Full wwPDB EM Validation Report

Mar 8, 2026 – 06:33 AM UTC

PDB ID	:	7MUC / pdb_00007muc
EMDB ID	:	EMD-24004
Title	:	Legionella pneumophila Dot/Icm T4SS C1 Reconstruction
Authors	:	Sheedlo, M.J.; Durie, C.L.; Swanson, M.; Lacy, D.B.; Ohi, M.D.
Deposited on	:	2021-05-14
Resolution	:	3.80 Å(reported)

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

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/EMValidationReportHelp
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

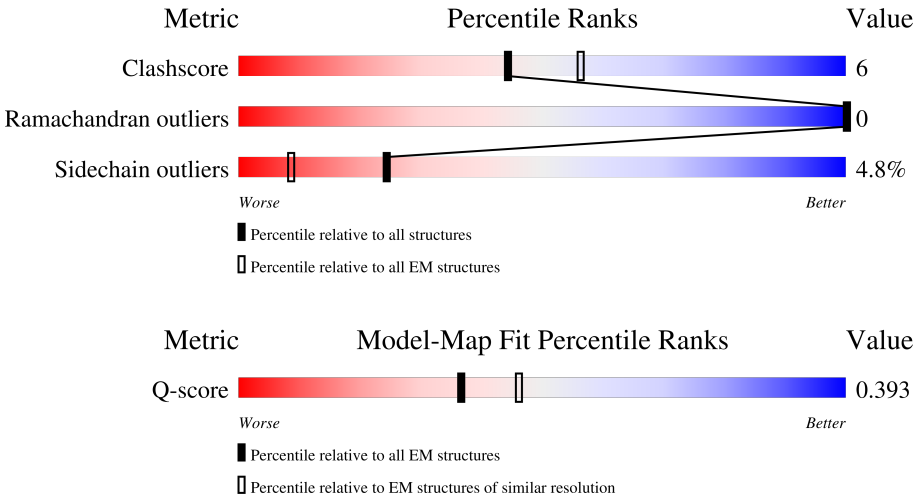
EMDB validation analysis	:	0.0.1.dev132
MolProbity	:	4-5-2 with Phenix2.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics	:	202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ	:	1.9.13
Ideal geometry (proteins)	:	Engh &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

1 Overall quality at a glance

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

The reported resolution of this entry is 3.80 Å.

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



Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	10198 (3.30 - 4.30)

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	AC	303	57% 11% 30%
1	BC	303	58% 10% 30%
1	CC	303	58% 10% 30%
1	DC	303	56% 11% 30%

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Mol	Chain	Length	Quality of chain
1	EC	303	58% 10% 30%
1	FC	303	57% 10% 30%
1	GC	303	57% 11% 30%
1	HC	303	57% 11% 30%
1	IC	303	57% 11% 30%
1	JC	303	57% 10% 30%
1	KC	303	56% 11% 30%
1	LC	303	58% 9% 30%
1	MC	303	57% 10% 30%
2	AD	163	75% 10% 14%
2	Ad	163	76% 7% 16%
2	BD	163	77% 7% 14%
2	Bd	163	71% 12% 16%
2	CD	163	69% 16% 14%
2	Cd	163	71% 12% 16%
2	DD	163	72% 13% 14%
2	Dd	163	77% 7% 16%
2	ED	163	74% 11% