS GLY LEU SER LYS GLU ALA ILE GLU ARG LEU LYS GLN GLN ALA LEU ASP CYS LEU LYS ASN ALA LYS THR ASP GLU GLU ARG ASN

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

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

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

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

CYS VAL SER GLN ALA LYS THR GLU ALA GLU LYS LYS GLU CYS GLU LYS LEU LEU THR PRO GLU ALA ARG LYS LEU LEU GLU GLU ALA LYS LYS SER VAL LYS ALA TYR LEU ASP CYS VAL SER GLN ALA LYS THR GLU ALA GLU LYS LYS GLU CYS GLU LYS LEU LEU THR PRO GLU ALA

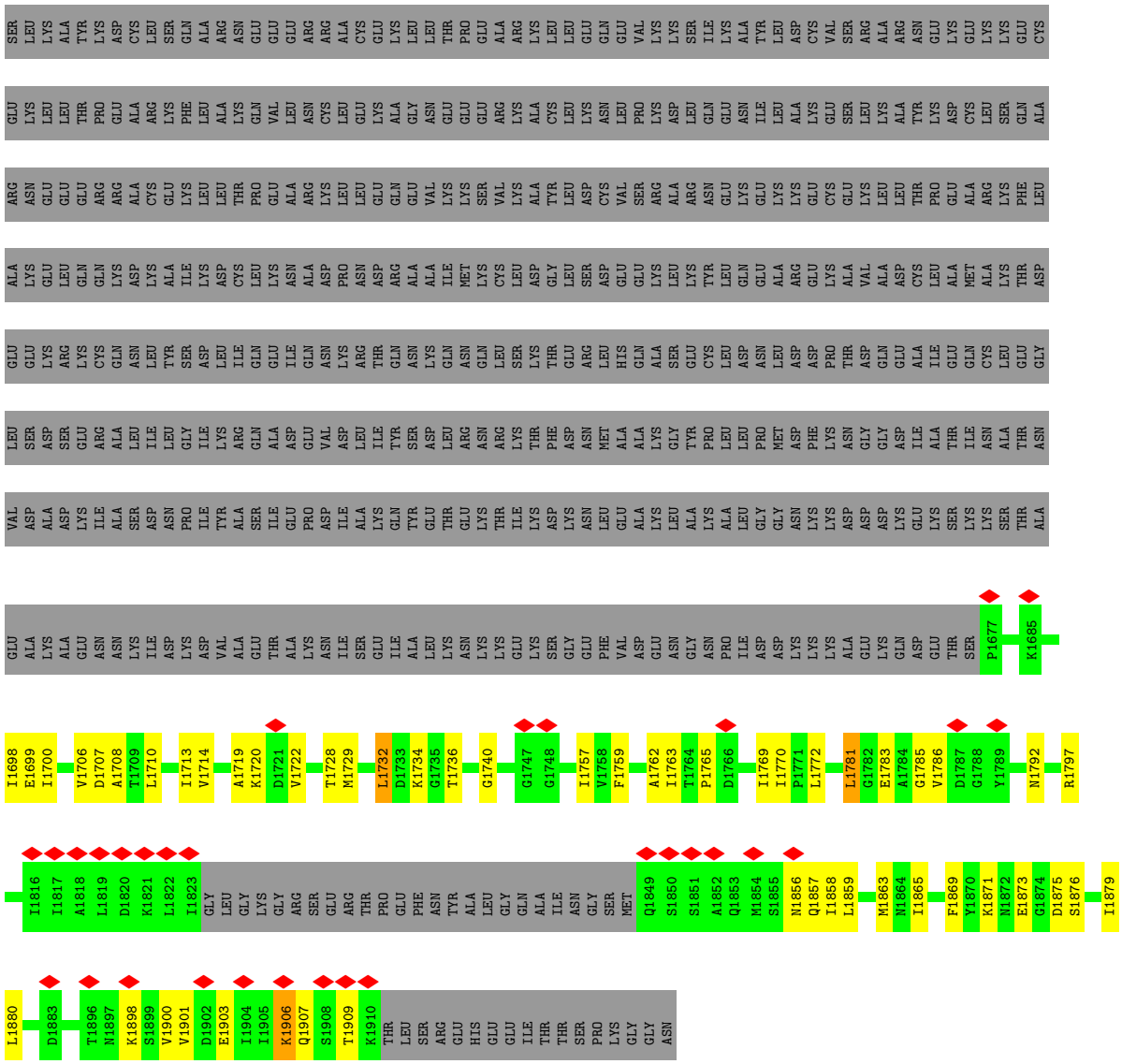
ARG LYS LEU LEU GLU GLU UNK ALA LYS GLU SER VAL LYS ALA TYR LEU ASP CYS VAL SER GLN ALA LYS ASN GLU ALA GLU LYS LYS GLU CYS GLU LYS LEU LEU THR LEU GLU SER LYS LYS LEU GLU GLU ALA LYS LYS SER VAL LYS ALA TYR LEU ASP CYS VAL SER GLN ALA

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

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

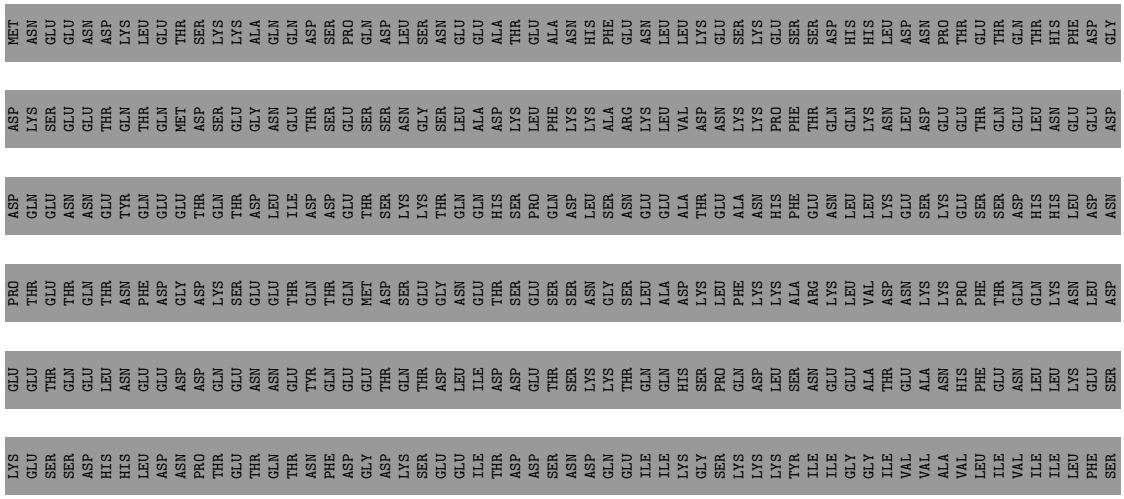
THR PRO GLU ALA ARG LYS LEU LEU GLU GLU ALA LYS GLU SER VAL LYS ALA TYR LYS ASP CYS VAL SER LYS ALA ARG ASN GLU LYS GLU LYS GLU CYS GLU LYS LEU LEU THR PRO GLU ALA LYS LYS LEU LEU GLU GLN GLN VAL LEU ASP CYS LEU LYS ASN ALA LYS THR GLU

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



• Molecule 3: Cag pathogenicity island protein (Cag7)

Chain IY: 8% 89%



ARG SER ILE PHE HIS TYR PHE MET PRO LEU GLU ASP LYS SER SER ARG PHE SER LYS ASP ARG ASN LEU TYR VAL ASN ASP GLU ILE GLN ILE ARG GLN GLU TYR ASN ARG LEU LEU LYS GLU ARG ASN GLU LYS GLY ASN MET ILE ASP LYS ASN LEU PHE PHE ASN ASP ASP PRO ASN

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

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

PHE ARG LYS GLU LEU GLU LEU GLN LYS GLU LEU GLN GLU TYR LYS ASP CYS ILE LYS ASN ALA LYS THR GLU ALA GLU LYS ASN LYS CYS LEU LYS GLY LEU SER LYS GLU ALA ILE GLU ARG LEU LYS GLN GLN ALA LEU ASP CYS LEU LYS ASN ALA LYS THR ASP GLU GLU ARG ASN

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

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

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

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

CYS VAL SER GLN ALA LYS THR GLU ALA GLU LYS LYS GLU CYS GLU LYS LEU LEU THR PRO GLU ALA ARG LYS LEU LEU GLU GLU ALA LYS LYS SER VAL LYS ALA TYR LEU ASP CYS VAL SER GLN ALA LYS THR GLU ALA GLU LYS LYS GLU CYS GLU LYS LEU LEU THR PRO GLU ALA

ARG LYS LEU LEU GLU GLU UNK ALA LYS GLU SER VAL LYS ALA TYR LEU ASP CYS VAL SER GLN ALA LYS ASN GLU ALA GLU LYS LYS GLU CYS GLU LYS LEU LEU THR LEU GLU SER LYS LYS LYS LEU GLU GLU ALA LYS LYS SER VAL LYS ALA TYR LEU ASP CYS VAL SER GLN ALA

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

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

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

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

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

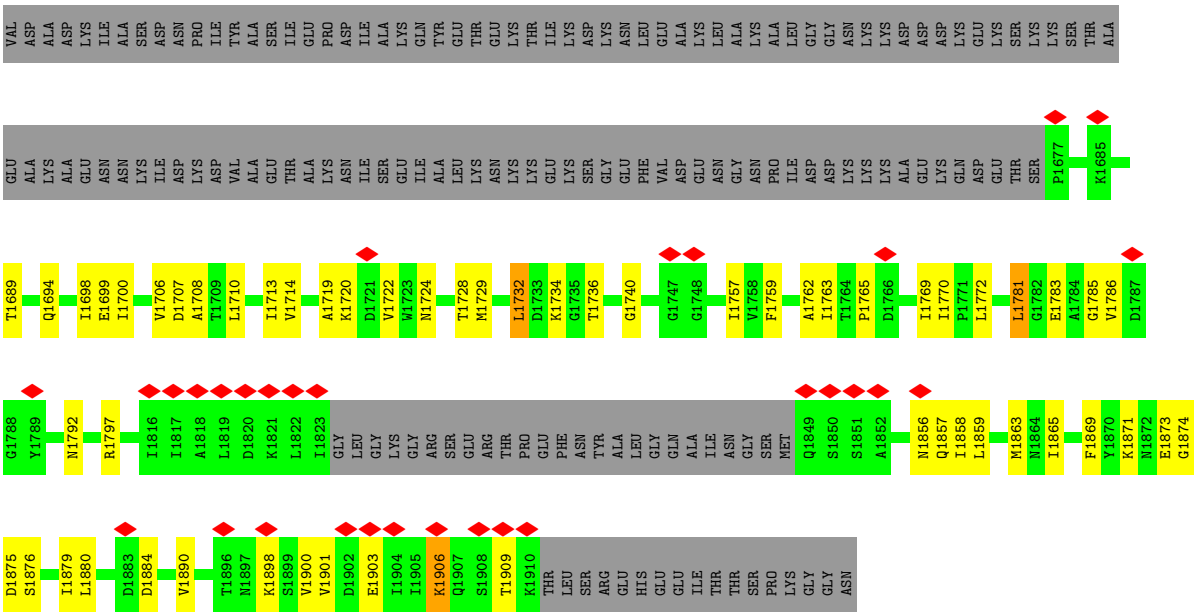
GLU LYS LEU LEU THR PRO GLU ALA ARG LYS PHE LEU ALA LYS GLN VAL LEU ASN CYS LEU GLU LYS ALA GLY ASN GLU GLU GLU ARG LYS ALA CYS LEU LYS ASN LEU PRO LYS ASP LEU GLN GLU ASN ILE LEU ALA LYS GLU SER LEU LYS ALA TYR LYS ASP CYS LEU SER GLN ALA

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

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

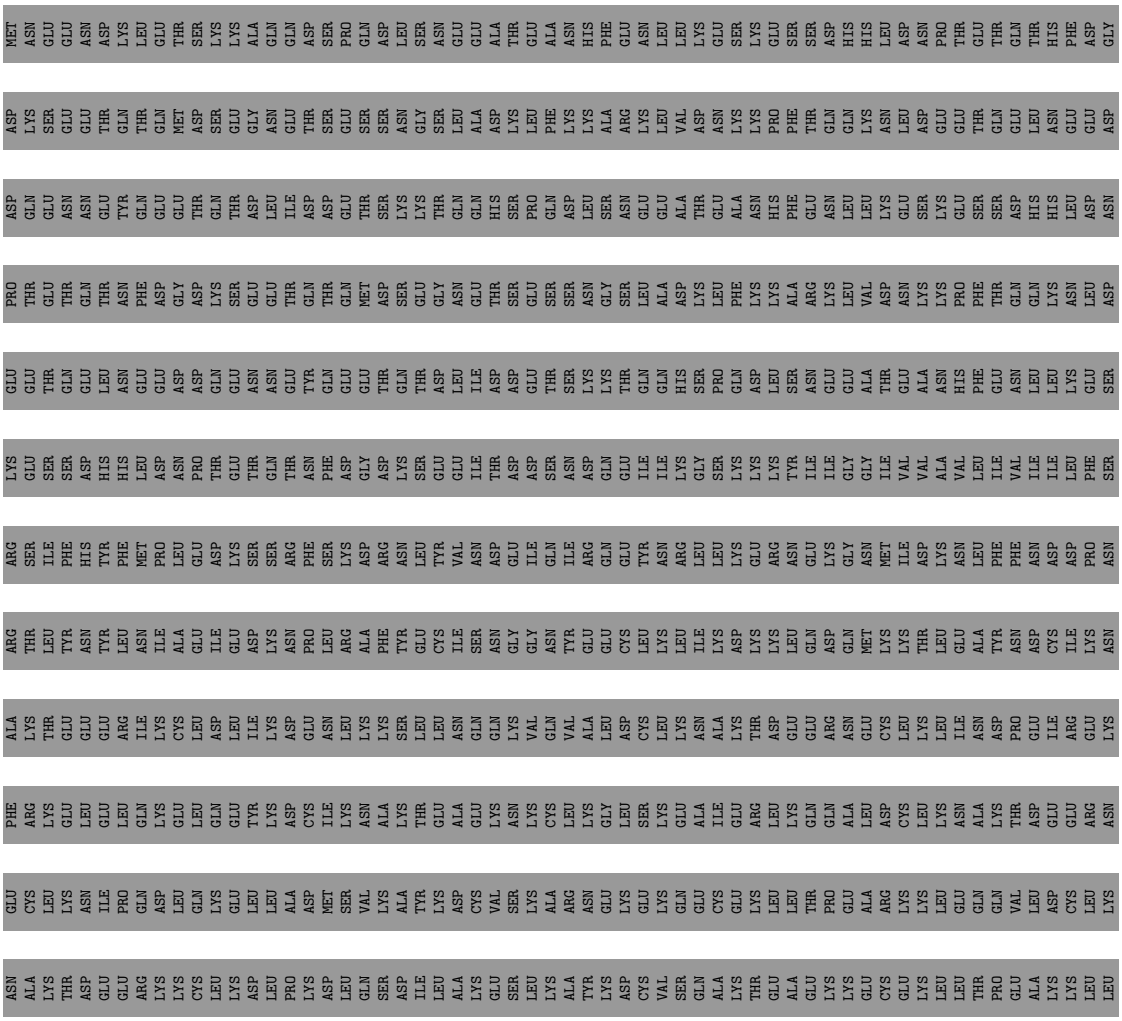
GLU GLU LYS ARG LYS CYS GLN ASN LEU TYR SER ASP LEU ILE GLN GLU ILE GLN ASN LYS ARG THR GLN ASN LYS GLN ASN GLN LEU SER LYS THR GLU ARG LEU HIS GLN ALA SER GLU CYS LEU ASP ASN LEU ASP ASP PRO THR ASP GLN GLU ALA ILE GLU GLN CYS LEU GLU GLY

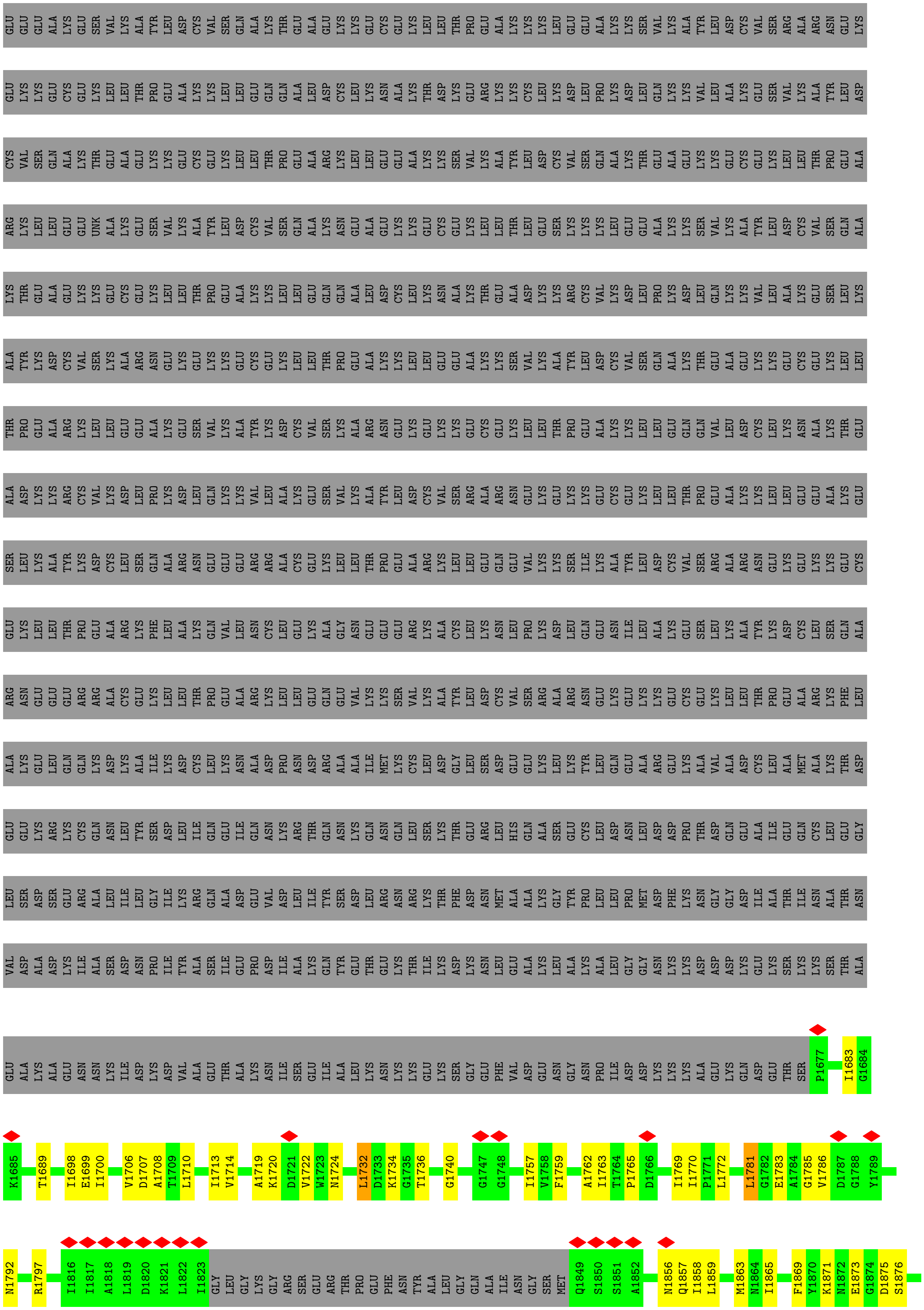
LEU SER ASP SER GLU ARG ALA LEU ILE LEU GLY ILE LYS ARG GLN ALA ASP GLU VAL ASP LEU ILE TYR SER ASP LEU ARG ASN ARG LYS THR PHE ASP ASN MET ALA ALA LYS GLY TYR PRO LEU LEU PRO MET ASP PHE LYS ASN GLY GLY ASP ILE ALA THR ILE ASN ALA THR ASN



- Molecule 3: Cag pathogenicity island protein (Cag7)

Chain JY:

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I1763
T1764
P1765
D1766
I1769
I1770
P1771
L1772
L1781
G1782
E1783
A1784
G1785
V1786
D1787
G1788
Y1789
N1792
R1797
I1816
I1817
A1818
L1819
D1820
K1821
L1822
I1823
Q1849
S1850
S1851
A1852
N1856
Q1857
I1858
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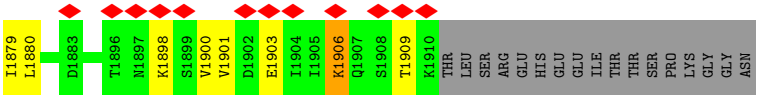
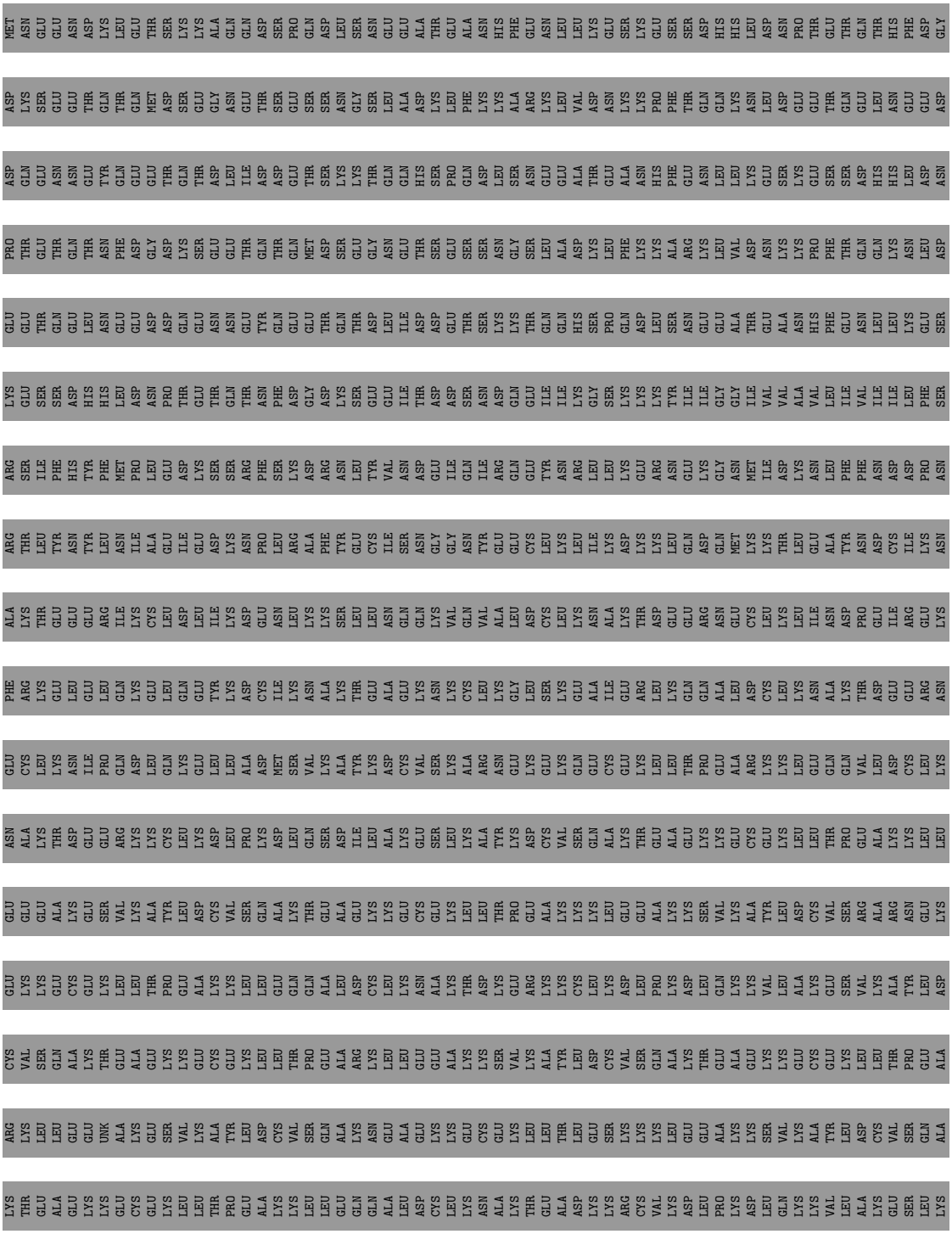
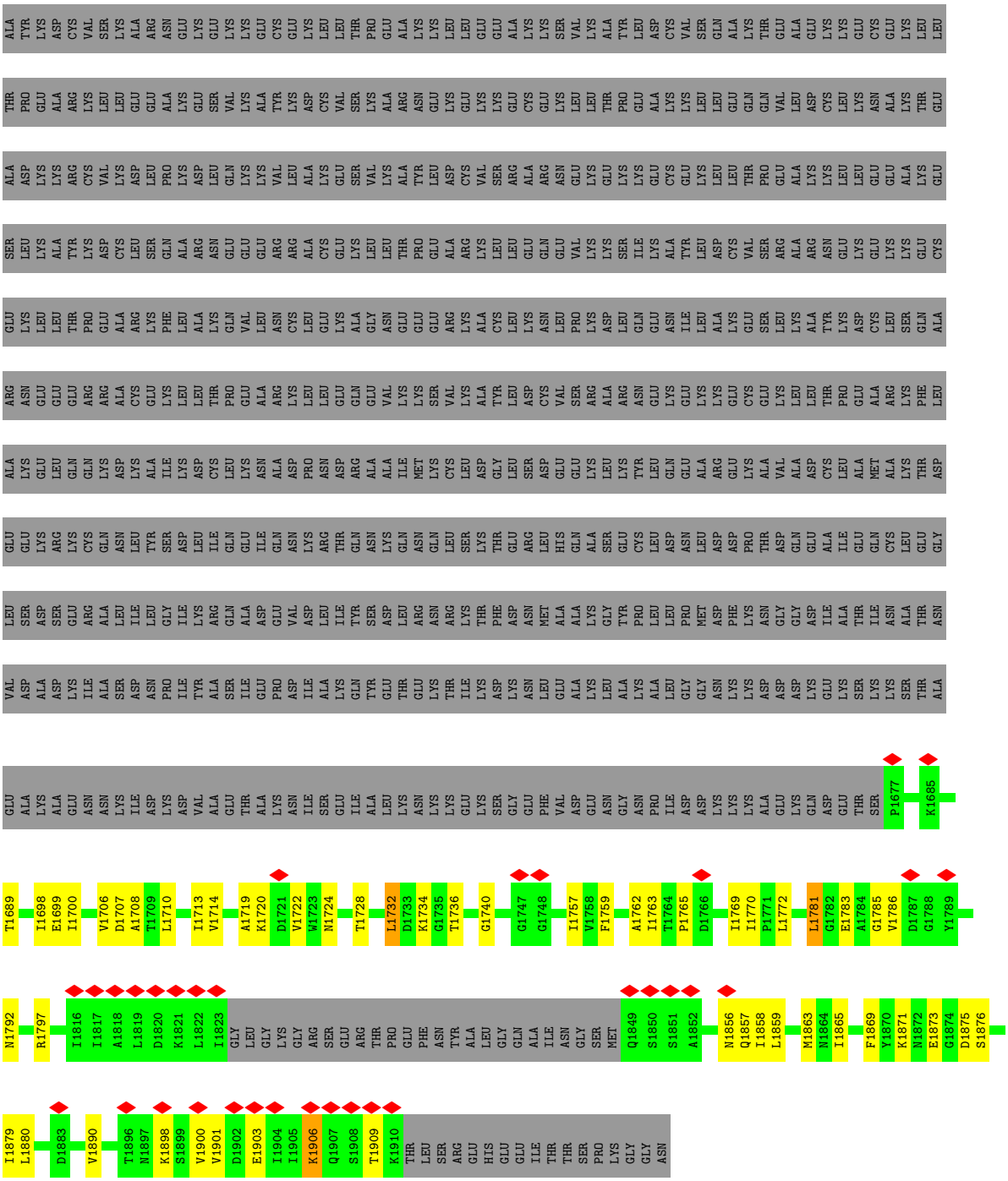
G1874
D1875
S1876



- Molecule 3: Cag pathogenicity island protein (Cag7)

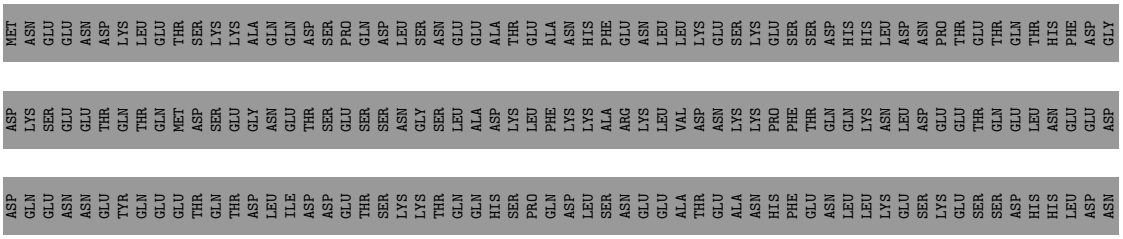
Chain KY:

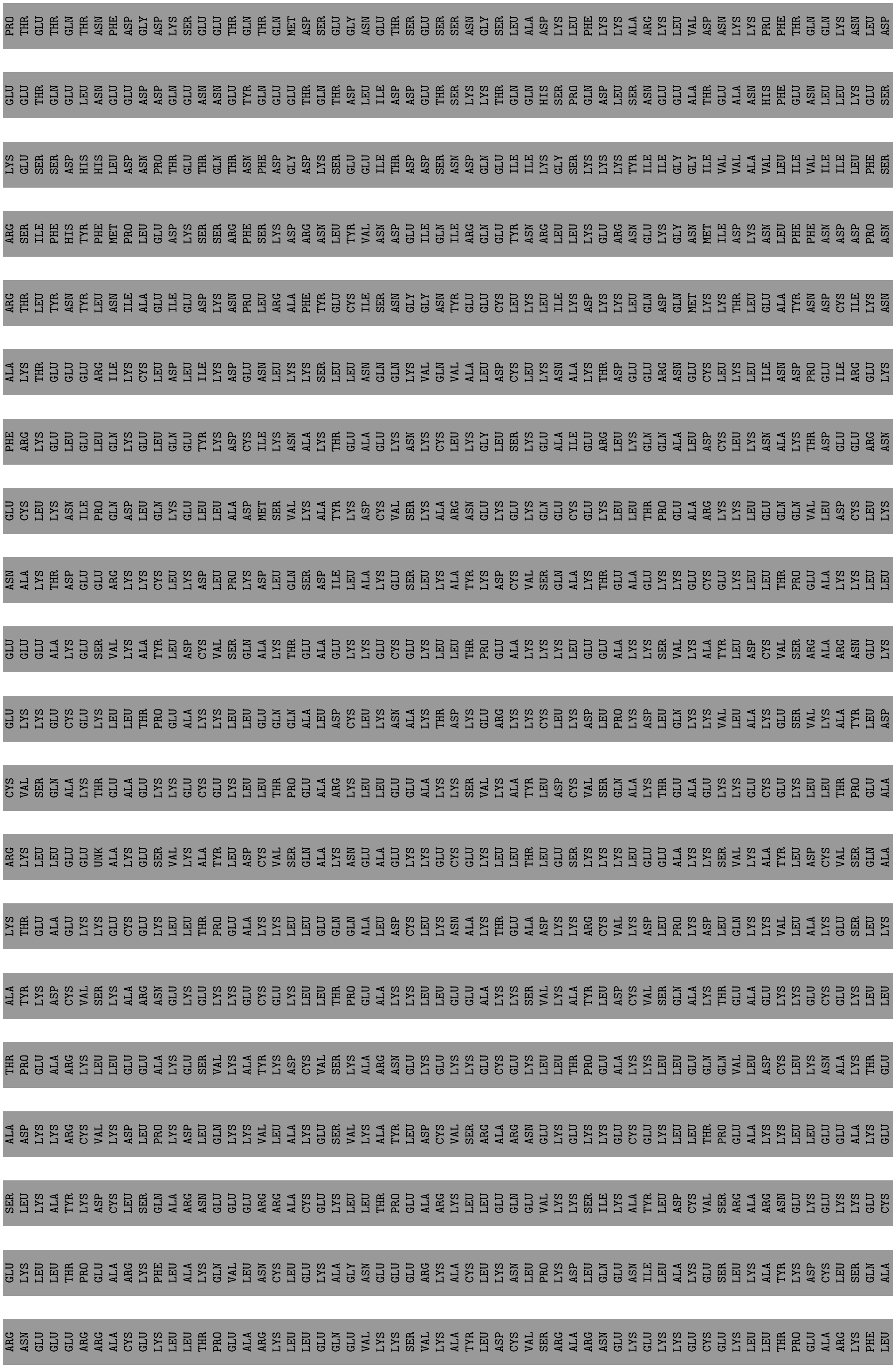


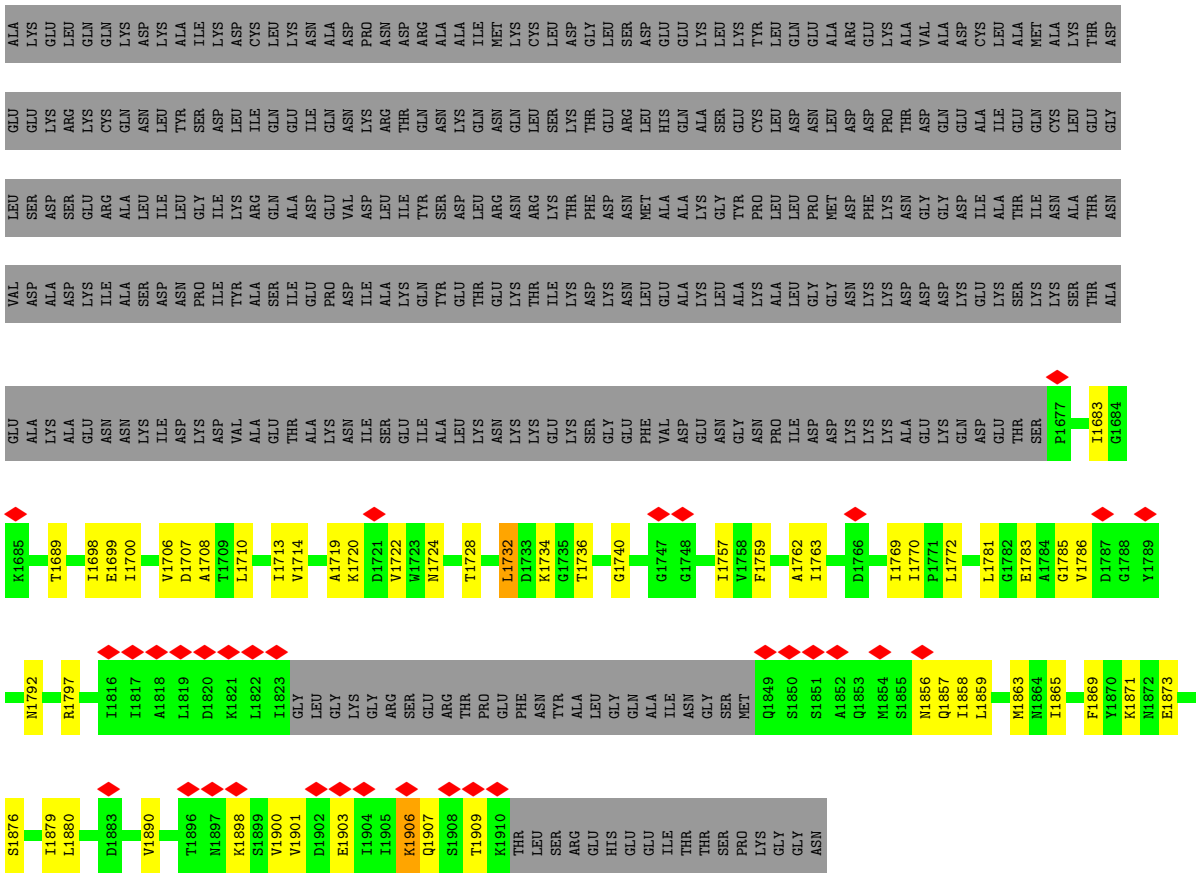


• Molecule 3: Cag pathogenicity island protein (Cag7)

Chain LY:

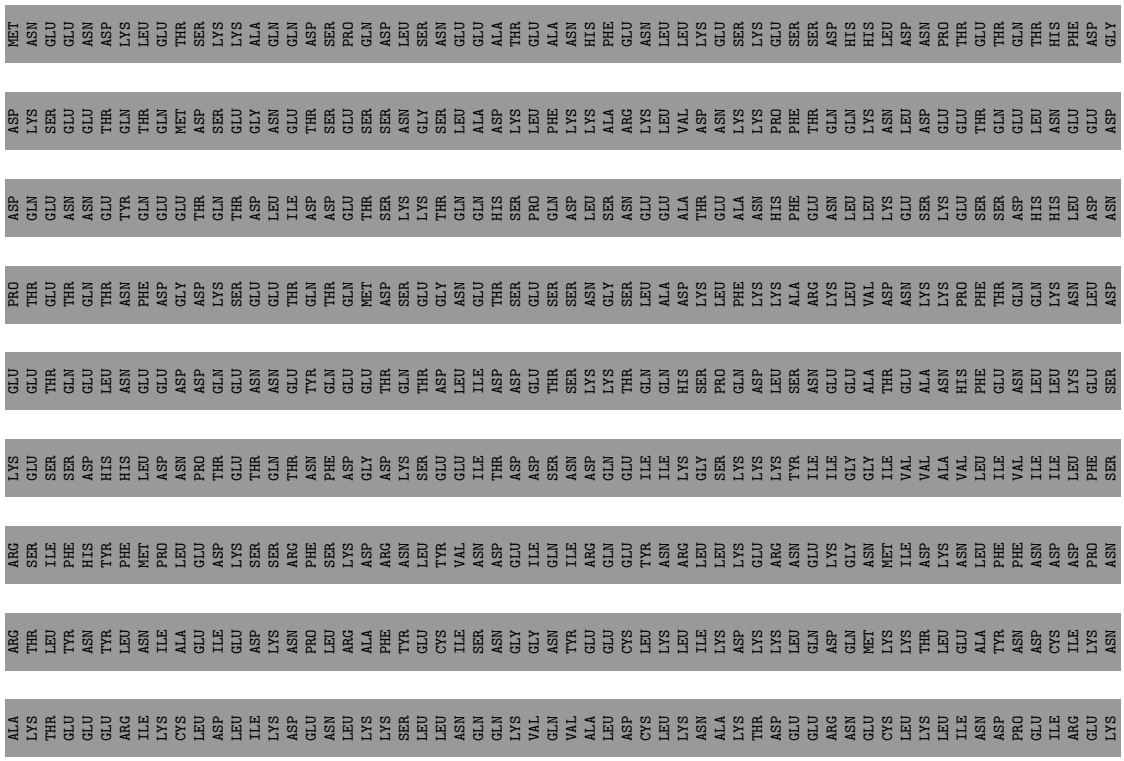



PRO THR GLU THR GLN THR ASN PHE ASP GLY ASP LYS SER GLU GLU THR GLN THR GLN MET ASP SER GLU GLY ASN GLU THR SER GLU SER SER ASN GLY SER LEU ALA ASP LYS LEU PHE LYS LYS ALA ARG LYS LEU VAL ASP ASN LYS LYS PRO PHE THR GLN GLN LYS ASN LEU ASP
GLU GLU THR GLN GLU LEU ASN GLU GLU ASP ASP GLN GLU ASN ASN GLU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

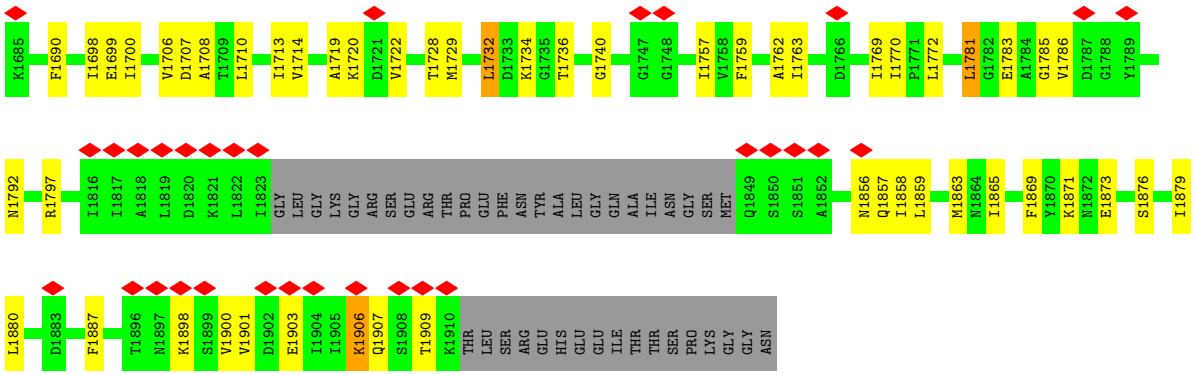


• Molecule 3: Cag pathogenicity island protein (Cag7)

Chain MY: 8% 89%

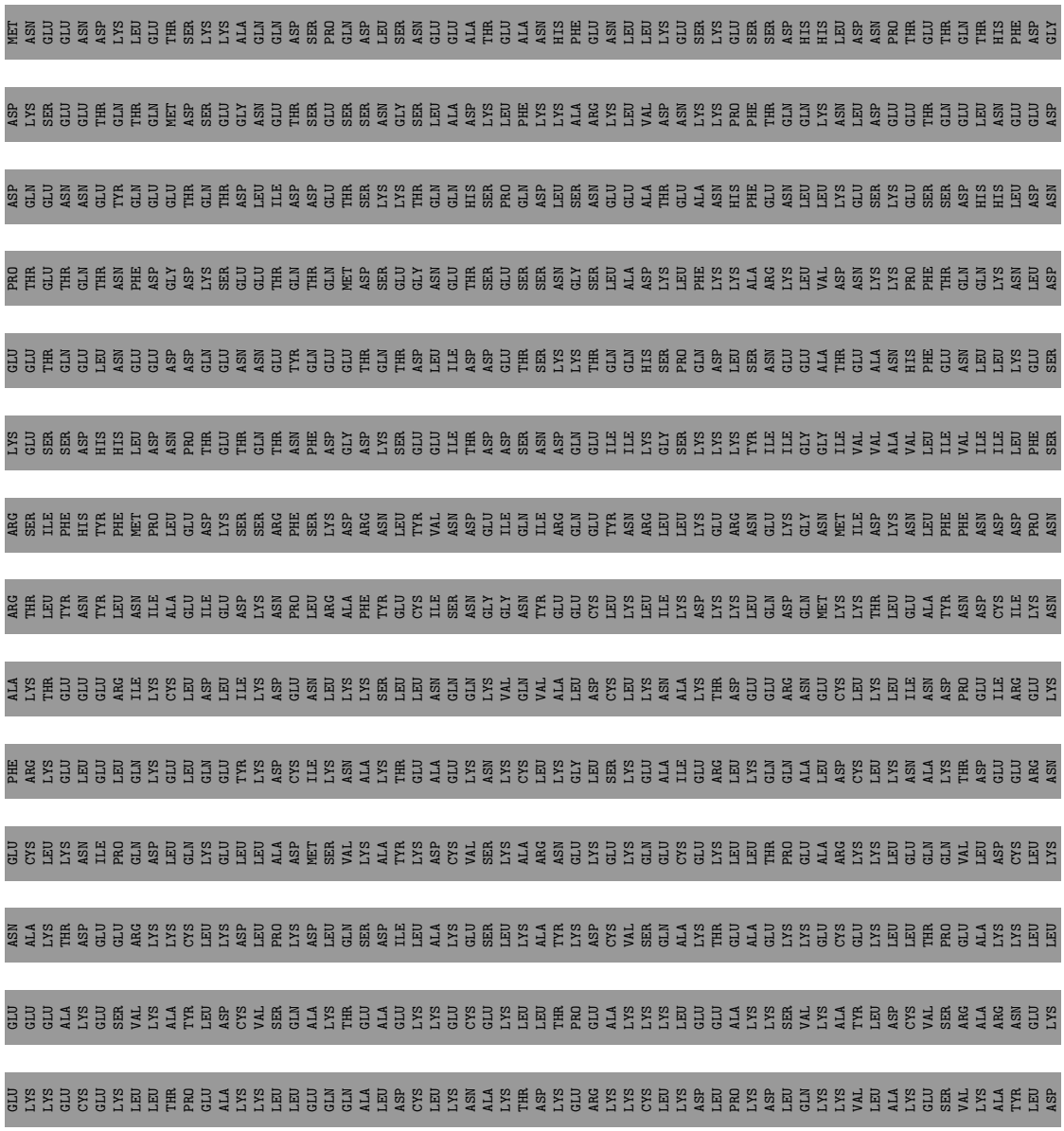


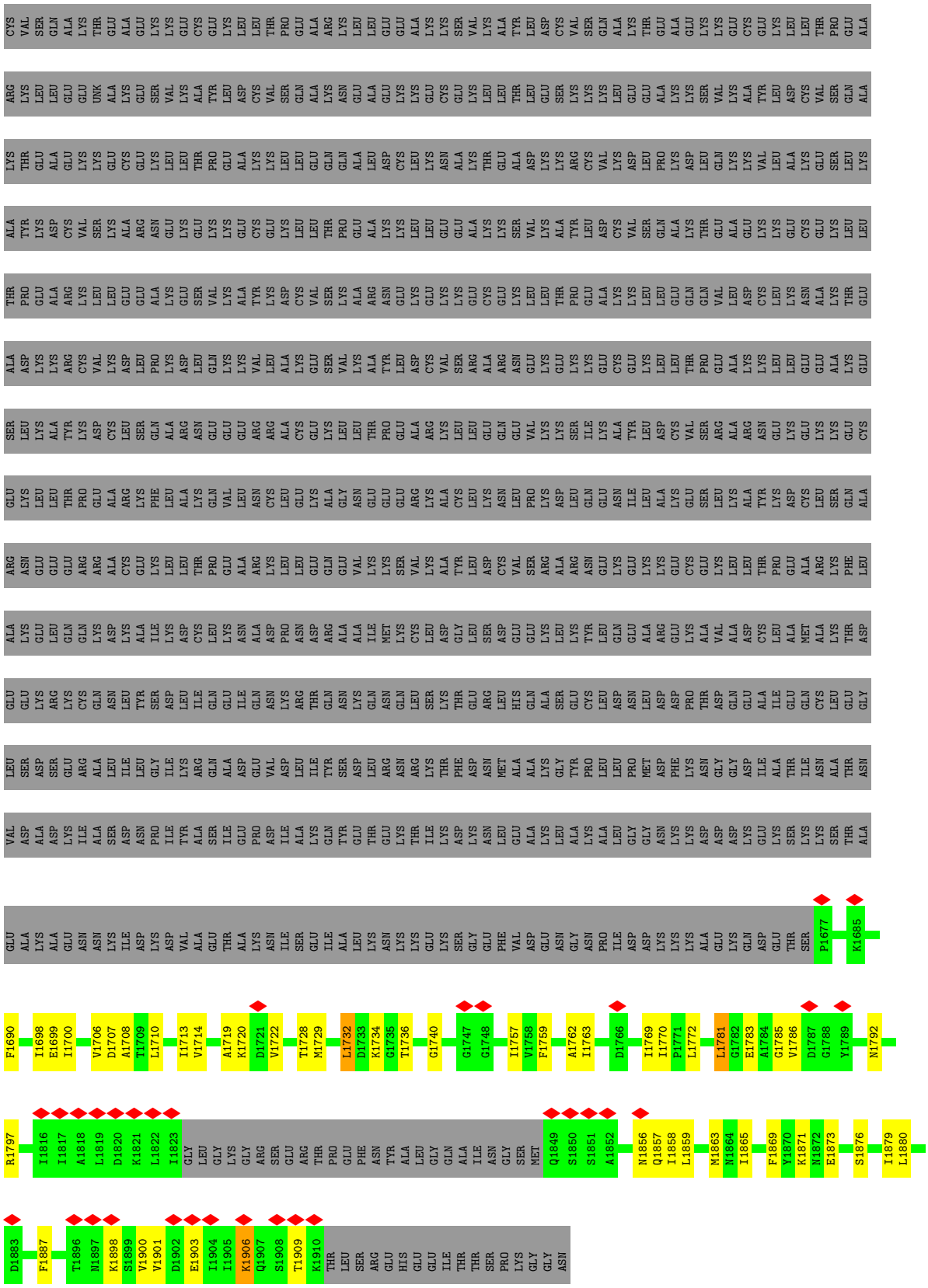
P1677
I1683
G1684



• Molecule 3: Cag pathogenicity island protein (Cag7)

Chain NY:




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

4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C14	Depositor
Number of particles used	7337	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 ($e^-/Å^2$)	59.7	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.049	Depositor
Minimum map value	-0.036	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.002	Depositor
Recommended contour level	0.015	Depositor
Map size (Å)	510.0, 510.0, 510.0	wwPDB
Map dimensions	510, 510, 510	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.0, 1.0, 1.0	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	AT	0.41	0/1525	0.71	4/2049 (0.2%)
1	BT	0.41	0/1525	0.71	4/2049 (0.2%)
1	CT	0.41	0/1525	0.71	4/2049 (0.2%)
1	DT	0.41	0/1525	0.71	4/2049 (0.2%)
1	ET	0.41	0/1525	0.71	4/2049 (0.2%)
1	FT	0.41	0/1525	0.71	4/2049 (0.2%)
1	GT	0.41	0/1525	0.71	4/2049 (0.2%)
1	HT	0.41	0/1525	0.71	4/2049 (0.2%)
1	IT	0.41	0/1525	0.71	4/2049 (0.2%)
1	JT	0.41	0/1525	0.71	4/2049 (0.2%)
1	KT	0.41	0/1525	0.71	4/2049 (0.2%)
1	LT	0.41	0/1525	0.71	4/2049 (0.2%)
1	MT	0.41	0/1525	0.71	4/2049 (0.2%)
1	NT	0.41	0/1525	0.71	4/2049 (0.2%)
2	AX	0.54	0/1301	0.69	0/1752
2	BX	0.54	0/1301	0.69	0/1752
2	CX	0.54	0/1301	0.69	0/1752
2	DX	0.54	0/1301	0.69	0/1752
2	EX	0.54	0/1301	0.69	0/1752
2	FX	0.54	0/1301	0.69	1/1752 (0.1%)
2	GX	0.54	0/1301	0.69	0/1752
2	HX	0.54	0/1301	0.69	0/1752
2	IX	0.54	0/1301	0.69	0/1752
2	JX	0.54	0/1301	0.69	0/1752
2	KX	0.54	0/1301	0.69	1/1752 (0.1%)
2	LX	0.54	0/1301	0.69	1/1752 (0.1%)
2	MX	0.54	0/1301	0.69	1/1752 (0.1%)
2	NX	0.54	0/1301	0.69	1/1752 (0.1%)
3	AY	0.55	0/1602	0.72	3/2170 (0.1%)
3	BY	0.55	0/1602	0.72	4/2170 (0.2%)
3	CY	0.55	0/1602	0.72	4/2170 (0.2%)
3	DY	0.55	0/1602	0.72	3/2170 (0.1%)
3	EY	0.55	0/1602	0.72	3/2170 (0.1%)
3	FY	0.55	0/1602	0.72	4/2170 (0.2%)

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	#|Z| >5	RMSZ	#|Z| >5
3	GY	0.55	0/1602	0.72	4/2170 (0.2%)
3	HY	0.55	0/1602	0.72	4/2170 (0.2%)
3	IY	0.55	0/1602	0.72	4/2170 (0.2%)
3	JY	0.55	0/1602	0.72	4/2170 (0.2%)
3	KY	0.55	0/1602	0.72	4/2170 (0.2%)
3	LY	0.55	0/1602	0.72	3/2170 (0.1%)
3	MY	0.55	0/1602	0.72	4/2170 (0.2%)
3	NY	0.55	0/1602	0.72	4/2170 (0.2%)
All	All	0.50	0/61992	0.71	113/83594 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
2	AX	0	1
2	BX	0	1
2	CX	0	1
2	DX	0	1
2	EX	0	1
2	FX	0	1
2	GX	0	1
2	HX	0	1
2	IX	0	1
2	JX	0	1
2	KX	0	1
2	LX	0	1
2	MX	0	1
2	NX	0	1
All	All	0	14

There are no bond length outliers.

All (113) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	MY	1772	LEU	CA-CB-CG	7.79	133.22	115.30
3	CY	1772	LEU	CA-CB-CG	7.79	133.21	115.30
3	KY	1772	LEU	CA-CB-CG	7.78	133.20	115.30
3	AY	1772	LEU	CA-CB-CG	7.78	133.19	115.30
3	HY	1772	LEU	CA-CB-CG	7.78	133.20	115.30
3	FY	1772	LEU	CA-CB-CG	7.78	133.19	115.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	LY	1772	LEU	CA-CB-CG	7.77	133.18	115.30
3	EY	1772	LEU	CA-CB-CG	7.77	133.18	115.30
3	JY	1772	LEU	CA-CB-CG	7.77	133.17	115.30
3	NY	1772	LEU	CA-CB-CG	7.77	133.17	115.30
3	DY	1772	LEU	CA-CB-CG	7.77	133.17	115.30
3	GY	1772	LEU	CA-CB-CG	7.76	133.16	115.30
3	BY	1772	LEU	CA-CB-CG	7.76	133.15	115.30
3	IY	1772	LEU	CA-CB-CG	7.76	133.15	115.30
1	BT	168	LEU	CA-CB-CG	7.12	131.68	115.30
1	LT	168	LEU	CA-CB-CG	7.12	131.68	115.30
1	GT	168	LEU	CA-CB-CG	7.12	131.67	115.30
1	IT	168	LEU	CA-CB-CG	7.12	131.67	115.30
1	DT	168	LEU	CA-CB-CG	7.11	131.66	115.30
1	NT	168	LEU	CA-CB-CG	7.11	131.66	115.30
1	AT	168	LEU	CA-CB-CG	7.11	131.65	115.30
1	FT	168	LEU	CA-CB-CG	7.11	131.65	115.30
1	HT	168	LEU	CA-CB-CG	7.11	131.65	115.30
1	JT	168	LEU	CA-CB-CG	7.11	131.64	115.30
1	KT	168	LEU	CA-CB-CG	7.10	131.64	115.30
1	ET	168	LEU	CA-CB-CG	7.10	131.63	115.30
1	MT	168	LEU	CA-CB-CG	7.10	131.63	115.30
1	CT	168	LEU	CA-CB-CG	7.09	131.61	115.30
1	KT	72	ASP	CB-CG-OD2	6.55	124.20	118.30
1	DT	72	ASP	CB-CG-OD2	6.50	124.15	118.30
1	ET	72	ASP	CB-CG-OD2	6.50	124.15	118.30
1	AT	72	ASP	CB-CG-OD2	6.50	124.15	118.30
1	CT	72	ASP	CB-CG-OD2	6.49	124.14	118.30
1	LT	72	ASP	CB-CG-OD2	6.49	124.14	118.30
1	FT	72	ASP	CB-CG-OD2	6.48	124.13	118.30
1	BT	72	ASP	CB-CG-OD2	6.48	124.13	118.30
1	JT	72	ASP	CB-CG-OD2	6.47	124.12	118.30
1	NT	72	ASP	CB-CG-OD2	6.46	124.11	118.30
1	MT	72	ASP	CB-CG-OD2	6.45	124.10	118.30
1	HT	72	ASP	CB-CG-OD2	6.45	124.10	118.30
1	IT	72	ASP	CB-CG-OD2	6.44	124.09	118.30
1	GT	72	ASP	CB-CG-OD2	6.43	124.08	118.30
3	HY	1732	LEU	CB-CG-CD2	-5.92	100.94	111.00
3	AY	1732	LEU	CB-CG-CD2	-5.91	100.95	111.00
3	JY	1732	LEU	CB-CG-CD2	-5.90	100.96	111.00
3	MY	1732	LEU	CB-CG-CD2	-5.90	100.97	111.00
3	LY	1732	LEU	CB-CG-CD2	-5.90	100.97	111.00
3	GY	1732	LEU	CB-CG-CD2	-5.90	100.97	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	KY	1732	LEU	CB-CG-CD2	-5.89	100.98	111.00
3	CY	1732	LEU	CB-CG-CD2	-5.89	100.98	111.00
3	BY	1732	LEU	CB-CG-CD2	-5.89	100.99	111.00
3	IY	1732	LEU	CB-CG-CD2	-5.89	100.99	111.00
3	NY	1732	LEU	CB-CG-CD2	-5.89	100.99	111.00
3	EY	1732	LEU	CB-CG-CD2	-5.89	100.99	111.00
3	FY	1732	LEU	CB-CG-CD2	-5.88	101.01	111.00
3	DY	1732	LEU	CB-CG-CD2	-5.87	101.02	111.00
3	KY	1772	LEU	CB-CG-CD1	5.49	120.34	111.00
3	GY	1772	LEU	CB-CG-CD1	5.49	120.33	111.00
3	CY	1772	LEU	CB-CG-CD1	5.49	120.33	111.00
3	JY	1772	LEU	CB-CG-CD1	5.49	120.33	111.00
3	EY	1772	LEU	CB-CG-CD1	5.48	120.31	111.00
3	HY	1772	LEU	CB-CG-CD1	5.47	120.30	111.00
3	AY	1772	LEU	CB-CG-CD1	5.46	120.29	111.00
3	NY	1772	LEU	CB-CG-CD1	5.46	120.29	111.00
3	MY	1772	LEU	CB-CG-CD1	5.46	120.29	111.00
3	LY	1772	LEU	CB-CG-CD1	5.46	120.28	111.00
3	BY	1772	LEU	CB-CG-CD1	5.46	120.28	111.00
3	FY	1772	LEU	CB-CG-CD1	5.46	120.28	111.00
3	IY	1772	LEU	CB-CG-CD1	5.44	120.25	111.00
3	DY	1772	LEU	CB-CG-CD1	5.44	120.24	111.00
1	MT	28	VAL	CG1-CB-CG2	-5.27	102.47	110.90
1	DT	28	VAL	CG1-CB-CG2	-5.27	102.47	110.90
1	BT	28	VAL	CG1-CB-CG2	-5.25	102.50	110.90
1	IT	28	VAL	CG1-CB-CG2	-5.25	102.50	110.90
1	AT	28	VAL	CG1-CB-CG2	-5.25	102.50	110.90
1	JT	28	VAL	CG1-CB-CG2	-5.24	102.52	110.90
1	LT	28	VAL	CG1-CB-CG2	-5.24	102.52	110.90
1	CT	28	VAL	CG1-CB-CG2	-5.24	102.52	110.90
1	NT	28	VAL	CG1-CB-CG2	-5.24	102.53	110.90
1	FT	28	VAL	CG1-CB-CG2	-5.23	102.53	110.90
1	ET	28	VAL	CG1-CB-CG2	-5.23	102.53	110.90
1	GT	28	VAL	CG1-CB-CG2	-5.23	102.53	110.90
1	KT	28	VAL	CG1-CB-CG2	-5.22	102.55	110.90
1	HT	28	VAL	CG1-CB-CG2	-5.21	102.56	110.90
1	IT	197	GLN	CA-CB-CG	5.20	124.84	113.40
1	AT	197	GLN	CA-CB-CG	5.19	124.83	113.40
1	DT	197	GLN	CA-CB-CG	5.19	124.81	113.40
1	ET	197	GLN	CA-CB-CG	5.18	124.81	113.40
1	LT	197	GLN	CA-CB-CG	5.18	124.80	113.40
1	FT	197	GLN	CA-CB-CG	5.18	124.80	113.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	JT	197	GLN	CA-CB-CG	5.18	124.79	113.40
1	HT	197	GLN	CA-CB-CG	5.17	124.79	113.40
1	MT	197	GLN	CA-CB-CG	5.17	124.78	113.40
1	BT	197	GLN	CA-CB-CG	5.17	124.78	113.40
1	KT	197	GLN	CA-CB-CG	5.17	124.78	113.40
1	CT	197	GLN	CA-CB-CG	5.17	124.77	113.40
1	NT	197	GLN	CA-CB-CG	5.16	124.76	113.40
1	GT	197	GLN	CA-CB-CG	5.16	124.75	113.40
2	KX	422	ASP	CB-CG-OD2	5.03	122.82	118.30
3	HY	1781	LEU	CA-CB-CG	-5.02	103.75	115.30
2	FX	422	ASP	CB-CG-OD2	5.02	122.82	118.30
3	JY	1781	LEU	CA-CB-CG	-5.02	103.76	115.30
3	NY	1781	LEU	CA-CB-CG	-5.02	103.76	115.30
2	NX	422	ASP	CB-CG-OD2	5.01	122.81	118.30
3	BY	1781	LEU	CA-CB-CG	-5.01	103.79	115.30
2	MX	422	ASP	CB-CG-OD2	5.01	122.81	118.30
3	FY	1781	LEU	CA-CB-CG	-5.00	103.79	115.30
3	MY	1781	LEU	CA-CB-CG	-5.00	103.79	115.30
3	GY	1781	LEU	CA-CB-CG	-5.00	103.79	115.30
3	IY	1781	LEU	CA-CB-CG	-5.00	103.79	115.30
3	KY	1781	LEU	CA-CB-CG	-5.00	103.79	115.30
3	CY	1781	LEU	CA-CB-CG	-5.00	103.80	115.30
2	LX	422	ASP	CB-CG-OD2	5.00	122.80	118.30

There are no chirality outliers.

All (14) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	AX	479	ILE	Peptide
2	BX	479	ILE	Peptide
2	CX	479	ILE	Peptide
2	DX	479	ILE	Peptide
2	EX	479	ILE	Peptide
2	FX	479	ILE	Peptide
2	GX	479	ILE	Peptide
2	HX	479	ILE	Peptide
2	IX	479	ILE	Peptide
2	JX	479	ILE	Peptide
2	KX	479	ILE	Peptide
2	LX	479	ILE	Peptide
2	MX	479	ILE	Peptide
2	NX	479	ILE	Peptide

5.2 Too-close contacts (i)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	AT	1502	0	1525	21	0
1	BT	1502	0	1525	23	0
1	CT	1502	0	1525	23	0
1	DT	1502	0	1525	23	0
1	ET	1502	0	1525	20	0
1	FT	1502	0	1525	27	0
1	GT	1502	0	1525	21	0
1	HT	1502	0	1525	23	0
1	IT	1502	0	1525	23	0
1	JT	1502	0	1525	23	0
1	KT	1502	0	1525	26	0
1	LT	1502	0	1525	26	0
1	MT	1502	0	1525	24	0
1	NT	1502	0	1525	20	0
2	AX	1275	0	1309	26	0
2	BX	1275	0	1309	28	0
2	CX	1275	0	1309	26	0
2	DX	1275	0	1309	24	0
2	EX	1275	0	1309	27	0
2	FX	1275	0	1309	29	0
2	GX	1275	0	1309	31	0
2	HX	1275	0	1309	29	0
2	IX	1275	0	1309	25	0
2	JX	1275	0	1309	27	0
2	KX	1275	0	1309	21	0
2	LX	1275	0	1309	26	0
2	MX	1275	0	1309	27	0
2	NX	1275	0	1309	28	0
3	AY	1578	0	1637	45	0
3	BY	1578	0	1637	45	0
3	CY	1578	0	1637	48	0
3	DY	1578	0	1637	43	0
3	EY	1578	0	1637	42	0
3	FY	1578	0	1637	45	0
3	GY	1578	0	1637	44	0
3	HY	1578	0	1637	44	0
3	IY	1578	0	1637	46	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	JY	1578	0	1637	42	0
3	KY	1578	0	1637	42	0
3	LY	1578	0	1637	41	0
3	MY	1578	0	1637	43	0
3	NY	1578	0	1637	42	0
All	All	60970	0	62594	984	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 8.

All (984) 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)
2:BX:444:PRO:O	2:CX:410:ARG:NH2	2.22	0.73
2:HX:444:PRO:O	2:IX:410:ARG:NH2	2.21	0.73
3:AY:1722:VAL:HB	3:AY:1732:LEU:HB2	1.72	0.72
3:DY:1722:VAL:HB	3:DY:1732:LEU:HB2	1.72	0.72
3:FY:1722:VAL:HB	3:FY:1732:LEU:HB2	1.72	0.72
3:CY:1722:VAL:HB	3:CY:1732:LEU:HB2	1.72	0.72
3:GY:1722:VAL:HB	3:GY:1732:LEU:HB2	1.72	0.72
3:BY:1722:VAL:HB	3:BY:1732:LEU:HB2	1.72	0.72
3:EY:1722:VAL:HB	3:EY:1732:LEU:HB2	1.72	0.72
3:IY:1722:VAL:HB	3:IY:1732:LEU:HB2	1.72	0.72
3:MY:1722:VAL:HB	3:MY:1732:LEU:HB2	1.72	0.72
3:NY:1722:VAL:HB	3:NY:1732:LEU:HB2	1.72	0.72
3:HY:1722:VAL:HB	3:HY:1732:LEU:HB2	1.72	0.71
3:KY:1700:ILE:O	3:KY:1876:SER:HA	1.90	0.71
3:NY:1700:ILE:O	3:NY:1876:SER:HA	1.91	0.71
3:KY:1722:VAL:HB	3:KY:1732:LEU:HB2	1.72	0.71
3:JY:1722:VAL:HB	3:JY:1732:LEU:HB2	1.72	0.71
3:LY:1700:ILE:O	3:LY:1876:SER:HA	1.91	0.71
3:LY:1722:VAL:HB	3:LY:1732:LEU:HB2	1.72	0.71
3:AY:1700:ILE:O	3:AY:1876:SER:HA	1.91	0.71
3:CY:1700:ILE:O	3:CY:1876:SER:HA	1.90	0.71
3:IY:1700:ILE:O	3:IY:1876:SER:HA	1.90	0.70
2:JX:480:LYS:HB2	2:JX:483:ALA:HB3	1.74	0.70
2:LX:480:LYS:HB2	2:LX:483:ALA:HB3	1.74	0.70
2:DX:480:LYS:HB2	2:DX:483:ALA:HB3	1.74	0.70
3:FY:1700:ILE:O	3:FY:1876:SER:HA	1.91	0.70
2:BX:480:LYS:HB2	2:BX:483:ALA:HB3	1.74	0.70
2:CX:480:LYS:HB2	2:CX:483:ALA:HB3	1.74	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:EY:1700:ILE:O	3:EY:1876:SER:HA	1.90	0.70
2:IX:480:LYS:HB2	2:IX:483:ALA:HB3	1.74	0.70
2:NX:480:LYS:HB2	2:NX:483:ALA:HB3	1.74	0.70
2:AX:480:LYS:HB2	2:AX:483:ALA:HB3	1.74	0.70
3:BY:1700:ILE:O	3:BY:1876:SER:HA	1.90	0.70
3:JY:1700:ILE:O	3:JY:1876:SER:HA	1.90	0.70
3:MY:1700:ILE:O	3:MY:1876:SER:HA	1.91	0.70
3:HY:1700:ILE:O	3:HY:1876:SER:HA	1.90	0.70
2:KX:480:LYS:HB2	2:KX:483:ALA:HB3	1.74	0.70
2:FX:480:LYS:HB2	2:FX:483:ALA:HB3	1.74	0.70
2:GX:480:LYS:HB2	2:GX:483:ALA:HB3	1.74	0.70
3:GY:1700:ILE:O	3:GY:1876:SER:HA	1.91	0.70
2:EX:480:LYS:HB2	2:EX:483:ALA:HB3	1.74	0.70
2:MX:480:LYS:HB2	2:MX:483:ALA:HB3	1.74	0.70
3:DY:1700:ILE:O	3:DY:1876:SER:HA	1.91	0.69
2:HX:480:LYS:HB2	2:HX:483:ALA:HB3	1.74	0.69
2:KX:439:ILE:HG12	2:KX:478:LEU:HG	1.77	0.67
2:EX:439:ILE:HG12	2:EX:478:LEU:HG	1.77	0.67
2:FX:439:ILE:HG12	2:FX:478:LEU:HG	1.77	0.67
2:LX:439:ILE:HG12	2:LX:478:LEU:HG	1.77	0.67
2:HX:439:ILE:HG12	2:HX:478:LEU:HG	1.77	0.67
2:IX:439:ILE:HG12	2:IX:478:LEU:HG	1.77	0.67
2:NX:439:ILE:HG12	2:NX:478:LEU:HG	1.77	0.67
2:DX:439:ILE:HG12	2:DX:478:LEU:HG	1.77	0.67
2:KX:444:PRO:O	2:LX:410:ARG:NH2	2.28	0.67
2:BX:439:ILE:HG12	2:BX:478:LEU:HG	1.77	0.66
2:CX:439:ILE:HG12	2:CX:478:LEU:HG	1.77	0.66
2:MX:439:ILE:HG12	2:MX:478:LEU:HG	1.77	0.66
2:GX:439:ILE:HG12	2:GX:478:LEU:HG	1.77	0.66
2:JX:439:ILE:HG12	2:JX:478:LEU:HG	1.77	0.66
2:AX:439:ILE:HG12	2:AX:478:LEU:HG	1.77	0.66
3:EY:1857:GLN:NE2	3:FY:1856:ASN:OD1	2.30	0.65
3:IY:1857:GLN:NE2	3:JY:1856:ASN:OD1	2.30	0.64
3:KY:1857:GLN:NE2	3:LY:1856:ASN:OD1	2.30	0.64
3:GY:1857:GLN:NE2	3:HY:1856:ASN:OD1	2.31	0.64
3:HY:1857:GLN:NE2	3:IY:1856:ASN:OD1	2.31	0.64
2:AX:444:PRO:O	2:BX:410:ARG:NH2	2.30	0.64
3:LY:1857:GLN:NE2	3:MY:1856:ASN:OD1	2.30	0.64
3:DY:1857:GLN:NE2	3:EY:1856:ASN:OD1	2.31	0.64
3:AY:1856:ASN:OD1	3:NY:1857:GLN:NE2	2.30	0.63
1:ET:36:TYR:HB3	2:FX:404:TYR:HB3	1.78	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:FX:399:ARG:NH2	1:GT:169:HIS:O	2.31	0.63
3:FY:1857:GLN:NE2	3:GY:1856:ASN:OD1	2.31	0.63
3:JY:1857:GLN:NE2	3:KY:1856:ASN:OD1	2.32	0.63
3:BY:1857:GLN:NE2	3:CY:1856:ASN:OD1	2.31	0.63
3:AY:1857:GLN:NE2	3:BY:1856:ASN:OD1	2.31	0.62
2:JX:444:PRO:O	2:KX:410:ARG:NH2	2.32	0.62
3:MY:1857:GLN:NE2	3:NY:1856:ASN:OD1	2.32	0.62
3:CY:1857:GLN:NE2	3:DY:1856:ASN:OD1	2.32	0.61
1:JT:213:ARG:HH12	2:JX:364:ASN:HB3	1.64	0.61
1:LT:39:VAL:HG11	2:MX:419:ILE:HG13	1.82	0.61
2:EX:444:PRO:O	2:FX:410:ARG:NH2	2.33	0.61
3:NY:1714:VAL:HG21	3:NY:1757:ILE:HD13	1.82	0.61
1:CT:36:TYR:HB3	2:DX:404:TYR:HB3	1.83	0.61
3:LY:1714:VAL:HG21	3:LY:1757:ILE:HD13	1.82	0.61
3:MY:1714:VAL:HG21	3:MY:1757:ILE:HD13	1.83	0.61
3:KY:1714:VAL:HG21	3:KY:1757:ILE:HD13	1.82	0.61
1:IT:36:TYR:HB3	2:JX:404:TYR:HB3	1.82	0.60
1:LT:39:VAL:HG21	2:MX:419:ILE:HD12	1.82	0.60
1:MT:36:TYR:HB3	2:NX:404:TYR:HB3	1.82	0.60
3:AY:1714:VAL:HG21	3:AY:1757:ILE:HD13	1.83	0.60
3:JY:1714:VAL:HG21	3:JY:1757:ILE:HD13	1.83	0.60
3:LY:1762:ALA:HB3	3:LY:1770:ILE:HB	1.84	0.60
3:BY:1714:VAL:HG21	3:BY:1757:ILE:HD13	1.83	0.60
1:GT:36:TYR:HB3	2:HX:404:TYR:HB3	1.81	0.60
2:MX:444:PRO:O	2:NX:410:ARG:NH2	2.35	0.60
3:NY:1762:ALA:HB3	3:NY:1770:ILE:HB	1.84	0.60
2:CX:444:PRO:O	2:DX:410:ARG:NH2	2.34	0.60
2:FX:444:PRO:O	2:GX:410:ARG:NH2	2.34	0.60
2:GX:444:PRO:O	2:HX:410:ARG:NH2	2.35	0.60
2:LX:444:PRO:O	2:MX:410:ARG:NH2	2.34	0.60
3:BY:1762:ALA:HB3	3:BY:1770:ILE:HB	1.84	0.60
3:IY:1714:VAL:HG21	3:IY:1757:ILE:HD13	1.83	0.60
1:AT:36:TYR:HB3	2:BX:404:TYR:HB3	1.83	0.60
3:JY:1762:ALA:HB3	3:JY:1770:ILE:HB	1.84	0.60
3:MY:1762:ALA:HB3	3:MY:1770:ILE:HB	1.84	0.60
3:CY:1714:VAL:HG21	3:CY:1757:ILE:HD13	1.83	0.59
3:EY:1714:VAL:HG21	3:EY:1757:ILE:HD13	1.82	0.59
3:HY:1714:VAL:HG21	3:HY:1757:ILE:HD13	1.83	0.59
3:KY:1762:ALA:HB3	3:KY:1770:ILE:HB	1.84	0.59
2:LX:399:ARG:NH2	1:MT:169:HIS:O	2.35	0.59
1:ET:39:VAL:HG11	2:FX:419:ILE:HG13	1.83	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:IY:1762:ALA:HB3	3:IY:1770:ILE:HB	1.84	0.59
1:LT:36:TYR:HB3	2:MX:404:TYR:HB3	1.84	0.59
3:AY:1762:ALA:HB3	3:AY:1770:ILE:HB	1.84	0.59
3:DY:1714:VAL:HG21	3:DY:1757:ILE:HD13	1.83	0.59
3:GY:1714:VAL:HG21	3:GY:1757:ILE:HD13	1.82	0.59
2:BX:399:ARG:NH2	1:CT:169:HIS:O	2.34	0.59
3:FY:1714:VAL:HG21	3:FY:1757:ILE:HD13	1.83	0.59
3:DY:1762:ALA:HB3	3:DY:1770:ILE:HB	1.84	0.59
3:GY:1762:ALA:HB3	3:GY:1770:ILE:HB	1.84	0.59
1:KT:36:TYR:HB3	2:LX:404:TYR:HB3	1.85	0.59
1:MT:39:VAL:HG11	2:NX:419:ILE:HG13	1.84	0.58
3:DY:1706:VAL:HB	3:DY:1871:LYS:HB3	1.86	0.58
2:AX:410:ARG:NH2	2:NX:444:PRO:O	2.36	0.58
3:AY:1706:VAL:HB	3:AY:1871:LYS:HB3	1.86	0.58
3:BY:1706:VAL:HB	3:BY:1871:LYS:HB3	1.86	0.58
3:EY:1762:ALA:HB3	3:EY:1770:ILE:HB	1.84	0.58
2:IX:444:PRO:O	2:JX:410:ARG:NH2	2.37	0.58
3:EY:1706:VAL:HB	3:EY:1871:LYS:HB3	1.86	0.58
3:FY:1762:ALA:HB3	3:FY:1770:ILE:HB	1.84	0.58
3:CY:1706:VAL:HB	3:CY:1871:LYS:HB3	1.86	0.58
1:DT:36:TYR:HB3	2:EX:404:TYR:HB3	1.85	0.58
3:FY:1706:VAL:HB	3:FY:1871:LYS:HB3	1.86	0.58
2:AX:404:TYR:HB3	1:NT:36:TYR:HB3	1.84	0.58
3:HY:1762:ALA:HB3	3:HY:1770:ILE:HB	1.84	0.58
1:IT:39:VAL:HG11	2:JX:419:ILE:HG13	1.84	0.57
3:MY:1706:VAL:HB	3:MY:1871:LYS:HB3	1.86	0.57
3:NY:1706:VAL:HB	3:NY:1871:LYS:HB3	1.86	0.57
1:AT:169:HIS:O	2:NX:399:ARG:NH2	2.37	0.57
1:ET:39:VAL:HG21	2:FX:419:ILE:HD12	1.86	0.57
3:GY:1706:VAL:HB	3:GY:1871:LYS:HB3	1.86	0.57
3:CY:1762:ALA:HB3	3:CY:1770:ILE:HB	1.84	0.57
1:BT:36:TYR:HB3	2:CX:404:TYR:HB3	1.86	0.57
1:AT:39:VAL:HG11	2:BX:419:ILE:HG13	1.87	0.57
1:MT:39:VAL:HG21	2:NX:419:ILE:HD12	1.86	0.57
2:DX:399:ARG:NH2	1:ET:169:HIS:O	2.37	0.57
3:KY:1706:VAL:HB	3:KY:1871:LYS:HB3	1.86	0.57
2:DX:444:PRO:O	2:EX:410:ARG:NH2	2.37	0.57
3:IY:1706:VAL:HB	3:IY:1871:LYS:HB3	1.86	0.57
3:HY:1706:VAL:HB	3:HY:1871:LYS:HB3	1.86	0.56
3:LY:1706:VAL:HB	3:LY:1871:LYS:HB3	1.86	0.56
3:JY:1706:VAL:HB	3:JY:1871:LYS:HB3	1.86	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:MX:399:ARG:NH2	1:NT:169:HIS:O	2.39	0.56
2:HX:399:ARG:NH2	1:IT:169:HIS:O	2.37	0.56
1:KT:39:VAL:HG11	2:LX:419:ILE:HG13	1.86	0.56
1:IT:39:VAL:HG21	2:JX:419:ILE:HD12	1.87	0.56
2:AX:419:ILE:HG13	1:NT:39:VAL:HG11	1.88	0.56
3:HY:1732:LEU:HD22	3:HY:1736:THR:HG21	1.89	0.55
3:IY:1732:LEU:HD22	3:IY:1736:THR:HG21	1.89	0.55
3:JY:1732:LEU:HD22	3:JY:1736:THR:HG21	1.89	0.55
2:KX:507:GLY:HA2	1:LT:168:LEU:HD11	1.87	0.55
2:AX:399:ARG:NH2	1:BT:169:HIS:O	2.38	0.55
2:EX:399:ARG:NH2	1:FT:169:HIS:O	2.39	0.55
3:LY:1732:LEU:HD22	3:LY:1736:THR:HG21	1.88	0.55
3:GY:1732:LEU:HD22	3:GY:1736:THR:HG21	1.89	0.55
3:KY:1732:LEU:HD22	3:KY:1736:THR:HG21	1.89	0.55
3:MY:1732:LEU:HD22	3:MY:1736:THR:HG21	1.89	0.55
3:FY:1732:LEU:HD22	3:FY:1736:THR:HG21	1.89	0.54
3:NY:1732:LEU:HD22	3:NY:1736:THR:HG21	1.89	0.54
1:LT:296:GLU:OE2	1:LT:300:ASN:ND2	2.41	0.54
3:AY:1732:LEU:HD22	3:AY:1736:THR:HG21	1.89	0.54
1:ET:296:GLU:OE2	1:ET:300:ASN:ND2	2.41	0.54
1:FT:296:GLU:OE2	1:FT:300:ASN:ND2	2.41	0.54
3:BY:1732:LEU:HD22	3:BY:1736:THR:HG21	1.88	0.54
1:IT:296:GLU:OE2	1:IT:300:ASN:ND2	2.41	0.54
1:MT:296:GLU:OE2	1:MT:300:ASN:ND2	2.41	0.54
3:DY:1732:LEU:HD22	3:DY:1736:THR:HG21	1.88	0.54
3:EY:1732:LEU:HD22	3:EY:1736:THR:HG21	1.89	0.54
1:KT:296:GLU:OE2	1:KT:300:ASN:ND2	2.41	0.54
1:FT:39:VAL:HG21	2:GX:419:ILE:HD12	1.89	0.54
1:CT:296:GLU:OE2	1:CT:300:ASN:ND2	2.41	0.54
3:CY:1732:LEU:HD22	3:CY:1736:THR:HG21	1.89	0.54
3:HY:1859:LEU:O	3:HY:1863:MET:CB	2.56	0.54
3:AY:1710:LEU:HD12	3:BY:1785:GLY:HA3	1.90	0.54
3:CY:1710:LEU:HD12	3:DY:1785:GLY:HA3	1.90	0.54
3:GY:1859:LEU:O	3:GY:1863:MET:CB	2.56	0.54
1:HT:296:GLU:OE2	1:HT:300:ASN:ND2	2.41	0.54
3:IY:1859:LEU:O	3:IY:1863:MET:CB	2.56	0.54
1:NT:296:GLU:OE2	1:NT:300:ASN:ND2	2.41	0.54
1:BT:296:GLU:OE2	1:BT:300:ASN:ND2	2.41	0.54
2:IX:399:ARG:NH2	1:JT:169:HIS:O	2.41	0.54
2:JX:399:ARG:NH2	1:KT:169:HIS:O	2.41	0.54
3:JY:1710:LEU:HD12	3:KY:1785:GLY:HA3	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:AY:1859:LEU:O	3:AY:1863:MET:CB	2.56	0.53
1:JT:296:GLU:OE2	1:JT:300:ASN:ND2	2.41	0.53
3:JY:1859:LEU:O	3:JY:1863:MET:CB	2.56	0.53
3:LY:1710:LEU:HD12	3:MY:1785:GLY:HA3	1.90	0.53
2:CX:399:ARG:NH2	1:DT:169:HIS:O	2.40	0.53
3:CY:1859:LEU:O	3:CY:1863:MET:CB	2.56	0.53
1:FT:39:VAL:HG11	2:GX:419:ILE:HG13	1.88	0.53
1:GT:296:GLU:OE2	1:GT:300:ASN:ND2	2.41	0.53
2:KX:399:ARG:NH2	1:LT:169:HIS:O	2.40	0.53
1:DT:39:VAL:HG11	2:EX:419:ILE:HG13	1.89	0.53
3:DY:1859:LEU:O	3:DY:1863:MET:CB	2.56	0.53
3:FY:1859:LEU:O	3:FY:1863:MET:CB	2.56	0.53
1:AT:296:GLU:OE2	1:AT:300:ASN:ND2	2.41	0.53
3:AY:1785:GLY:HA3	3:NY:1710:LEU:HD12	1.91	0.53
3:BY:1859:LEU:O	3:BY:1863:MET:CB	2.56	0.53
1:DT:296:GLU:OE2	1:DT:300:ASN:ND2	2.41	0.53
1:KT:39:VAL:HG21	2:LX:419:ILE:HD12	1.89	0.53
2:BX:507:GLY:HA2	1:CT:168:LEU:HD11	1.90	0.53
3:KY:1710:LEU:HD12	3:LY:1785:GLY:HA3	1.91	0.53
3:KY:1859:LEU:O	3:KY:1863:MET:CB	2.56	0.53
2:GX:376:ILE:HA	2:GX:379:GLN:HG2	1.91	0.53
3:MY:1859:LEU:O	3:MY:1863:MET:CB	2.56	0.53
2:AX:376:ILE:HA	2:AX:379:GLN:HG2	1.91	0.53
2:BX:376:ILE:HA	2:BX:379:GLN:HG2	1.91	0.53
2:FX:376:ILE:HA	2:FX:379:GLN:HG2	1.91	0.53
2:GX:399:ARG:NH2	1:HT:169:HIS:O	2.42	0.53
3:GY:1710:LEU:HD12	3:HY:1785:GLY:HA3	1.90	0.53
1:HT:39:VAL:HG11	2:IX:419:ILE:HG13	1.90	0.53
3:HY:1858:ILE:HG13	3:IY:1859:LEU:HD21	1.91	0.53
2:MX:376:ILE:HA	2:MX:379:GLN:HG2	1.91	0.53
3:MY:1710:LEU:HD12	3:NY:1785:GLY:HA3	1.91	0.53
2:DX:376:ILE:HA	2:DX:379:GLN:HG2	1.91	0.53
2:EX:376:ILE:HA	2:EX:379:GLN:HG2	1.91	0.53
1:HT:111:ALA:O	1:IT:167:ASN:ND2	2.42	0.53
1:LT:114:ALA:HB1	1:LT:118:ILE:HD11	1.91	0.53
1:CT:114:ALA:HB1	1:CT:118:ILE:HD11	1.91	0.52
1:DT:114:ALA:HB1	1:DT:118:ILE:HD11	1.92	0.52
3:HY:1710:LEU:HD12	3:IY:1785:GLY:HA3	1.90	0.52
2:IX:376:ILE:HA	2:IX:379:GLN:HG2	1.91	0.52
1:JT:210:TYR:CD1	2:JX:368:ILE:HG13	2.45	0.52
2:KX:376:ILE:HA	2:KX:379:GLN:HG2	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:LX:376:ILE:HA	2:LX:379:GLN:HG2	1.91	0.52
2:CX:376:ILE:HA	2:CX:379:GLN:HG2	1.91	0.52
3:DY:1710:LEU:HD12	3:EY:1785:GLY:HA3	1.90	0.52
3:EY:1859:LEU:O	3:EY:1863:MET:CB	2.56	0.52
1:FT:114:ALA:HB1	1:FT:118:ILE:HD11	1.92	0.52
2:HX:376:ILE:HA	2:HX:379:GLN:HG2	1.91	0.52
3:LY:1859:LEU:O	3:LY:1863:MET:CB	2.56	0.52
1:NT:114:ALA:HB1	1:NT:118:ILE:HD11	1.92	0.52
2:NX:376:ILE:HA	2:NX:379:GLN:HG2	1.91	0.52
3:NY:1859:LEU:O	3:NY:1863:MET:CB	2.56	0.52
1:AT:114:ALA:HB1	1:AT:118:ILE:HD11	1.92	0.52
1:BT:111:ALA:O	1:CT:167:ASN:ND2	2.43	0.52
1:ET:114:ALA:HB1	1:ET:118:ILE:HD11	1.91	0.52
1:FT:36:TYR:HB3	2:GX:404:TYR:HB3	1.91	0.52
1:GT:90:SER:HA	1:GT:137:ASN:HB2	1.92	0.52
1:HT:90:SER:HA	1:HT:137:ASN:HB2	1.92	0.52
1:MT:114:ALA:HB1	1:MT:118:ILE:HD11	1.91	0.52
3:EY:1858:ILE:HG13	3:FY:1859:LEU:HD21	1.92	0.52
1:GT:39:VAL:HG11	2:HX:419:ILE:HG13	1.91	0.52
1:HT:114:ALA:HB1	1:HT:118:ILE:HD11	1.91	0.52
1:JT:90:SER:HA	1:JT:137:ASN:HB2	1.92	0.52
1:JT:114:ALA:HB1	1:JT:118:ILE:HD11	1.92	0.52
2:JX:376:ILE:HA	2:JX:379:GLN:HG2	1.91	0.52
1:KT:90:SER:HA	1:KT:137:ASN:HB2	1.92	0.52
3:LY:1858:ILE:HG13	3:MY:1859:LEU:HD21	1.92	0.52
3:EY:1710:LEU:HD12	3:FY:1785:GLY:HA3	1.91	0.52
3:AY:1859:LEU:HD21	3:NY:1858:ILE:HG13	1.92	0.52
3:AY:1858:ILE:HG13	3:BY:1859:LEU:HD21	1.92	0.52
3:BY:1710:LEU:HD12	3:CY:1785:GLY:HA3	1.92	0.52
1:GT:114:ALA:HB1	1:GT:118:ILE:HD11	1.91	0.52
1:KT:114:ALA:HB1	1:KT:118:ILE:HD11	1.92	0.52
1:BT:114:ALA:HB1	1:BT:118:ILE:HD11	1.92	0.52
1:ET:111:ALA:O	1:FT:167:ASN:ND2	2.43	0.52
1:IT:114:ALA:HB1	1:IT:118:ILE:HD11	1.91	0.52
1:MT:90:SER:HA	1:MT:137:ASN:HB2	1.92	0.52
1:FT:90:SER:HA	1:FT:137:ASN:HB2	1.92	0.52
1:IT:90:SER:HA	1:IT:137:ASN:HB2	1.92	0.52
1:ET:90:SER:HA	1:ET:137:ASN:HB2	1.92	0.52
1:ET:193:GLU:HA	1:ET:196:ASN:HD22	1.75	0.52
2:KX:439:ILE:HD12	2:KX:467:TYR:HD1	1.75	0.52
1:LT:90:SER:HA	1:LT:137:ASN:HB2	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AT:39:VAL:HG21	2:BX:419:ILE:HD12	1.91	0.51
3:KY:1858:ILE:HG13	3:LY:1859:LEU:HD21	1.92	0.51
2:MX:439:ILE:HD12	2:MX:467:TYR:HD1	1.76	0.51
1:FT:193:GLU:HA	1:FT:196:ASN:HD22	1.75	0.51
1:IT:213:ARG:HH12	2:IX:364:ASN:HB3	1.76	0.51
1:KT:193:GLU:HA	1:KT:196:ASN:HD22	1.75	0.51
1:DT:193:GLU:HA	1:DT:196:ASN:HD22	1.75	0.51
3:GY:1858:ILE:HG13	3:HY:1859:LEU:HD21	1.93	0.51
3:IY:1710:LEU:HD12	3:JY:1785:GLY:HA3	1.92	0.51
1:KT:111:ALA:O	1:LT:167:ASN:ND2	2.44	0.51
2:AX:439:ILE:HD12	2:AX:467:TYR:HD1	1.76	0.51
1:DT:90:SER:HA	1:DT:137:ASN:HB2	1.92	0.51
1:JT:193:GLU:HA	1:JT:196:ASN:HD22	1.75	0.51
1:NT:90:SER:HA	1:NT:137:ASN:HB2	1.92	0.51
1:CT:90:SER:HA	1:CT:137:ASN:HB2	1.92	0.51
3:FY:1710:LEU:HD12	3:GY:1785:GLY:HA3	1.92	0.51
3:FY:1858:ILE:HG13	3:GY:1859:LEU:HD21	1.93	0.51
1:GT:193:GLU:HA	1:GT:196:ASN:HD22	1.75	0.51
2:IX:436:GLN:HB2	3:IY:1734:LYS:HG2	1.93	0.51
3:IY:1858:ILE:HG13	3:JY:1859:LEU:HD21	1.93	0.51
1:AT:193:GLU:HA	1:AT:196:ASN:HD22	1.75	0.51
3:DY:1858:ILE:HG13	3:EY:1859:LEU:HD21	1.92	0.51
2:IX:439:ILE:HD12	2:IX:467:TYR:HD1	1.76	0.51
1:AT:90:SER:HA	1:AT:137:ASN:HB2	1.92	0.51
2:AX:419:ILE:HD12	1:NT:39:VAL:HG21	1.91	0.51
1:BT:193:GLU:HA	1:BT:196:ASN:HD22	1.75	0.51
1:CT:193:GLU:HA	1:CT:196:ASN:HD22	1.75	0.51
1:IT:210:TYR:CD1	2:IX:368:ILE:HG13	2.46	0.51
1:MT:111:ALA:O	1:NT:167:ASN:ND2	2.44	0.51
3:MY:1858:ILE:HG13	3:NY:1859:LEU:HD21	1.93	0.51
1:DT:39:VAL:HG21	2:EX:419:ILE:HD12	1.92	0.51
1:LT:193:GLU:HA	1:LT:196:ASN:HD22	1.75	0.51
1:AT:111:ALA:O	1:BT:167:ASN:ND2	2.44	0.50
1:IT:193:GLU:HA	1:IT:196:ASN:HD22	1.75	0.50
1:BT:90:SER:HA	1:BT:137:ASN:HB2	1.92	0.50
3:CY:1858:ILE:HG13	3:DY:1859:LEU:HD21	1.93	0.50
1:DT:111:ALA:O	1:ET:167:ASN:ND2	2.44	0.50
2:EX:439:ILE:HD12	2:EX:467:TYR:HD1	1.75	0.50
2:FX:439:ILE:HD12	2:FX:467:TYR:HD1	1.76	0.50
2:HX:439:ILE:HD12	2:HX:467:TYR:HD1	1.76	0.50
3:JY:1858:ILE:HG13	3:KY:1859:LEU:HD21	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:CX:439:ILE:HD12	2:CX:467:TYR:HD1	1.76	0.50
3:CY:1859:LEU:O	3:CY:1863:MET:HB2	2.11	0.50
1:HT:39:VAL:HG21	2:IX:419:ILE:HD12	1.92	0.50
2:JX:439:ILE:HD12	2:JX:467:TYR:HD1	1.76	0.50
2:LX:439:ILE:HD12	2:LX:467:TYR:HD1	1.76	0.50
3:BY:1858:ILE:HG13	3:CY:1859:LEU:HD21	1.93	0.50
1:HT:193:GLU:HA	1:HT:196:ASN:HD22	1.75	0.50
2:IX:507:GLY:HA2	1:JT:168:LEU:HD11	1.93	0.50
3:NY:1859:LEU:O	3:NY:1863:MET:HB2	2.11	0.50
2:AX:497:LEU:O	2:BX:432:ASN:ND2	2.37	0.50
3:BY:1859:LEU:O	3:BY:1863:MET:HB2	2.11	0.50
1:FT:111:ALA:O	1:GT:167:ASN:ND2	2.45	0.50
3:FY:1740:GLY:HA2	3:FY:1759:PHE:HA	1.94	0.50
3:HY:1859:LEU:O	3:HY:1863:MET:HB2	2.11	0.50
3:IY:1763:ILE:HG12	3:IY:1769:ILE:HG22	1.94	0.50
3:JY:1859:LEU:O	3:JY:1863:MET:HB2	2.11	0.50
3:KY:1859:LEU:O	3:KY:1863:MET:HB2	2.11	0.50
1:NT:193:GLU:HA	1:NT:196:ASN:HD22	1.75	0.50
3:AY:1859:LEU:O	3:AY:1863:MET:HB2	2.11	0.50
3:EY:1740:GLY:HA2	3:EY:1759:PHE:HA	1.94	0.50
3:GY:1740:GLY:HA2	3:GY:1759:PHE:HA	1.94	0.50
3:GY:1763:ILE:HG12	3:GY:1769:ILE:HG22	1.94	0.50
3:KY:1763:ILE:HG12	3:KY:1769:ILE:HG22	1.94	0.50
3:AY:1740:GLY:HA2	3:AY:1759:PHE:HA	1.94	0.50
1:BT:39:VAL:HG11	2:CX:419:ILE:HG13	1.93	0.50
2:BX:436:GLN:HB2	3:BY:1734:LYS:HG2	1.93	0.50
2:DX:439:ILE:HD12	2:DX:467:TYR:HD1	1.76	0.50
3:EY:1763:ILE:HG12	3:EY:1769:ILE:HG22	1.94	0.50
2:NX:439:ILE:HD12	2:NX:467:TYR:HD1	1.76	0.50
3:NY:1740:GLY:HA2	3:NY:1759:PHE:HA	1.94	0.50
1:AT:167:ASN:ND2	1:NT:111:ALA:O	2.45	0.49
3:BY:1740:GLY:HA2	3:BY:1759:PHE:HA	1.94	0.49
3:DY:1740:GLY:HA2	3:DY:1759:PHE:HA	1.94	0.49
3:HY:1740:GLY:HA2	3:HY:1759:PHE:HA	1.94	0.49
3:HY:1763:ILE:HG12	3:HY:1769:ILE:HG22	1.94	0.49
3:IY:1859:LEU:O	3:IY:1863:MET:HB2	2.12	0.49
3:LY:1859:LEU:O	3:LY:1863:MET:HB2	2.12	0.49
3:MY:1740:GLY:HA2	3:MY:1759:PHE:HA	1.94	0.49
1:AT:28:VAL:HG23	1:AT:29:VAL:H	1.78	0.49
3:CY:1740:GLY:HA2	3:CY:1759:PHE:HA	1.94	0.49
2:FX:399:ARG:NH1	2:FX:495:ASN:OD1	2.46	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:HT:36:TYR:HB3	2:IX:404:TYR:HB3	1.92	0.49
2:IX:455:ILE:HB	3:IY:1765:PRO:HB3	1.93	0.49
2:IX:481:ASP:HB3	3:IY:1875:ASP:HA	1.95	0.49
3:JY:1763:ILE:HG12	3:JY:1769:ILE:HG22	1.94	0.49
3:LY:1763:ILE:HG12	3:LY:1769:ILE:HG22	1.94	0.49
2:MX:399:ARG:NH1	2:MX:495:ASN:OD1	2.46	0.49
2:EX:399:ARG:NH1	2:EX:495:ASN:OD1	2.46	0.49
3:EY:1859:LEU:O	3:EY:1863:MET:HB2	2.11	0.49
2:GX:439:ILE:HD12	2:GX:467:TYR:HD1	1.75	0.49
3:LY:1740:GLY:HA2	3:LY:1759:PHE:HA	1.94	0.49
3:FY:1763:ILE:HG12	3:FY:1769:ILE:HG22	1.94	0.49
2:GX:399:ARG:NH1	2:GX:495:ASN:OD1	2.46	0.49
1:IT:111:ALA:O	1:JT:167:ASN:ND2	2.45	0.49
3:IY:1740:GLY:HA2	3:IY:1759:PHE:HA	1.94	0.49
1:LT:111:ALA:O	1:MT:167:ASN:ND2	2.45	0.49
3:NY:1763:ILE:HG12	3:NY:1769:ILE:HG22	1.94	0.49
1:BT:28:VAL:HG23	1:BT:29:VAL:H	1.78	0.49
2:BX:399:ARG:NH1	2:BX:495:ASN:OD1	2.46	0.49
2:BX:439:ILE:HD12	2:BX:467:TYR:HD1	1.76	0.49
2:CX:399:ARG:NH1	2:CX:495:ASN:OD1	2.46	0.49
3:CY:1763:ILE:HG12	3:CY:1769:ILE:HG22	1.94	0.49
1:FT:28:VAL:HG23	1:FT:29:VAL:H	1.78	0.49
1:MT:193:GLU:HA	1:MT:196:ASN:HD22	1.75	0.49
1:NT:28:VAL:HG23	1:NT:29:VAL:H	1.78	0.49
2:NX:399:ARG:NH1	2:NX:495:ASN:OD1	2.46	0.49
2:GX:400:ASN:HB3	2:GX:421:ASP:OD1	2.13	0.49
2:JX:399:ARG:NH1	2:JX:495:ASN:OD1	2.46	0.49
3:KY:1740:GLY:HA2	3:KY:1759:PHE:HA	1.94	0.49
3:MY:1859:LEU:O	3:MY:1863:MET:HB2	2.12	0.49
2:DX:400:ASN:HB3	2:DX:421:ASP:OD1	2.13	0.49
1:GT:39:VAL:HG21	2:HX:419:ILE:HD12	1.95	0.49
3:GY:1859:LEU:O	3:GY:1863:MET:HB2	2.11	0.49
2:HX:400:ASN:HB3	2:HX:421:ASP:OD1	2.13	0.49
3:JY:1740:GLY:HA2	3:JY:1759:PHE:HA	1.94	0.49
1:KT:196:ASN:OD1	3:LY:1907:GLN:NE2	2.46	0.49
2:DX:399:ARG:NH1	2:DX:495:ASN:OD1	2.46	0.49
1:ET:28:VAL:HG23	1:ET:29:VAL:H	1.78	0.49
2:CX:481:ASP:HB3	3:CY:1875:ASP:HA	1.94	0.49
3:EY:1699:GLU:HG2	3:EY:1719:ALA:HB3	1.95	0.49
3:FY:1699:GLU:HG2	3:FY:1719:ALA:HB3	1.95	0.49
3:GY:1699:GLU:HG2	3:GY:1719:ALA:HB3	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:KX:399:ARG:NH1	2:KX:495:ASN:OD1	2.46	0.49
1:CT:28:VAL:HG23	1:CT:29:VAL:H	1.78	0.49
3:FY:1859:LEU:O	3:FY:1863:MET:HB2	2.11	0.49
1:GT:28:VAL:HG23	1:GT:29:VAL:H	1.78	0.49
2:HX:399:ARG:NH1	2:HX:495:ASN:OD1	2.46	0.49
1:JT:99:ALA:HB2	1:JT:124:PRO:HB3	1.95	0.49
1:JT:213:ARG:NH1	2:JX:364:ASN:HB3	2.28	0.49
1:KT:28:VAL:HG23	1:KT:29:VAL:H	1.78	0.49
3:MY:1698:ILE:HG23	3:MY:1722:VAL:HG21	1.95	0.49
3:AY:1698:ILE:HG23	3:AY:1722:VAL:HG21	1.95	0.48
2:CX:400:ASN:HB3	2:CX:421:ASP:OD1	2.13	0.48
3:DY:1699:GLU:HG2	3:DY:1719:ALA:HB3	1.95	0.48
3:HY:1699:GLU:HG2	3:HY:1719:ALA:HB3	1.95	0.48
2:JX:507:GLY:HA2	1:KT:168:LEU:HD11	1.94	0.48
2:LX:399:ARG:NH1	2:LX:495:ASN:OD1	2.46	0.48
3:CY:1698:ILE:HG23	3:CY:1722:VAL:HG21	1.95	0.48
3:EY:1698:ILE:HG23	3:EY:1722:VAL:HG21	1.95	0.48
1:HT:99:ALA:HB2	1:HT:124:PRO:HB3	1.95	0.48
3:HY:1698:ILE:HG23	3:HY:1722:VAL:HG21	1.95	0.48
1:JT:28:VAL:HG23	1:JT:29:VAL:H	1.78	0.48
3:KY:1698:ILE:HG23	3:KY:1722:VAL:HG21	1.95	0.48
2:MX:400:ASN:HB3	2:MX:421:ASP:OD1	2.13	0.48
2:NX:400:ASN:HB3	2:NX:421:ASP:OD1	2.13	0.48
3:DY:1859:LEU:O	3:DY:1863:MET:HB2	2.12	0.48
2:EX:400:ASN:HB3	2:EX:421:ASP:OD1	2.13	0.48
3:FY:1698:ILE:HG23	3:FY:1722:VAL:HG21	1.95	0.48
1:IT:28:VAL:HG23	1:IT:29:VAL:H	1.78	0.48
3:IY:1699:GLU:HG2	3:IY:1719:ALA:HB3	1.95	0.48
2:JX:364:ASN:HA	2:JX:367:LYS:HG2	1.96	0.48
2:LX:400:ASN:HB3	2:LX:421:ASP:OD1	2.13	0.48
3:LY:1873:GLU:HG2	3:MY:1781:LEU:HD22	1.95	0.48
2:AX:400:ASN:HB3	2:AX:421:ASP:OD1	2.13	0.48
3:BY:1763:ILE:HG12	3:BY:1769:ILE:HG22	1.94	0.48
2:DX:364:ASN:HA	2:DX:367:LYS:HG2	1.96	0.48
1:FT:99:ALA:HB2	1:FT:124:PRO:HB3	1.95	0.48
2:KX:400:ASN:HB3	2:KX:421:ASP:OD1	2.13	0.48
2:AX:364:ASN:HA	2:AX:367:LYS:HG2	1.96	0.48
2:AX:399:ARG:NH1	2:AX:495:ASN:OD1	2.46	0.48
3:CY:1699:GLU:HG2	3:CY:1719:ALA:HB3	1.95	0.48
2:FX:364:ASN:HA	2:FX:367:LYS:HG2	1.96	0.48
2:HX:364:ASN:HA	2:HX:367:LYS:HG2	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:HX:497:LEU:O	2:IX:432:ASN:ND2	2.38	0.48
2:IX:400:ASN:HB3	2:IX:421:ASP:OD1	2.13	0.48
3:IY:1873:GLU:HG2	3:JY:1781:LEU:HD22	1.95	0.48
3:JY:1698:ILE:HG23	3:JY:1722:VAL:HG21	1.95	0.48
1:MT:28:VAL:HG23	1:MT:29:VAL:H	1.78	0.48
3:EY:1792:ASN:O	3:EY:1797:ARG:NH1	2.46	0.48
2:FX:400:ASN:HB3	2:FX:421:ASP:OD1	2.13	0.48
3:FY:1792:ASN:O	3:FY:1797:ARG:NH1	2.46	0.48
3:JY:1699:GLU:HG2	3:JY:1719:ALA:HB3	1.95	0.48
1:LT:99:ALA:HB2	1:LT:124:PRO:HB3	1.95	0.48
2:LX:364:ASN:HA	2:LX:367:LYS:HG2	1.96	0.48
3:MY:1763:ILE:HG12	3:MY:1769:ILE:HG22	1.94	0.48
2:BX:400:ASN:HB3	2:BX:421:ASP:OD1	2.13	0.48
3:JY:1792:ASN:O	3:JY:1797:ARG:NH1	2.46	0.48
3:AY:1713:ILE:HD11	3:BY:1880:LEU:HD23	1.96	0.48
3:CY:1873:GLU:HG2	3:DY:1781:LEU:HD22	1.95	0.48
3:GY:1698:ILE:HG23	3:GY:1722:VAL:HG21	1.95	0.48
3:HY:1713:ILE:HD11	3:IY:1880:LEU:HD23	1.96	0.48
2:IX:399:ARG:NH1	2:IX:495:ASN:OD1	2.46	0.48
2:JX:400:ASN:HB3	2:JX:421:ASP:OD1	2.13	0.48
3:MY:1792:ASN:O	3:MY:1797:ARG:NH1	2.46	0.48
3:BY:1699:GLU:HG2	3:BY:1719:ALA:HB3	1.95	0.48
2:CX:364:ASN:HA	2:CX:367:LYS:HG2	1.96	0.48
3:DY:1698:ILE:HG23	3:DY:1722:VAL:HG21	1.95	0.48
3:GY:1792:ASN:O	3:GY:1797:ARG:NH1	2.46	0.48
2:MX:364:ASN:HA	2:MX:367:LYS:HG2	1.96	0.48
3:NY:1792:ASN:O	3:NY:1797:ARG:NH1	2.46	0.48
3:AY:1763:ILE:HG12	3:AY:1769:ILE:HG22	1.94	0.48
3:KY:1699:GLU:HG2	3:KY:1719:ALA:HB3	1.95	0.48
3:LY:1792:ASN:O	3:LY:1797:ARG:NH1	2.46	0.48
3:DY:1792:ASN:O	3:DY:1797:ARG:NH1	2.46	0.47
3:EY:1873:GLU:HG2	3:FY:1781:LEU:HD22	1.96	0.47
1:LT:28:VAL:HG23	1:LT:29:VAL:H	1.78	0.47
3:LY:1707:ASP:HB2	3:MY:1783:GLU:HB2	1.96	0.47
3:AY:1781:LEU:HD22	3:NY:1873:GLU:HG2	1.95	0.47
3:AY:1792:ASN:O	3:AY:1797:ARG:NH1	2.46	0.47
3:KY:1792:ASN:O	3:KY:1797:ARG:NH1	2.46	0.47
2:BX:364:ASN:HA	2:BX:367:LYS:HG2	1.96	0.47
3:BY:1698:ILE:HG23	3:BY:1722:VAL:HG21	1.95	0.47
1:DT:28:VAL:HG23	1:DT:29:VAL:H	1.78	0.47
3:DY:1763:ILE:HG12	3:DY:1769:ILE:HG22	1.94	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:HT:28:VAL:HG23	1:HT:29:VAL:H	1.78	0.47
2:IX:364:ASN:HA	2:IX:367:LYS:HG2	1.96	0.47
3:IY:1698:ILE:HG23	3:IY:1722:VAL:HG21	1.95	0.47
3:LY:1698:ILE:HG23	3:LY:1722:VAL:HG21	1.95	0.47
1:NT:52:LEU:HD12	1:NT:89:HIS:CD2	2.49	0.47
3:NY:1698:ILE:HG23	3:NY:1722:VAL:HG21	1.95	0.47
1:AT:99:ALA:HB2	1:AT:124:PRO:HB3	1.95	0.47
3:AY:1699:GLU:HG2	3:AY:1719:ALA:HB3	1.95	0.47
1:BT:48:SER:OG	1:CT:170:GLY:O	2.22	0.47
3:BY:1873:GLU:HG2	3:CY:1781:LEU:HD22	1.95	0.47
1:CT:99:ALA:HB2	1:CT:124:PRO:HB3	1.95	0.47
1:IT:306:LEU:HD22	2:IX:384:GLU:HB2	1.97	0.47
3:IY:1792:ASN:O	3:IY:1797:ARG:NH1	2.46	0.47
3:LY:1699:GLU:HG2	3:LY:1719:ALA:HB3	1.95	0.47
1:MT:52:LEU:HD12	1:MT:89:HIS:CD2	2.49	0.47
3:AY:1783:GLU:HB2	3:NY:1707:ASP:HB2	1.96	0.47
1:CT:52:LEU:HD12	1:CT:89:HIS:CD2	2.49	0.47
2:GX:364:ASN:HA	2:GX:367:LYS:HG2	1.96	0.47
3:GY:1713:ILE:HD11	3:HY:1880:LEU:HD23	1.97	0.47
3:JY:1873:GLU:HG2	3:KY:1781:LEU:HD22	1.96	0.47
1:LT:33:ASN:ND2	2:MX:407:PRO:O	2.47	0.47
1:AT:52:LEU:HD12	1:AT:89:HIS:CD2	2.49	0.47
3:BY:1713:ILE:HD11	3:CY:1880:LEU:HD23	1.97	0.47
3:BY:1792:ASN:O	3:BY:1797:ARG:NH1	2.46	0.47
1:DT:52:LEU:HD12	1:DT:89:HIS:CD2	2.49	0.47
2:EX:364:ASN:HA	2:EX:367:LYS:HG2	1.96	0.47
1:LT:52:LEU:HD12	1:LT:89:HIS:CD2	2.49	0.47
1:BT:52:LEU:HD12	1:BT:89:HIS:CD2	2.49	0.47
1:BT:99:ALA:HB2	1:BT:124:PRO:HB3	1.95	0.47
1:ET:52:LEU:HD12	1:ET:89:HIS:CD2	2.49	0.47
1:FT:199:TYR:HB3	2:FX:379:GLN:OE1	2.15	0.47
3:FY:1873:GLU:HG2	3:GY:1781:LEU:HD22	1.96	0.47
1:GT:99:ALA:HB2	1:GT:124:PRO:HB3	1.95	0.47
1:IT:52:LEU:HD12	1:IT:89:HIS:CD2	2.49	0.47
1:IT:99:ALA:HB2	1:IT:124:PRO:HB3	1.95	0.47
1:KT:52:LEU:HD12	1:KT:89:HIS:CD2	2.50	0.47
3:MY:1699:GLU:HG2	3:MY:1719:ALA:HB3	1.95	0.47
1:NT:99:ALA:HB2	1:NT:124:PRO:HB3	1.95	0.47
3:NY:1699:GLU:HG2	3:NY:1719:ALA:HB3	1.95	0.47
1:DT:99:ALA:HB2	1:DT:124:PRO:HB3	1.95	0.47
1:FT:52:LEU:HD12	1:FT:89:HIS:CD2	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:FT:60:PRO:HA	1:FT:137:ASN:O	2.15	0.47
1:GT:60:PRO:HA	1:GT:137:ASN:O	2.15	0.47
1:HT:52:LEU:HD12	1:HT:89:HIS:CD2	2.49	0.47
1:IT:60:PRO:HA	1:IT:137:ASN:O	2.15	0.47
1:JT:52:LEU:HD12	1:JT:89:HIS:CD2	2.49	0.47
1:KT:99:ALA:HB2	1:KT:124:PRO:HB3	1.95	0.47
2:BX:480:LYS:NZ	3:BY:1699:GLU:OE2	2.38	0.47
3:DY:1713:ILE:HD11	3:EY:1880:LEU:HD23	1.96	0.47
1:ET:60:PRO:HA	1:ET:137:ASN:O	2.15	0.47
1:ET:99:ALA:HB2	1:ET:124:PRO:HB3	1.95	0.47
2:GX:477:LYS:HD3	2:GX:484:LEU:HD11	1.96	0.47
1:HT:60:PRO:HA	1:HT:137:ASN:O	2.15	0.47
3:HY:1792:ASN:O	3:HY:1797:ARG:NH1	2.46	0.47
1:MT:99:ALA:HB2	1:MT:124:PRO:HB3	1.95	0.47
3:MY:1873:GLU:HG2	3:NY:1781:LEU:HD22	1.97	0.47
3:CY:1713:ILE:HD11	3:DY:1880:LEU:HD23	1.97	0.47
3:CY:1792:ASN:O	3:CY:1797:ARG:NH1	2.46	0.47
1:DT:60:PRO:HA	1:DT:137:ASN:O	2.15	0.47
2:EX:507:GLY:HA2	1:FT:168:LEU:HD11	1.97	0.47
1:GT:52:LEU:HD12	1:GT:89:HIS:CD2	2.49	0.47
2:HX:477:LYS:HD3	2:HX:484:LEU:HD11	1.96	0.47
1:JT:60:PRO:HA	1:JT:137:ASN:O	2.15	0.47
3:JY:1707:ASP:HB2	3:KY:1783:GLU:HB2	1.96	0.47
2:KX:364:ASN:HA	2:KX:367:LYS:HG2	1.96	0.47
2:NX:364:ASN:HA	2:NX:367:LYS:HG2	1.96	0.47
1:CT:60:PRO:HA	1:CT:137:ASN:O	2.15	0.46
3:GY:1707:ASP:HB2	3:HY:1783:GLU:HB2	1.96	0.46
2:IX:477:LYS:HD3	2:IX:484:LEU:HD11	1.96	0.46
1:LT:60:PRO:HA	1:LT:137:ASN:O	2.15	0.46
1:GT:291:GLU:O	1:GT:295:ASN:HB2	2.16	0.46
2:JX:477:LYS:HD3	2:JX:484:LEU:HD11	1.96	0.46
3:MY:1707:ASP:HB2	3:NY:1783:GLU:HB2	1.97	0.46
2:FX:417:SER:N	2:FX:429:GLY:O	2.44	0.46
2:FX:477:LYS:HD3	2:FX:484:LEU:HD11	1.96	0.46
3:FY:1713:ILE:HD11	3:GY:1880:LEU:HD23	1.97	0.46
3:IY:1713:ILE:HD11	3:JY:1880:LEU:HD23	1.97	0.46
1:KT:291:GLU:O	1:KT:295:ASN:HB2	2.16	0.46
1:LT:291:GLU:O	1:LT:295:ASN:HB2	2.16	0.46
1:AT:60:PRO:HA	1:AT:137:ASN:O	2.15	0.46
1:CT:111:ALA:O	1:DT:167:ASN:ND2	2.49	0.46
3:CY:1707:ASP:HB2	3:DY:1783:GLU:HB2	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:EX:477:LYS:HD3	2:EX:484:LEU:HD11	1.96	0.46
1:FT:291:GLU:O	1:FT:295:ASN:HB2	2.16	0.46
3:GY:1873:GLU:HG2	3:HY:1781:LEU:HD22	1.96	0.46
1:HT:199:TYR:HB3	2:HX:379:GLN:OE1	2.16	0.46
1:KT:60:PRO:HA	1:KT:137:ASN:O	2.15	0.46
2:KX:477:LYS:HD3	2:KX:484:LEU:HD11	1.96	0.46
1:BT:39:VAL:HG21	2:CX:419:ILE:HD12	1.97	0.46
1:ET:291:GLU:O	1:ET:295:ASN:HB2	2.16	0.46
2:MX:477:LYS:HD3	2:MX:484:LEU:HD11	1.96	0.46
2:NX:477:LYS:HD3	2:NX:484:LEU:HD11	1.96	0.46
1:HT:291:GLU:O	1:HT:295:ASN:HB2	2.16	0.46
1:KT:51:PRO:HD3	2:KX:495:ASN:HB2	1.97	0.46
1:KT:200:LYS:HD3	1:KT:200:LYS:HA	1.76	0.46
2:LX:477:LYS:HD3	2:LX:484:LEU:HD11	1.96	0.46
2:BX:455:ILE:HB	3:BY:1765:PRO:HB3	1.98	0.46
3:FY:1707:ASP:HB2	3:GY:1783:GLU:HB2	1.97	0.46
3:KY:1707:ASP:HB2	3:LY:1783:GLU:HB2	1.98	0.46
3:KY:1873:GLU:HG2	3:LY:1781:LEU:HD22	1.98	0.46
1:MT:291:GLU:O	1:MT:295:ASN:HB2	2.16	0.46
2:MX:392:LEU:HD22	3:NY:1728:THR:HB	1.98	0.46
2:AX:392:LEU:HD22	3:BY:1728:THR:HB	1.98	0.46
1:BT:60:PRO:HA	1:BT:137:ASN:O	2.15	0.46
2:DX:477:LYS:HD3	2:DX:484:LEU:HD11	1.96	0.46
3:DY:1873:GLU:HG2	3:EY:1781:LEU:HD22	1.98	0.46
1:MT:60:PRO:HA	1:MT:137:ASN:O	2.15	0.46
2:AX:477:LYS:HD3	2:AX:484:LEU:HD11	1.96	0.46
2:CX:436:GLN:HB2	3:CY:1734:LYS:HG2	1.98	0.46
2:CX:477:LYS:HD3	2:CX:484:LEU:HD11	1.96	0.46
1:DT:291:GLU:O	1:DT:295:ASN:HB2	2.16	0.46
3:IY:1707:ASP:HB2	3:JY:1783:GLU:HB2	1.97	0.46
3:JY:1713:ILE:HD11	3:KY:1880:LEU:HD23	1.97	0.46
2:BX:477:LYS:HD3	2:BX:484:LEU:HD11	1.96	0.46
3:KY:1713:ILE:HD11	3:LY:1880:LEU:HD23	1.97	0.46
1:JT:291:GLU:O	1:JT:295:ASN:HB2	2.16	0.45
1:NT:60:PRO:HA	1:NT:137:ASN:O	2.15	0.45
3:AY:1880:LEU:HD23	3:NY:1713:ILE:HD11	1.98	0.45
1:CT:291:GLU:O	1:CT:295:ASN:HB2	2.16	0.45
3:MY:1713:ILE:HD11	3:NY:1880:LEU:HD23	1.97	0.45
3:AY:1707:ASP:HB2	3:BY:1783:GLU:HB2	1.98	0.45
3:DY:1707:ASP:HB2	3:EY:1783:GLU:HB2	1.98	0.45
1:NT:200:LYS:HA	1:NT:200:LYS:HD3	1.76	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:BY:1707:ASP:HB2	3:CY:1783:GLU:HB2	1.97	0.45
3:EY:1707:ASP:HB2	3:FY:1783:GLU:HB2	1.96	0.45
3:EY:1713:ILE:HD11	3:FY:1880:LEU:HD23	1.98	0.45
1:NT:291:GLU:O	1:NT:295:ASN:HB2	2.16	0.45
1:BT:291:GLU:O	1:BT:295:ASN:HB2	2.16	0.45
2:HX:392:LEU:HD22	3:IY:1728:THR:HB	1.97	0.45
1:IT:291:GLU:O	1:IT:295:ASN:HB2	2.16	0.45
1:LT:200:LYS:HD3	1:LT:200:LYS:HA	1.76	0.45
2:AX:432:ASN:ND2	2:NX:497:LEU:O	2.38	0.45
3:AY:1873:GLU:HG2	3:BY:1781:LEU:HD22	1.98	0.45
1:CT:38:PRO:HG2	1:DT:55:LYS:HB2	1.97	0.45
3:FY:1690:PHE:O	3:FY:1887:PHE:N	2.45	0.45
3:HY:1873:GLU:HG2	3:IY:1781:LEU:HD22	1.97	0.45
2:JX:481:ASP:HB3	3:JY:1875:ASP:HA	1.98	0.45
1:CT:51:PRO:HD3	2:CX:495:ASN:HB2	1.99	0.45
3:EY:1900:VAL:HA	3:EY:1903:GLU:HG3	1.99	0.45
1:AT:291:GLU:O	1:AT:295:ASN:HB2	2.16	0.45
3:AY:1900:VAL:HA	3:AY:1903:GLU:HG3	1.99	0.45
1:GT:111:ALA:O	1:HT:167:ASN:ND2	2.50	0.45
1:JT:36:TYR:HB3	2:KX:404:TYR:HB3	1.98	0.45
3:FY:1900:VAL:HA	3:FY:1903:GLU:HG3	1.99	0.45
2:LX:392:LEU:HD22	3:MY:1728:THR:HB	1.99	0.45
3:NY:1690:PHE:O	3:NY:1887:PHE:N	2.45	0.45
3:NY:1900:VAL:HA	3:NY:1903:GLU:HG3	1.99	0.45
1:AT:200:LYS:HD3	1:AT:200:LYS:HA	1.76	0.45
2:BX:392:LEU:HD22	3:CY:1728:THR:HB	1.99	0.44
2:KX:417:SER:N	2:KX:429:GLY:O	2.44	0.44
2:LX:497:LEU:O	2:MX:432:ASN:ND2	2.40	0.44
3:LY:1713:ILE:HD11	3:MY:1880:LEU:HD23	1.98	0.44
3:AY:1690:PHE:O	3:AY:1887:PHE:N	2.45	0.44
3:DY:1900:VAL:HA	3:DY:1903:GLU:HG3	1.99	0.44
3:HY:1707:ASP:HB2	3:IY:1783:GLU:HB2	1.98	0.44
3:MY:1900:VAL:HA	3:MY:1903:GLU:HG3	1.99	0.44
1:AT:75:LEU:HD23	1:AT:75:LEU:HA	1.86	0.44
3:AY:1728:THR:HB	2:NX:392:LEU:HD22	1.99	0.44
2:EX:497:LEU:O	2:FX:432:ASN:ND2	2.41	0.44
3:KY:1732:LEU:HA	3:KY:1732:LEU:HD23	1.84	0.44
2:NX:480:LYS:NZ	3:NY:1699:GLU:OE2	2.48	0.44
2:AX:507:GLY:HA2	1:BT:168:LEU:HD11	1.99	0.44
1:FT:33:ASN:ND2	2:GX:407:PRO:O	2.50	0.44
2:FX:392:LEU:HD22	3:GY:1728:THR:HB	1.98	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:LT:196:ASN:OD1	3:MY:1907:GLN:NE2	2.50	0.44
1:MT:55:LYS:O	1:MT:59:THR:OG1	2.31	0.44
1:MT:75:LEU:HA	1:MT:75:LEU:HD23	1.86	0.44
3:LY:1900:VAL:HA	3:LY:1903:GLU:HG3	1.99	0.44
3:BY:1900:VAL:HA	3:BY:1903:GLU:HG3	1.99	0.44
2:GX:417:SER:N	2:GX:429:GLY:O	2.45	0.44
1:HT:213:ARG:HH12	2:HX:364:ASN:HB3	1.83	0.44
2:JX:436:GLN:HB2	3:JY:1734:LYS:HG2	1.99	0.44
2:CX:455:ILE:HB	3:CY:1765:PRO:HB3	1.99	0.44
3:GY:1900:VAL:HA	3:GY:1903:GLU:HG3	1.99	0.44
1:AT:33:ASN:HD22	2:BX:406:ALA:HB1	1.83	0.44
2:EX:481:ASP:OD1	2:EX:481:ASP:N	2.50	0.44
1:MT:200:LYS:HD3	1:MT:200:LYS:HA	1.76	0.44
3:MY:1690:PHE:O	3:MY:1887:PHE:N	2.45	0.44
3:AY:1729:MET:HB3	2:NX:503:ILE:HG13	2.00	0.44
1:FT:75:LEU:HD23	1:FT:75:LEU:HA	1.86	0.44
1:JT:38:PRO:HG2	1:KT:55:LYS:HB2	2.00	0.44
3:JY:1900:VAL:HA	3:JY:1903:GLU:HG3	1.99	0.44
3:KY:1900:VAL:HA	3:KY:1903:GLU:HG3	1.99	0.44
3:IY:1900:VAL:HA	3:IY:1903:GLU:HG3	1.99	0.43
1:NT:75:LEU:HD23	1:NT:75:LEU:HA	1.86	0.43
3:CY:1900:VAL:HA	3:CY:1903:GLU:HG3	1.99	0.43
2:DX:503:ILE:HG13	3:EY:1729:MET:HB3	2.01	0.43
1:FT:200:LYS:HA	1:FT:200:LYS:HD3	1.76	0.43
2:GX:392:LEU:HD22	3:HY:1728:THR:HB	1.99	0.43
2:DX:455:ILE:HB	3:DY:1765:PRO:HB3	2.00	0.43
1:ET:200:LYS:HA	1:ET:200:LYS:HD3	1.76	0.43
3:EY:1690:PHE:O	3:EY:1887:PHE:N	2.45	0.43
2:FX:507:GLY:HA2	1:GT:168:LEU:HD11	2.00	0.43
3:HY:1900:VAL:HA	3:HY:1903:GLU:HG3	1.99	0.43
1:JT:111:ALA:O	1:KT:167:ASN:ND2	2.52	0.43
3:JY:1732:LEU:HA	3:JY:1732:LEU:HD23	1.84	0.43
1:LT:38:PRO:HA	2:MX:404:TYR:HA	2.00	0.43
3:DY:1690:PHE:O	3:DY:1887:PHE:N	2.45	0.43
3:GY:1708:ALA:HB3	3:GY:1869:PHE:HB3	2.01	0.43
1:JT:199:TYR:HB3	2:JX:379:GLN:OE1	2.19	0.43
3:MY:1708:ALA:HB3	3:MY:1869:PHE:HB3	2.01	0.43
2:DX:503:ILE:HA	3:EY:1729:MET:HA	2.00	0.43
1:FT:54:ASP:OD2	2:FX:404:TYR:OH	2.26	0.43
1:FT:196:ASN:OD1	3:GY:1907:GLN:NE2	2.52	0.43
3:FY:1732:LEU:HD23	3:FY:1732:LEU:HA	1.84	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:GT:200:LYS:HA	1:GT:200:LYS:HD3	1.76	0.43
2:KX:481:ASP:HB3	3:KY:1875:ASP:HA	2.01	0.43
3:MY:1906:LYS:HA	3:MY:1909:THR:HG22	2.01	0.43
3:AY:1729:MET:HA	2:NX:503:ILE:HA	2.00	0.43
3:AY:1906:LYS:HA	3:AY:1909:THR:HG22	2.01	0.43
3:BY:1708:ALA:HB3	3:BY:1869:PHE:HB3	2.01	0.43
3:DY:1708:ALA:HB3	3:DY:1869:PHE:HB3	2.01	0.43
3:KY:1708:ALA:HB3	3:KY:1869:PHE:HB3	2.01	0.43
3:KY:1906:LYS:HA	3:KY:1909:THR:HG22	2.01	0.43
2:AX:394:ARG:HE	2:AX:394:ARG:HB3	1.67	0.43
3:BY:1690:PHE:O	3:BY:1887:PHE:N	2.45	0.43
3:CY:1690:PHE:O	3:CY:1887:PHE:N	2.45	0.43
3:CY:1906:LYS:HA	3:CY:1909:THR:HG22	2.01	0.43
2:DX:497:LEU:O	2:EX:432:ASN:ND2	2.39	0.43
2:FX:481:ASP:OD1	2:FX:481:ASP:N	2.50	0.43
1:GT:196:ASN:OD1	3:HY:1907:GLN:NE2	2.51	0.43
2:GX:497:LEU:O	2:HX:432:ASN:ND2	2.39	0.43
2:GX:507:GLY:HA2	1:HT:168:LEU:HD11	2.01	0.43
1:HT:200:LYS:HA	1:HT:200:LYS:HD3	1.76	0.43
3:IY:1708:ALA:HB3	3:IY:1869:PHE:HB3	2.01	0.43
3:AY:1708:ALA:HB3	3:AY:1869:PHE:HB3	2.01	0.43
1:BT:306:LEU:HD22	2:BX:384:GLU:HB2	2.00	0.43
3:EY:1708:ALA:HB3	3:EY:1869:PHE:HB3	2.01	0.43
1:HT:48:SER:OG	1:IT:170:GLY:O	2.34	0.43
2:HX:507:GLY:HA2	1:IT:168:LEU:HD11	2.01	0.43
2:LX:417:SER:N	2:LX:429:GLY:O	2.44	0.43
1:BT:200:LYS:HA	1:BT:200:LYS:HD3	1.76	0.43
1:DT:200:LYS:HA	1:DT:200:LYS:HD3	1.76	0.43
2:DX:392:LEU:HD22	3:EY:1728:THR:HB	1.99	0.43
1:ET:51:PRO:HD3	2:EX:495:ASN:HB2	2.01	0.43
3:GY:1732:LEU:HA	3:GY:1732:LEU:HD23	1.84	0.43
2:IX:480:LYS:NZ	3:IY:1699:GLU:OE2	2.41	0.43
3:NY:1906:LYS:HA	3:NY:1909:THR:HG22	2.01	0.43
3:BY:1906:LYS:HA	3:BY:1909:THR:HG22	2.01	0.42
1:CT:39:VAL:HG11	2:DX:419:ILE:HG13	2.00	0.42
3:EY:1906:LYS:HA	3:EY:1909:THR:HG22	2.01	0.42
3:HY:1906:LYS:HA	3:HY:1909:THR:HG22	2.01	0.42
3:IY:1732:LEU:HA	3:IY:1732:LEU:HD23	1.84	0.42
2:JX:455:ILE:HB	3:JY:1765:PRO:HB3	2.00	0.42
3:JY:1906:LYS:HA	3:JY:1909:THR:HG22	2.01	0.42
3:LY:1906:LYS:HA	3:LY:1909:THR:HG22	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:MX:503:ILE:HA	3:NY:1729:MET:HA	2.01	0.42
3:FY:1720:LYS:HA	3:FY:1734:LYS:HE2	2.01	0.42
3:HY:1732:LEU:HA	3:HY:1732:LEU:HD23	1.84	0.42
3:NY:1708:ALA:HB3	3:NY:1869:PHE:HB3	2.01	0.42
2:AX:436:GLN:HB2	3:AY:1734:LYS:HG2	2.01	0.42
3:EY:1720:LYS:HA	3:EY:1734:LYS:HE2	2.01	0.42
2:FX:503:ILE:HA	3:GY:1729:MET:HA	2.01	0.42
3:FY:1708:ALA:HB3	3:FY:1869:PHE:HB3	2.01	0.42
3:GY:1720:LYS:HA	3:GY:1734:LYS:HE2	2.01	0.42
2:LX:507:GLY:HA2	1:MT:168:LEU:HD11	2.00	0.42
3:DY:1720:LYS:HA	3:DY:1734:LYS:HE2	2.00	0.42
2:EX:392:LEU:HD22	3:FY:1728:THR:HB	1.99	0.42
3:GY:1903:GLU:HA	3:GY:1906:LYS:HG3	2.02	0.42
3:HY:1720:LYS:HA	3:HY:1734:LYS:HE2	2.00	0.42
2:JX:392:LEU:HD22	3:KY:1728:THR:HB	2.01	0.42
1:KT:81:VAL:HG21	2:KX:459:MET:HE3	2.01	0.42
2:LX:394:ARG:HE	2:LX:394:ARG:HB3	1.68	0.42
2:MX:497:LEU:O	2:NX:432:ASN:ND2	2.40	0.42
2:NX:440:PHE:CZ	3:NY:1873:GLU:HB3	2.55	0.42
3:NY:1786:VAL:HG22	3:NY:1871:LYS:HE2	2.01	0.42
3:IY:1903:GLU:HA	3:IY:1906:LYS:HG3	2.02	0.42
3:KY:1786:VAL:HG22	3:KY:1871:LYS:HE2	2.01	0.42
3:KY:1903:GLU:HA	3:KY:1906:LYS:HG3	2.02	0.42
3:MY:1786:VAL:HG22	3:MY:1871:LYS:HE2	2.01	0.42
1:NT:54:ASP:OD2	2:NX:404:TYR:OH	2.32	0.42
2:BX:503:ILE:HG13	3:CY:1729:MET:HB3	2.02	0.42
2:DX:436:GLN:HB2	3:DY:1734:LYS:HG2	2.01	0.42
2:DX:507:GLY:HA2	1:ET:168:LEU:HD11	2.02	0.42
2:EX:436:GLN:HB2	3:EY:1734:LYS:HG2	2.02	0.42
3:EY:1903:GLU:HA	3:EY:1906:LYS:HG3	2.02	0.42
2:FX:503:ILE:HG13	3:GY:1729:MET:HB3	2.02	0.42
3:JY:1689:THR:O	3:JY:1724:ASN:ND2	2.34	0.42
3:MY:1720:LYS:HA	3:MY:1734:LYS:HE2	2.00	0.42
3:NY:1720:LYS:HA	3:NY:1734:LYS:HE2	2.00	0.42
2:CX:440:PHE:CZ	3:CY:1873:GLU:HB3	2.54	0.42
2:EX:503:ILE:HG13	3:FY:1729:MET:HB3	2.02	0.42
2:GX:440:PHE:CZ	3:GY:1873:GLU:HB3	2.55	0.42
3:IY:1720:LYS:HA	3:IY:1734:LYS:HE2	2.00	0.42
1:JT:48:SER:OG	1:KT:170:GLY:O	2.28	0.42
3:LY:1786:VAL:HG22	3:LY:1871:LYS:HE2	2.01	0.42
2:AX:455:ILE:HB	3:AY:1765:PRO:HB3	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:IT:200:LYS:HA	1:IT:200:LYS:HD3	1.76	0.42
2:IX:479:ILE:HG22	3:IY:1874:GLY:HA3	2.02	0.42
3:AY:1720:LYS:HA	3:AY:1734:LYS:HE2	2.01	0.42
3:CY:1720:LYS:HA	3:CY:1734:LYS:HE2	2.01	0.42
2:DX:481:ASP:HB3	3:DY:1875:ASP:HA	2.01	0.42
3:FY:1906:LYS:HA	3:FY:1909:THR:HG22	2.01	0.42
2:GX:436:GLN:HB2	3:GY:1734:LYS:HG2	2.02	0.42
1:JT:56:LEU:HD12	1:JT:59:THR:HB	2.02	0.42
3:KY:1689:THR:O	3:KY:1724:ASN:ND2	2.34	0.42
1:LT:199:TYR:HB3	2:LX:379:GLN:OE1	2.20	0.42
3:LY:1708:ALA:HB3	3:LY:1869:PHE:HB3	2.01	0.42
2:MX:503:ILE:HG13	3:NY:1729:MET:HB3	2.02	0.42
2:AX:503:ILE:HA	3:BY:1729:MET:HA	2.01	0.42
3:BY:1903:GLU:HA	3:BY:1906:LYS:HG3	2.01	0.42
3:DY:1903:GLU:HA	3:DY:1906:LYS:HG3	2.02	0.42
3:HY:1903:GLU:HA	3:HY:1906:LYS:HG3	2.02	0.42
1:KT:75:LEU:HD23	1:KT:75:LEU:HA	1.86	0.42
2:LX:503:ILE:HG13	3:MY:1729:MET:HB3	2.02	0.42
3:LY:1720:LYS:HA	3:LY:1734:LYS:HE2	2.01	0.42
3:MY:1903:GLU:HA	3:MY:1906:LYS:HG3	2.02	0.42
3:AY:1786:VAL:HG22	3:AY:1871:LYS:HE2	2.01	0.41
3:AY:1859:LEU:O	3:AY:1863:MET:HB3	2.20	0.41
3:BY:1786:VAL:HG22	3:BY:1871:LYS:HE2	2.01	0.41
1:CT:200:LYS:HA	1:CT:200:LYS:HD3	1.76	0.41
2:CX:392:LEU:HD22	3:DY:1728:THR:HB	2.01	0.41
2:CX:479:ILE:HG22	3:CY:1874:GLY:HA3	2.02	0.41
2:GX:481:ASP:OD1	2:GX:481:ASP:N	2.50	0.41
2:GX:503:ILE:HG13	3:HY:1729:MET:HB3	2.02	0.41
3:HY:1708:ALA:HB3	3:HY:1869:PHE:HB3	2.01	0.41
3:HY:1859:LEU:O	3:HY:1863:MET:HB3	2.20	0.41
3:IY:1859:LEU:O	3:IY:1863:MET:HB3	2.20	0.41
3:JY:1786:VAL:HG22	3:JY:1871:LYS:HE2	2.01	0.41
3:JY:1903:GLU:HA	3:JY:1906:LYS:HG3	2.02	0.41
2:KX:392:LEU:HD22	3:LY:1728:THR:HB	2.02	0.41
2:LX:440:PHE:CZ	3:LY:1873:GLU:HB3	2.55	0.41
2:LX:480:LYS:NZ	3:LY:1699:GLU:OE2	2.48	0.41
3:MY:1898:LYS:HA	3:MY:1901:VAL:HG12	2.02	0.41
3:NY:1859:LEU:O	3:NY:1863:MET:HB3	2.20	0.41
3:AY:1732:LEU:HA	3:AY:1732:LEU:HD23	1.84	0.41
2:BX:394:ARG:HE	2:BX:394:ARG:HB3	1.68	0.41
3:CY:1708:ALA:HB3	3:CY:1869:PHE:HB3	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:FX:371:GLU:HA	2:FX:374:LYS:HG2	2.02	0.41
3:FY:1903:GLU:HA	3:FY:1906:LYS:HG3	2.02	0.41
3:IY:1890:VAL:HG12	3:JY:1683:ILE:HG13	2.02	0.41
3:IY:1906:LYS:HA	3:IY:1909:THR:HG22	2.01	0.41
3:JY:1720:LYS:HA	3:JY:1734:LYS:HE2	2.01	0.41
1:KT:56:LEU:HD12	1:KT:59:THR:HB	2.02	0.41
1:BT:51:PRO:HD3	2:BX:495:ASN:HB2	2.02	0.41
3:BY:1720:LYS:HA	3:BY:1734:LYS:HE2	2.00	0.41
2:CX:435:LEU:HD11	3:CY:1720:LYS:HE3	2.02	0.41
3:EY:1859:LEU:O	3:EY:1863:MET:HB3	2.20	0.41
2:FX:501:TYR:OH	1:GT:168:LEU:HB2	2.20	0.41
1:GT:56:LEU:HD12	1:GT:59:THR:HB	2.02	0.41
1:HT:56:LEU:HD12	1:HT:59:THR:HB	2.02	0.41
2:HX:371:GLU:HA	2:HX:374:LYS:HG2	2.03	0.41
1:IT:56:LEU:HD12	1:IT:59:THR:HB	2.02	0.41
3:IY:1689:THR:O	3:IY:1724:ASN:ND2	2.34	0.41
1:KT:48:SER:OG	1:LT:170:GLY:O	2.34	0.41
2:MX:417:SER:N	2:MX:429:GLY:O	2.44	0.41
3:NY:1797:ARG:HG3	3:NY:1865:ILE:HB	2.03	0.41
1:BT:167:ASN:HB3	1:BT:173:VAL:HG11	2.03	0.41
3:BY:1859:LEU:O	3:BY:1863:MET:HB3	2.20	0.41
3:BY:1898:LYS:HA	3:BY:1901:VAL:HG12	2.02	0.41
2:EX:503:ILE:HA	3:FY:1729:MET:HA	2.02	0.41
1:FT:213:ARG:HH12	2:FX:364:ASN:HB3	1.86	0.41
3:FY:1859:LEU:O	3:FY:1863:MET:HB3	2.20	0.41
3:IY:1786:VAL:HG22	3:IY:1871:LYS:HE2	2.01	0.41
3:JY:1708:ALA:HB3	3:JY:1869:PHE:HB3	2.01	0.41
3:JY:1859:LEU:O	3:JY:1863:MET:HB3	2.20	0.41
3:KY:1797:ARG:HG3	3:KY:1865:ILE:HB	2.03	0.41
1:LT:48:SER:OG	1:MT:170:GLY:O	2.34	0.41
1:MT:51:PRO:HD3	2:MX:495:ASN:HB2	2.03	0.41
2:MX:507:GLY:HA2	1:NT:168:LEU:HD11	2.02	0.41
1:CT:167:ASN:HB3	1:CT:173:VAL:HG11	2.03	0.41
2:CX:450:MET:HE3	3:DY:1880:LEU:HD11	2.01	0.41
1:DT:213:ARG:HH12	2:DX:364:ASN:HB3	1.85	0.41
1:FT:56:LEU:HD12	1:FT:59:THR:HB	2.02	0.41
2:GX:394:ARG:HE	2:GX:394:ARG:HB3	1.68	0.41
3:GY:1906:LYS:HA	3:GY:1909:THR:HG22	2.01	0.41
1:JT:298:LEU:HD13	2:JX:373:GLN:HE22	1.86	0.41
3:JY:1797:ARG:HG3	3:JY:1865:ILE:HB	2.03	0.41
3:KY:1720:LYS:HA	3:KY:1734:LYS:HE2	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:MY:1797:ARG:HG3	3:MY:1865:ILE:HB	2.03	0.41
1:AT:167:ASN:HB3	1:AT:173:VAL:HG11	2.03	0.41
3:CY:1797:ARG:HG3	3:CY:1865:ILE:HB	2.03	0.41
2:HX:417:SER:N	2:HX:429:GLY:O	2.45	0.41
3:KY:1898:LYS:HA	3:KY:1901:VAL:HG12	2.03	0.41
1:LT:56:LEU:HD12	1:LT:59:THR:HB	2.02	0.41
2:LX:503:ILE:HA	3:MY:1729:MET:HA	2.02	0.41
3:LY:1890:VAL:HG12	3:MY:1683:ILE:HG13	2.03	0.41
3:NY:1700:ILE:HG13	3:NY:1879:ILE:HG12	2.02	0.41
3:NY:1898:LYS:HA	3:NY:1901:VAL:HG12	2.02	0.41
3:AY:1797:ARG:HG3	3:AY:1865:ILE:HB	2.03	0.41
2:BX:503:ILE:HA	3:CY:1729:MET:HA	2.02	0.41
3:BY:1700:ILE:HG13	3:BY:1879:ILE:HG12	2.02	0.41
3:BY:1797:ARG:HG3	3:BY:1865:ILE:HB	2.03	0.41
3:CY:1732:LEU:HD23	3:CY:1732:LEU:HA	1.84	0.41
1:DT:167:ASN:HB3	1:DT:173:VAL:HG11	2.03	0.41
2:DX:371:GLU:HA	2:DX:374:LYS:HG2	2.02	0.41
2:EX:440:PHE:CZ	3:EY:1873:GLU:HB3	2.56	0.41
2:FX:440:PHE:CZ	3:FY:1873:GLU:HB3	2.55	0.41
2:GX:503:ILE:HA	3:HY:1729:MET:HA	2.02	0.41
3:GY:1786:VAL:HG22	3:GY:1871:LYS:HE2	2.01	0.41
3:HY:1797:ARG:HG3	3:HY:1865:ILE:HB	2.03	0.41
1:IT:51:PRO:HD3	2:IX:495:ASN:HB2	2.03	0.41
2:JX:371:GLU:HA	2:JX:374:LYS:HG2	2.02	0.41
3:JY:1700:ILE:HG13	3:JY:1879:ILE:HG12	2.02	0.41
3:LY:1797:ARG:HG3	3:LY:1865:ILE:HB	2.03	0.41
2:MX:440:PHE:CZ	3:MY:1873:GLU:HB3	2.55	0.41
3:MY:1859:LEU:O	3:MY:1863:MET:HB3	2.20	0.41
3:AY:1700:ILE:HG13	3:AY:1879:ILE:HG12	2.02	0.41
3:AY:1898:LYS:HA	3:AY:1901:VAL:HG12	2.02	0.41
2:BX:435:LEU:HD11	3:BY:1720:LYS:HE3	2.02	0.41
1:CT:75:LEU:HD23	1:CT:75:LEU:HA	1.86	0.41
1:CT:213:ARG:HH12	2:CX:364:ASN:HB3	1.85	0.41
3:CY:1786:VAL:HG22	3:CY:1871:LYS:HE2	2.01	0.41
3:DY:1859:LEU:O	3:DY:1863:MET:HB3	2.20	0.41
3:DY:1906:LYS:HA	3:DY:1909:THR:HG22	2.01	0.41
2:EX:455:ILE:HB	3:EY:1765:PRO:HB3	2.03	0.41
3:EY:1890:VAL:HG12	3:FY:1683:ILE:HG13	2.02	0.41
2:FX:436:GLN:HB2	3:FY:1734:LYS:HG2	2.02	0.41
3:FY:1786:VAL:HG22	3:FY:1871:LYS:HE2	2.01	0.41
3:IY:1797:ARG:HG3	3:IY:1865:ILE:HB	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:KY:1700:ILE:HG13	3:KY:1879:ILE:HG12	2.02	0.41
3:LY:1700:ILE:HG13	3:LY:1879:ILE:HG12	2.02	0.41
1:NT:167:ASN:HB3	1:NT:173:VAL:HG11	2.03	0.41
2:NX:394:ARG:HE	2:NX:394:ARG:HB3	1.68	0.41
1:BT:75:LEU:HA	1:BT:75:LEU:HD23	1.86	0.41
2:BX:440:PHE:CZ	3:BY:1873:GLU:HB3	2.56	0.41
3:CY:1694:GLN:N	3:CY:1884:ASP:OD2	2.54	0.41
3:CY:1700:ILE:HG13	3:CY:1879:ILE:HG12	2.02	0.41
3:CY:1903:GLU:HA	3:CY:1906:LYS:HG3	2.02	0.41
1:DT:298:LEU:HD12	1:DT:298:LEU:HA	1.88	0.41
3:DY:1694:GLN:N	3:DY:1884:ASP:OD2	2.54	0.41
1:ET:167:ASN:HB3	1:ET:173:VAL:HG11	2.03	0.41
3:EY:1786:VAL:HG22	3:EY:1871:LYS:HE2	2.01	0.41
1:FT:38:PRO:HA	2:GX:404:TYR:HA	2.03	0.41
3:FY:1689:THR:O	3:FY:1724:ASN:ND2	2.34	0.41
1:GT:199:TYR:HB3	2:GX:379:GLN:OE1	2.21	0.41
2:GX:455:ILE:HB	3:GY:1765:PRO:HB3	2.02	0.41
2:GX:481:ASP:HB3	3:GY:1875:ASP:HA	2.02	0.41
3:GY:1694:GLN:N	3:GY:1884:ASP:OD2	2.54	0.41
1:HT:298:LEU:HA	1:HT:298:LEU:HD12	1.88	0.41
2:HX:436:GLN:HB2	3:HY:1734:LYS:HG2	2.02	0.41
2:HX:440:PHE:CZ	3:HY:1873:GLU:HB3	2.56	0.41
2:HX:503:ILE:HA	3:IY:1729:MET:HA	2.02	0.41
3:HY:1786:VAL:HG22	3:HY:1871:LYS:HE2	2.01	0.41
3:IY:1700:ILE:HG13	3:IY:1879:ILE:HG12	2.02	0.41
1:JT:200:LYS:HA	1:JT:200:LYS:HD3	1.76	0.41
2:JX:440:PHE:CZ	3:JY:1873:GLU:HB3	2.56	0.41
2:JX:497:LEU:O	2:KX:432:ASN:ND2	2.40	0.41
3:JY:1898:LYS:HA	3:JY:1901:VAL:HG12	2.02	0.41
2:KX:436:GLN:HB2	3:KY:1734:LYS:HG2	2.02	0.41
2:KX:455:ILE:HB	3:KY:1765:PRO:HB3	2.01	0.41
3:KY:1859:LEU:O	3:KY:1863:MET:HB3	2.20	0.41
1:MT:56:LEU:HD12	1:MT:59:THR:HB	2.02	0.41
1:MT:167:ASN:HB3	1:MT:173:VAL:HG11	2.03	0.41
1:MT:199:TYR:HB3	2:MX:379:GLN:OE1	2.21	0.41
3:MY:1700:ILE:HG13	3:MY:1879:ILE:HG12	2.02	0.41
2:NX:436:GLN:HB2	3:NY:1734:LYS:HG2	2.02	0.41
3:NY:1903:GLU:HA	3:NY:1906:LYS:HG3	2.02	0.41
2:AX:440:PHE:CZ	3:AY:1873:GLU:HB3	2.56	0.41
1:BT:56:LEU:HD12	1:BT:59:THR:HB	2.02	0.41
3:BY:1890:VAL:HG12	3:CY:1683:ILE:HG13	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:CX:480:LYS:NZ	3:CY:1699:GLU:OE2	2.45	0.41
2:CX:507:GLY:HA2	1:DT:168:LEU:HD11	2.03	0.41
3:DY:1797:ARG:HG3	3:DY:1865:ILE:HB	2.03	0.41
1:ET:56:LEU:HD12	1:ET:59:THR:HB	2.02	0.41
3:FY:1694:GLN:N	3:FY:1884:ASP:OD2	2.54	0.41
3:FY:1797:ARG:HG3	3:FY:1865:ILE:HB	2.03	0.41
2:HX:365:LYS:HE2	2:HX:365:LYS:HB2	1.98	0.41
1:LT:75:LEU:HD23	1:LT:75:LEU:HA	1.86	0.41
3:LY:1689:THR:O	3:LY:1724:ASN:ND2	2.34	0.41
2:AX:503:ILE:HG13	3:BY:1729:MET:HB3	2.03	0.40
1:CT:56:LEU:HD12	1:CT:59:THR:HB	2.02	0.40
1:DT:75:LEU:HD23	1:DT:75:LEU:HA	1.86	0.40
3:DY:1898:LYS:HA	3:DY:1901:VAL:HG12	2.03	0.40
2:EX:365:LYS:HE2	2:EX:365:LYS:HB2	1.98	0.40
2:HX:455:ILE:HB	3:HY:1765:PRO:HB3	2.02	0.40
1:LT:167:ASN:HB3	1:LT:173:VAL:HG11	2.03	0.40
2:MX:436:GLN:HB2	3:MY:1734:LYS:HG2	2.02	0.40
2:NX:394:ARG:HD2	2:NX:394:ARG:HA	1.93	0.40
2:NX:417:SER:N	2:NX:429:GLY:O	2.45	0.40
1:AT:168:LEU:HD11	2:NX:507:GLY:HA2	2.03	0.40
3:CY:1859:LEU:O	3:CY:1863:MET:HB3	2.20	0.40
3:EY:1797:ARG:HG3	3:EY:1865:ILE:HB	2.03	0.40
3:GY:1859:LEU:O	3:GY:1863:MET:HB3	2.20	0.40
3:HY:1700:ILE:HG13	3:HY:1879:ILE:HG12	2.02	0.40
1:KT:167:ASN:HB3	1:KT:173:VAL:HG11	2.03	0.40
1:AT:56:LEU:HD12	1:AT:59:THR:HB	2.02	0.40
2:AX:417:SER:N	2:AX:429:GLY:O	2.45	0.40
1:DT:56:LEU:HD12	1:DT:59:THR:HB	2.02	0.40
3:DY:1786:VAL:HG22	3:DY:1871:LYS:HE2	2.01	0.40
3:FY:1879:ILE:HD13	3:FY:1879:ILE:HA	1.86	0.40
2:GX:365:LYS:HE2	2:GX:365:LYS:HB2	1.98	0.40
3:GY:1797:ARG:HG3	3:GY:1865:ILE:HB	2.03	0.40
3:IY:1694:GLN:N	3:IY:1884:ASP:OD2	2.54	0.40
3:KY:1890:VAL:HG12	3:LY:1683:ILE:HG13	2.03	0.40
2:LX:436:GLN:HB2	3:LY:1734:LYS:HG2	2.02	0.40
2:AX:481:ASP:HB3	3:AY:1875:ASP:HA	2.02	0.40
3:AY:1694:GLN:N	3:AY:1884:ASP:OD2	2.54	0.40
2:DX:440:PHE:CZ	3:DY:1873:GLU:HB3	2.57	0.40
2:EX:394:ARG:HE	2:EX:394:ARG:HB3	1.68	0.40
3:EY:1879:ILE:HD13	3:EY:1879:ILE:HA	1.86	0.40
1:FT:167:ASN:HB3	1:FT:173:VAL:HG11	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:FT:202:LEU:HD13	2:FX:375:ILE:HD11	2.04	0.40
2:GX:371:GLU:HA	2:GX:374:LYS:HG2	2.03	0.40
2:HX:481:ASP:HB3	3:HY:1875:ASP:HA	2.02	0.40
3:LY:1898:LYS:HA	3:LY:1901:VAL:HG12	2.02	0.40
3:AY:1903:GLU:HA	3:AY:1906:LYS:HG3	2.02	0.40
3:DY:1700:ILE:HG13	3:DY:1879:ILE:HG12	2.02	0.40
3:FY:1898:LYS:HA	3:FY:1901:VAL:HG12	2.02	0.40
3:GY:1879:ILE:HD13	3:GY:1879:ILE:HA	1.86	0.40
1:HT:51:PRO:HD3	2:HX:495:ASN:HB2	2.04	0.40
3:HY:1898:LYS:HA	3:HY:1901:VAL:HG12	2.03	0.40
3:IY:1898:LYS:HA	3:IY:1901:VAL:HG12	2.03	0.40
2:LX:394:ARG:HD2	2:LX:394:ARG:HA	1.93	0.40
2:NX:371:GLU:HA	2:NX:374:LYS:HG2	2.02	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	AT	171/278 (62%)	160 (94%)	11 (6%)	0	100	100
1	BT	171/278 (62%)	160 (94%)	11 (6%)	0	100	100
1	CT	171/278 (62%)	160 (94%)	11 (6%)	0	100	100
1	DT	171/278 (62%)	160 (94%)	11 (6%)	0	100	100
1	ET	171/278 (62%)	160 (94%)	11 (6%)	0	100	100
1	FT	171/278 (62%)	160 (94%)	11 (6%)	0	100	100
1	GT	171/278 (62%)	160 (94%)	11 (6%)	0	100	100
1	HT	171/278 (62%)	160 (94%)	11 (6%)	0	100	100
1	IT	171/278 (62%)	160 (94%)	11 (6%)	0	100	100
1	JT	171/278 (62%)	160 (94%)	11 (6%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	KT	171/278 (62%)	160 (94%)	11 (6%)	0	100	100
1	LT	171/278 (62%)	160 (94%)	11 (6%)	0	100	100
1	MT	171/278 (62%)	160 (94%)	11 (6%)	0	100	100
1	NT	171/278 (62%)	160 (94%)	11 (6%)	0	100	100
2	AX	153/522 (29%)	145 (95%)	8 (5%)	0	100	100
2	BX	153/522 (29%)	145 (95%)	8 (5%)	0	100	100
2	CX	153/522 (29%)	145 (95%)	8 (5%)	0	100	100
2	DX	153/522 (29%)	145 (95%)	8 (5%)	0	100	100
2	EX	153/522 (29%)	145 (95%)	8 (5%)	0	100	100
2	FX	153/522 (29%)	145 (95%)	8 (5%)	0	100	100
2	GX	153/522 (29%)	145 (95%)	8 (5%)	0	100	100
2	HX	153/522 (29%)	145 (95%)	8 (5%)	0	100	100
2	IX	153/522 (29%)	145 (95%)	8 (5%)	0	100	100
2	JX	153/522 (29%)	145 (95%)	8 (5%)	0	100	100
2	KX	153/522 (29%)	145 (95%)	8 (5%)	0	100	100
2	LX	153/522 (29%)	145 (95%)	8 (5%)	0	100	100
2	MX	153/522 (29%)	145 (95%)	8 (5%)	0	100	100
2	NX	153/522 (29%)	145 (95%)	8 (5%)	0	100	100
3	AY	205/1927 (11%)	195 (95%)	10 (5%)	0	100	100
3	BY	205/1927 (11%)	195 (95%)	10 (5%)	0	100	100
3	CY	205/1927 (11%)	195 (95%)	10 (5%)	0	100	100
3	DY	205/1927 (11%)	195 (95%)	10 (5%)	0	100	100
3	EY	205/1927 (11%)	195 (95%)	10 (5%)	0	100	100
3	FY	205/1927 (11%)	195 (95%)	10 (5%)	0	100	100
3	GY	205/1927 (11%)	195 (95%)	10 (5%)	0	100	100
3	HY	205/1927 (11%)	195 (95%)	10 (5%)	0	100	100
3	IY	205/1927 (11%)	195 (95%)	10 (5%)	0	100	100
3	JY	205/1927 (11%)	195 (95%)	10 (5%)	0	100	100
3	KY	205/1927 (11%)	195 (95%)	10 (5%)	0	100	100
3	LY	205/1927 (11%)	195 (95%)	10 (5%)	0	100	100
3	MY	205/1927 (11%)	195 (95%)	10 (5%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	NY	205/1927 (11%)	195 (95%)	10 (5%)	0	100	100
All	All	7406/38178 (19%)	7000 (94%)	406 (6%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains (i)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	AT	169/254 (66%)	167 (99%)	2 (1%)	71	88
1	BT	169/254 (66%)	167 (99%)	2 (1%)	71	88
1	CT	169/254 (66%)	167 (99%)	2 (1%)	71	88
1	DT	169/254 (66%)	167 (99%)	2 (1%)	71	88
1	ET	169/254 (66%)	167 (99%)	2 (1%)	71	88
1	FT	169/254 (66%)	167 (99%)	2 (1%)	71	88
1	GT	169/254 (66%)	167 (99%)	2 (1%)	71	88
1	HT	169/254 (66%)	167 (99%)	2 (1%)	71	88
1	IT	169/254 (66%)	167 (99%)	2 (1%)	71	88
1	JT	169/254 (66%)	167 (99%)	2 (1%)	71	88
1	KT	169/254 (66%)	167 (99%)	2 (1%)	71	88
1	LT	169/254 (66%)	167 (99%)	2 (1%)	71	88
1	MT	169/254 (66%)	167 (99%)	2 (1%)	71	88
1	NT	169/254 (66%)	167 (99%)	2 (1%)	71	88
2	AX	139/470 (30%)	139 (100%)	0	100	100
2	BX	139/470 (30%)	139 (100%)	0	100	100
2	CX	139/470 (30%)	139 (100%)	0	100	100
2	DX	139/470 (30%)	139 (100%)	0	100	100
2	EX	139/470 (30%)	139 (100%)	0	100	100
2	FX	139/470 (30%)	139 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	GX	139/470 (30%)	139 (100%)	0	100	100
2	HX	139/470 (30%)	139 (100%)	0	100	100
2	IX	139/470 (30%)	139 (100%)	0	100	100
2	JX	139/470 (30%)	139 (100%)	0	100	100
2	KX	139/470 (30%)	139 (100%)	0	100	100
2	LX	139/470 (30%)	139 (100%)	0	100	100
2	MX	139/470 (30%)	139 (100%)	0	100	100
2	NX	139/470 (30%)	139 (100%)	0	100	100
3	AY	177/1724 (10%)	176 (99%)	1 (1%)	86	94
3	BY	177/1724 (10%)	176 (99%)	1 (1%)	86	94
3	CY	177/1724 (10%)	176 (99%)	1 (1%)	86	94
3	DY	177/1724 (10%)	176 (99%)	1 (1%)	86	94
3	EY	177/1724 (10%)	176 (99%)	1 (1%)	86	94
3	FY	177/1724 (10%)	176 (99%)	1 (1%)	86	94
3	GY	177/1724 (10%)	176 (99%)	1 (1%)	86	94
3	HY	177/1724 (10%)	176 (99%)	1 (1%)	86	94
3	IY	177/1724 (10%)	176 (99%)	1 (1%)	86	94
3	JY	177/1724 (10%)	176 (99%)	1 (1%)	86	94
3	KY	177/1724 (10%)	176 (99%)	1 (1%)	86	94
3	LY	177/1724 (10%)	176 (99%)	1 (1%)	86	94
3	MY	177/1724 (10%)	176 (99%)	1 (1%)	86	94
3	NY	177/1724 (10%)	176 (99%)	1 (1%)	86	94
All	All	6790/34272 (20%)	6748 (99%)	42 (1%)	86	94

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	AT	206	LYS
1	AT	214	LYS
3	AY	1906	LYS
1	BT	206	LYS
1	BT	214	LYS
3	BY	1906	LYS
1	CT	206	LYS

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Mol	Chain	Res	Type
1	CT	214	LYS
3	CY	1906	LYS
1	DT	206	LYS
1	DT	214	LYS
3	DY	1906	LYS
1	ET	206	LYS
1	ET	214	LYS
3	EY	1906	LYS
1	FT	206	LYS
1	FT	214	LYS
3	FY	1906	LYS
1	GT	206	LYS
1	GT	214	LYS
3	GY	1906	LYS
1	HT	206	LYS
1	HT	214	LYS
3	HY	1906	LYS
1	IT	206	LYS
1	IT	214	LYS
3	IY	1906	LYS
1	JT	206	LYS
1	JT	214	LYS
3	JY	1906	LYS
1	KT	206	LYS
1	KT	214	LYS
3	KY	1906	LYS
1	LT	206	LYS
1	LT	214	LYS
3	LY	1906	LYS
1	MT	206	LYS
1	MT	214	LYS
3	MY	1906	LYS
1	NT	206	LYS
1	NT	214	LYS
3	NY	1906	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

There are no ligands in this entry.

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

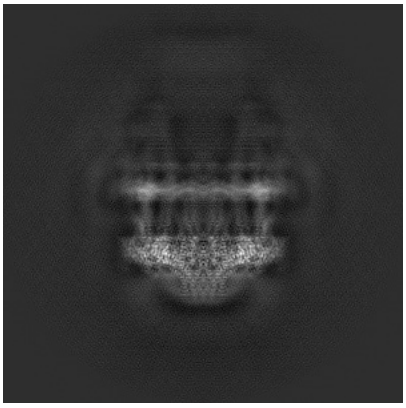
6 Map visualisation (i)

This section contains visualisations of the EMDB entry EMD-22076. These allow visual inspection of the internal detail of the map and identification of artifacts.

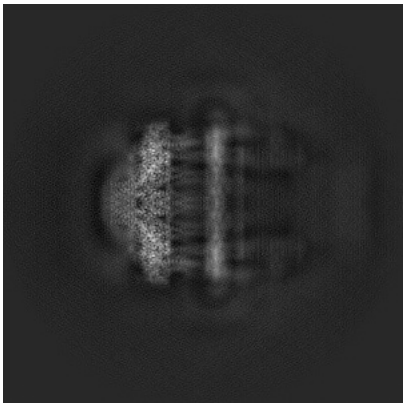
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections (i)

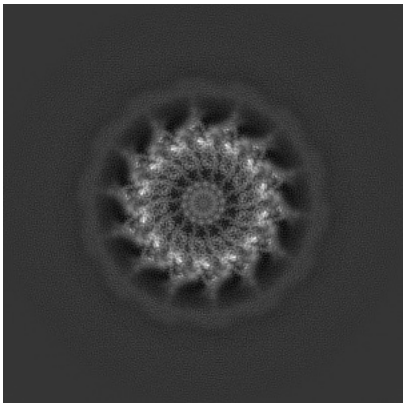
6.1.1 Primary map



X



Y

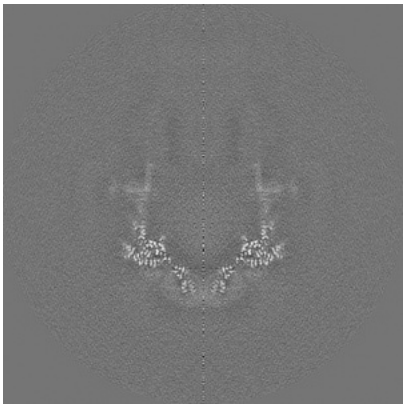


Z

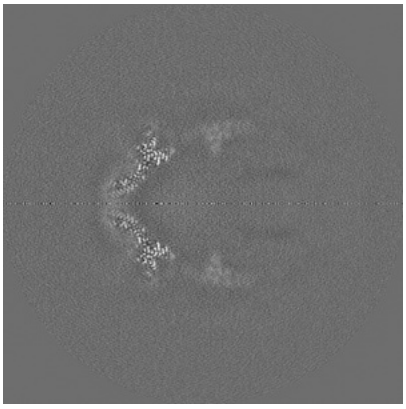
The images above show the map projected in three orthogonal directions.

6.2 Central slices (i)

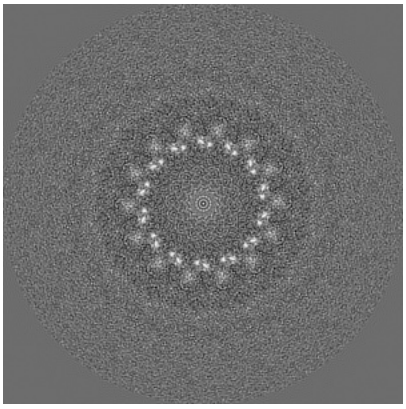
6.2.1 Primary map



X Index: 255



Y Index: 255

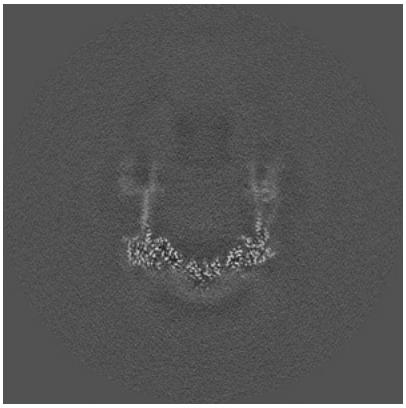


Z Index: 255

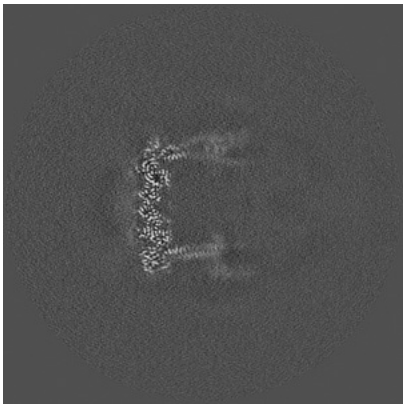
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices (i)

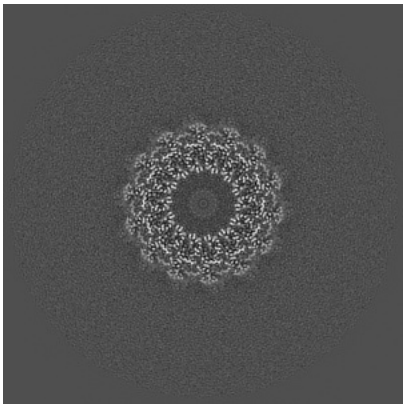
6.3.1 Primary map



X Index: 224



Y Index: 303

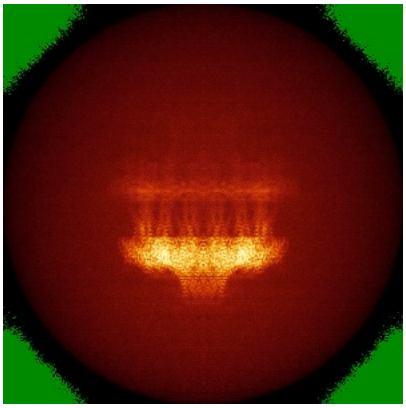


Z Index: 188

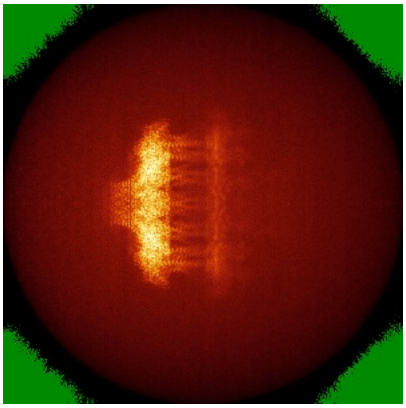
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) (i)

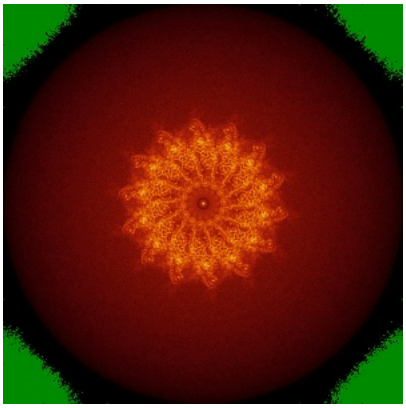
6.4.1 Primary map



X



Y



Z

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

6.5 Orthogonal surface views (i)

6.5.1 Primary map

X Y Z

The images above show the 3D surface view of the map at the recommended contour level 0.015. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

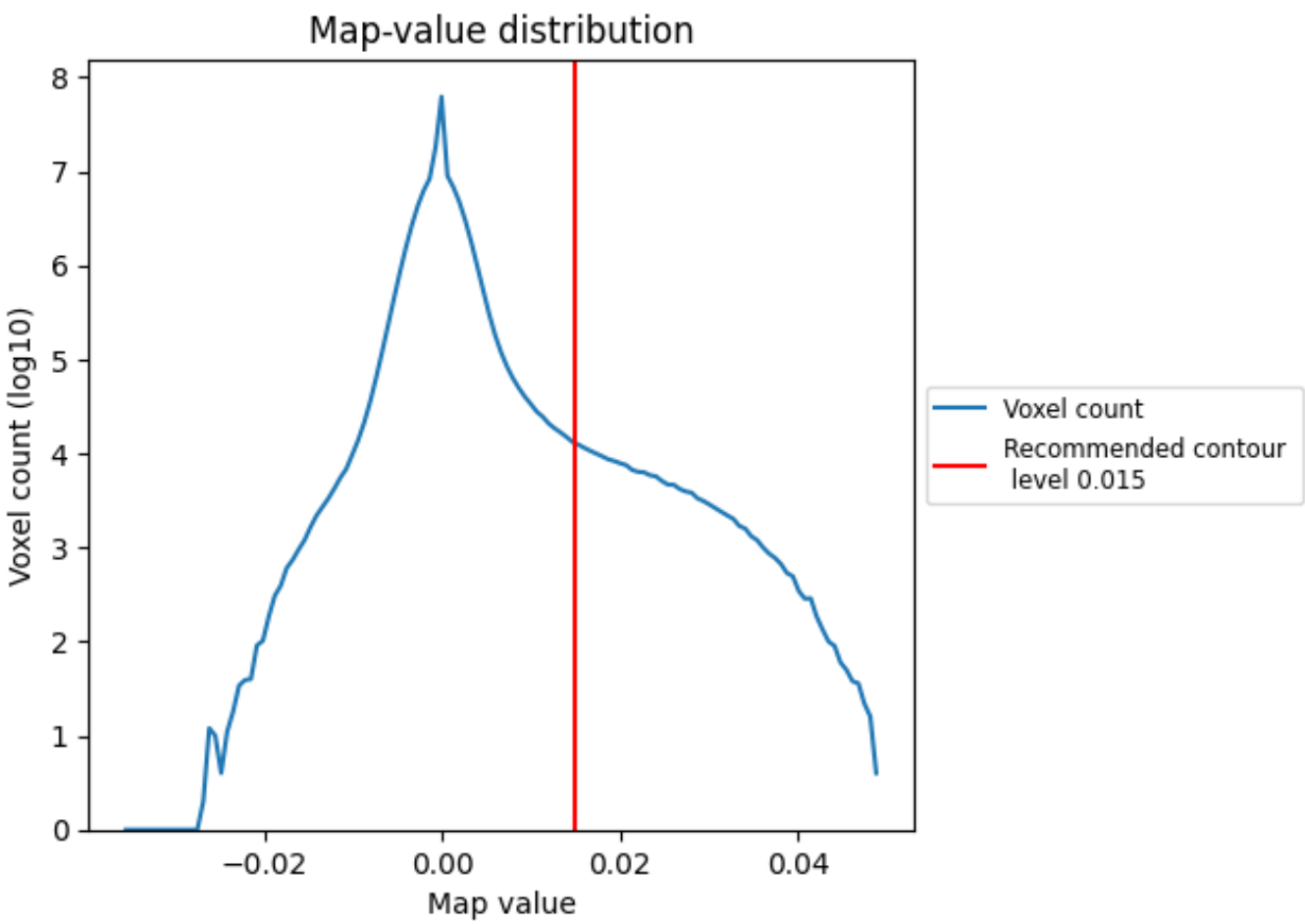
6.6 Mask visualisation (i)

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

7 Map analysis (i)

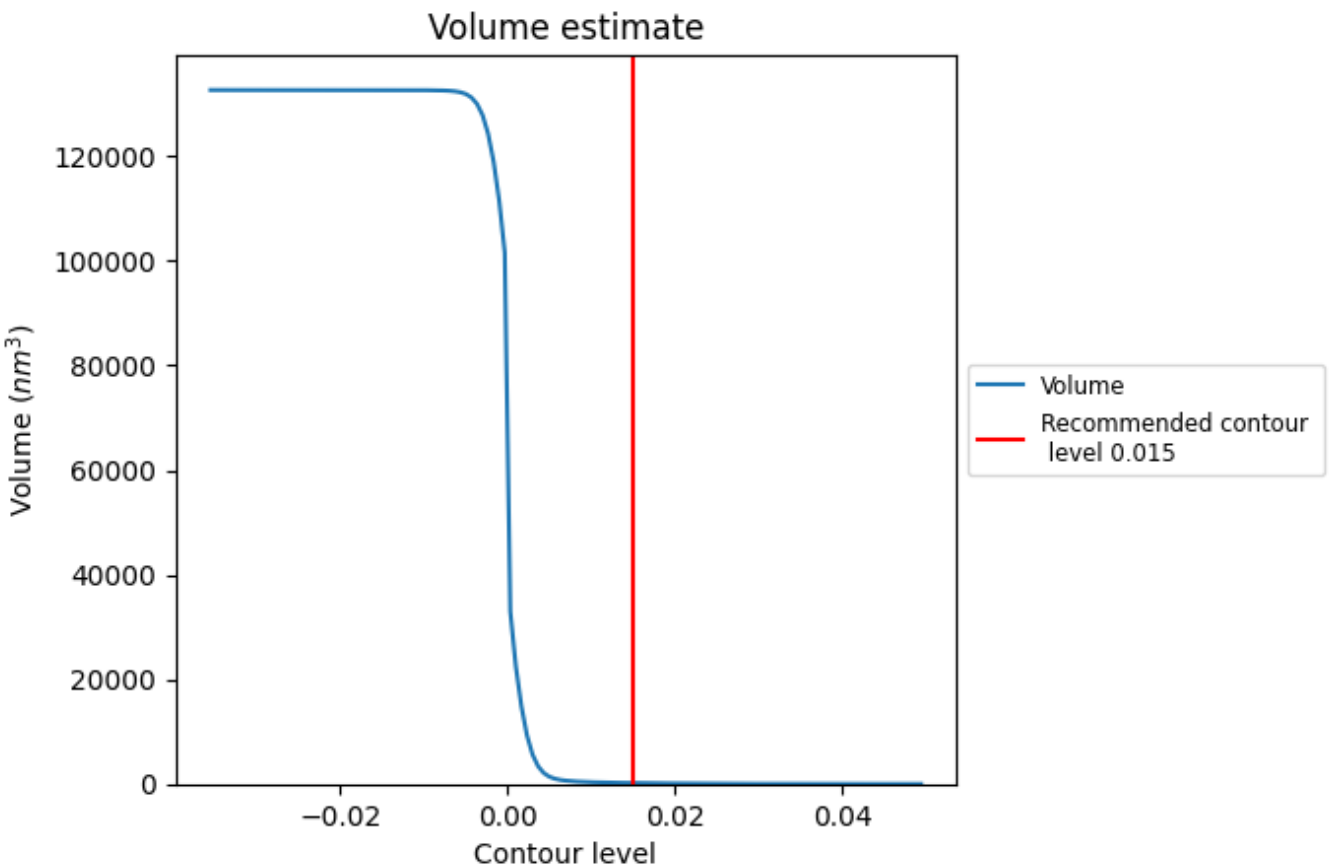
This section contains the results of statistical analysis of the map.

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.

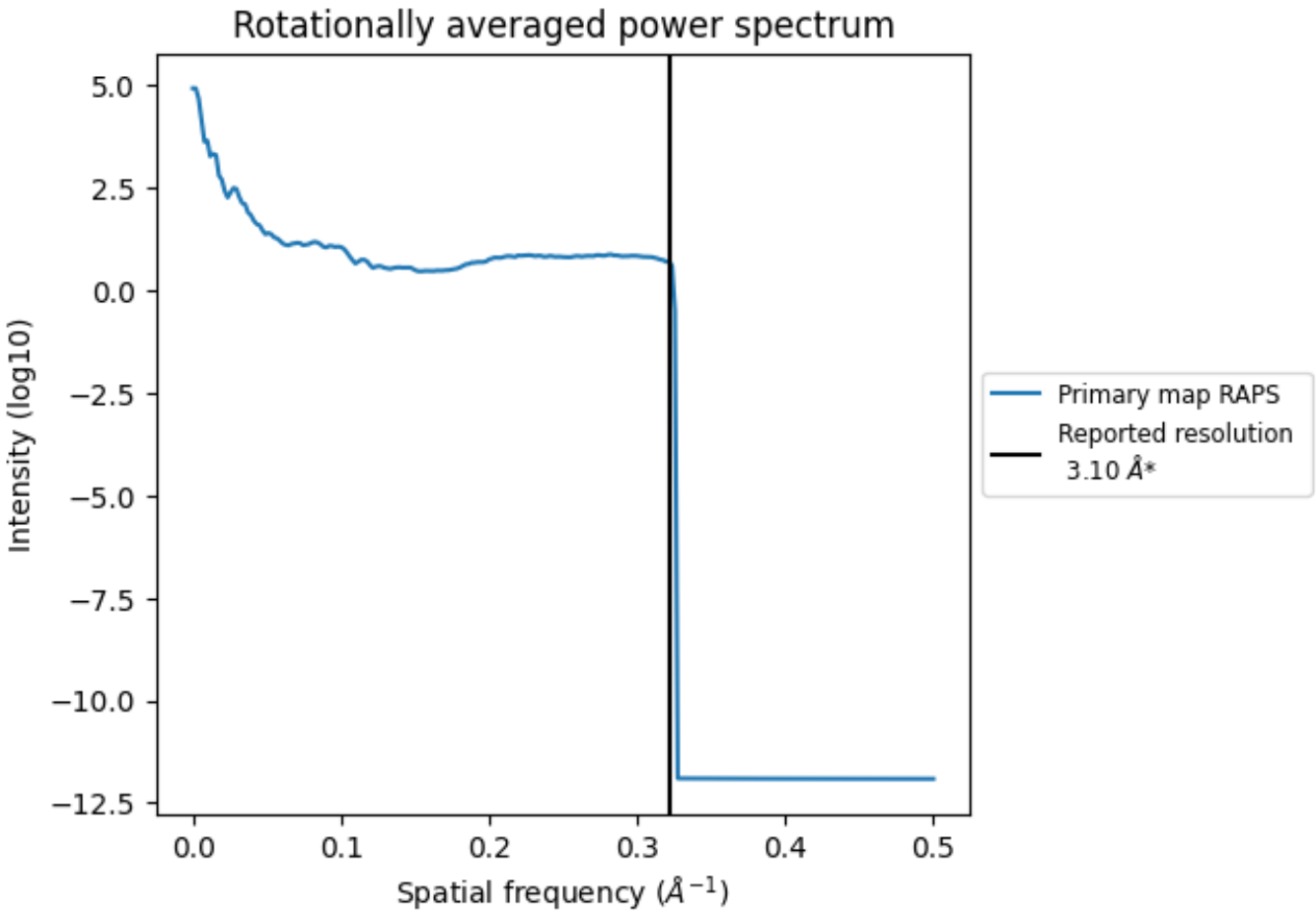
7.2 Volume estimate (i)



The volume at the recommended contour level is 180 nm^3; this corresponds to an approximate mass of 163 kDa.

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

7.3 Rotationally averaged power spectrum (i)

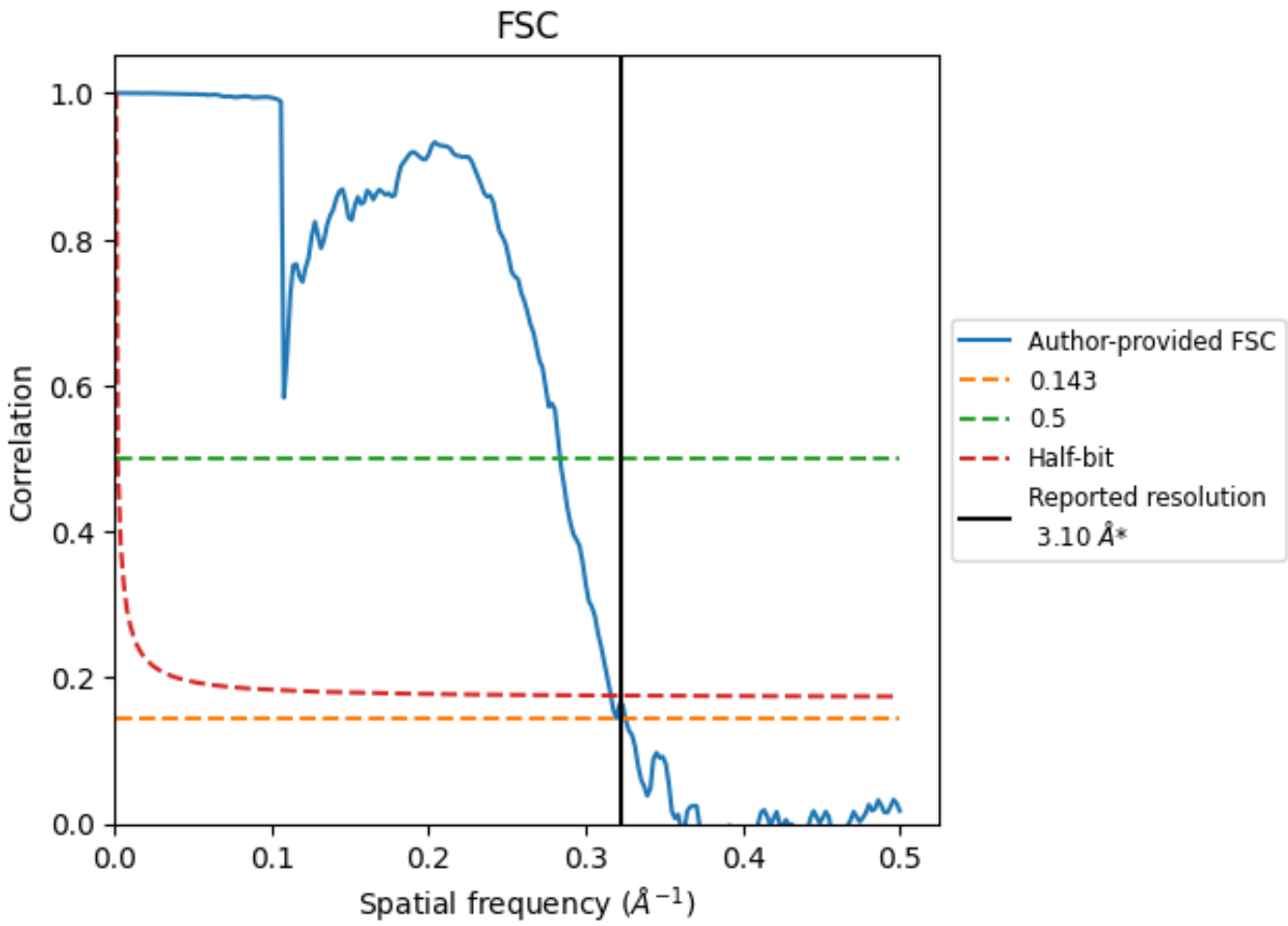


*Reported resolution corresponds to spatial frequency of 0.323 Å^{-1}

8 Fourier-Shell correlation (i)

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

8.1 FSC (i)



*Reported resolution corresponds to spatial frequency of 0.323 Å^{-1}

8.2 Resolution estimates (i)

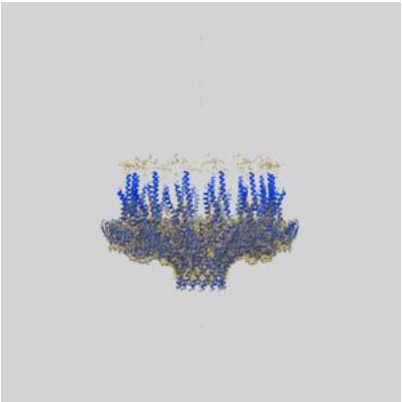
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.10	-	-
Author-provided FSC curve	3.07	3.52	3.16
Unmasked-calculated*	-	-	-

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps.

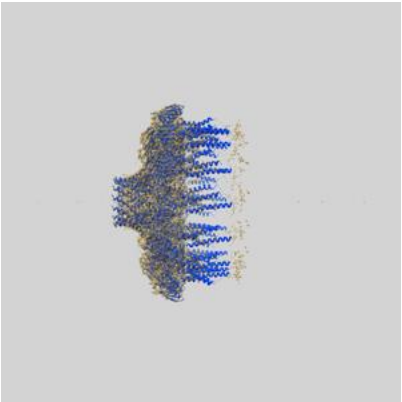
9 Map-model fit (i)

This section contains information regarding the fit between EMDB map EMD-22076 and PDB model 6X6K. Per-residue inclusion information can be found in section 3 on page 8.

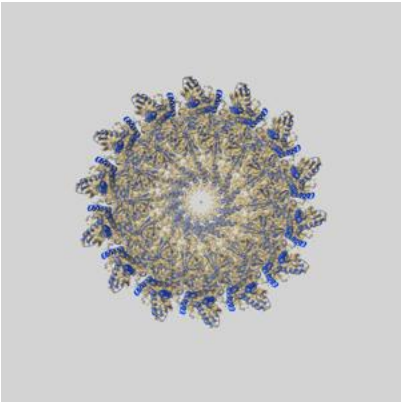
9.1 Map-model overlay (i)



X



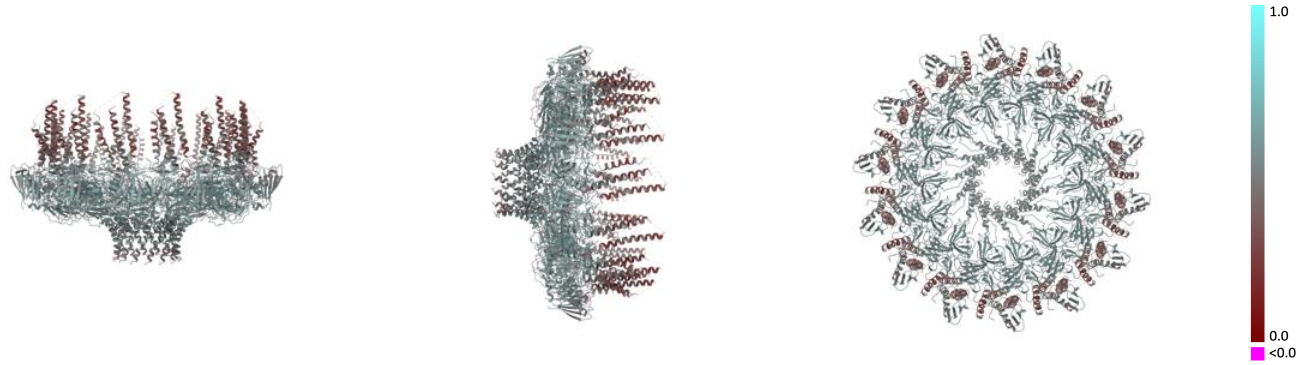
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.015 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

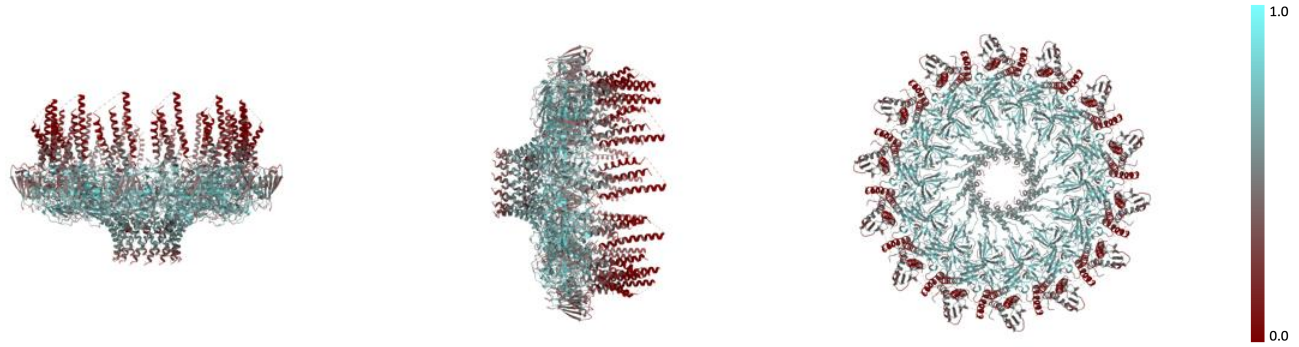
9.2 Q-score mapped to coordinate model (i)



X Y Z

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

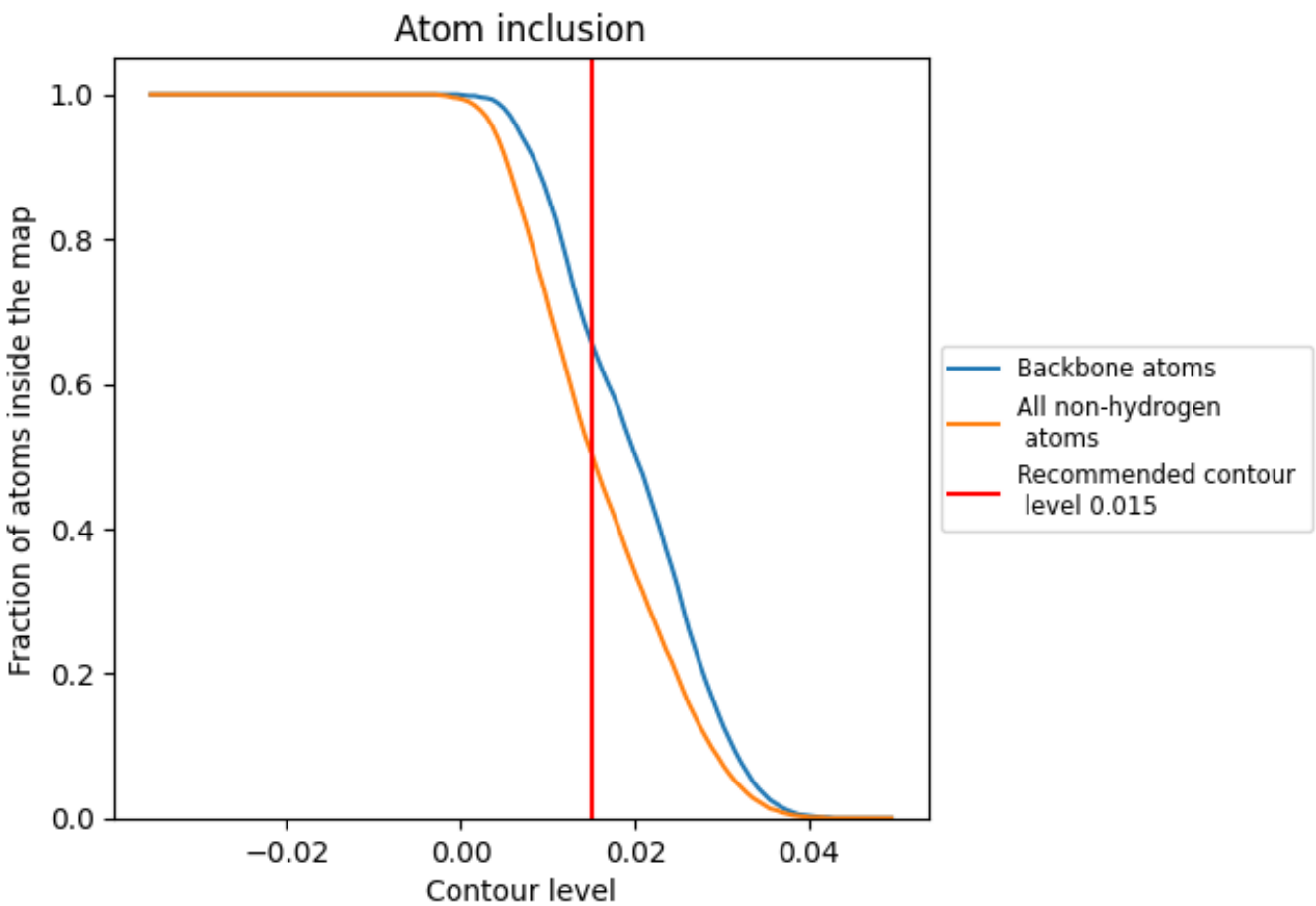
9.3 Atom inclusion mapped to coordinate model (i)



X Y Z

The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.015).

9.4 Atom inclusion (i)



At the recommended contour level, 66% of all backbone atoms, 50% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary (i)

The table lists the average atom inclusion at the recommended contour level (0.015) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.5050	0.5040
AT	0.3380	0.4550
AX	0.5770	0.5170
AY	0.6070	0.5400
BT	0.3390	0.4580
BX	0.5750	0.5160
BY	0.6070	0.5410
CT	0.3350	0.4550
CX	0.5740	0.5180
CY	0.6040	0.5410
DT	0.3390	0.4550
DX	0.5720	0.5190
DY	0.6020	0.5420
ET	0.3400	0.4550
EX	0.5790	0.5170
EY	0.6040	0.5430
FT	0.3360	0.4540
FX	0.5780	0.5190
FY	0.6050	0.5400
GT	0.3330	0.4530
GX	0.5800	0.5190
GY	0.6090	0.5410
HT	0.3370	0.4560
HX	0.5770	0.5170
HY	0.6060	0.5400
IT	0.3370	0.4560
IX	0.5710	0.5190
IY	0.6070	0.5400
JT	0.3350	0.4520
JX	0.5740	0.5170
JY	0.6050	0.5390
KT	0.3360	0.4550
KX	0.5730	0.5170
KY	0.6020	0.5400
LT	0.3370	0.4560

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Chain	Atom inclusion	Q-score
LX	0.5790	0.5170
LY	0.6050	0.5390
MT	0.3350	0.4570
MX	0.5770	0.5190
MY	0.6050	0.5390
NT	0.3390	0.4560
NX	0.5790	0.5170
NY	0.6090	0.5400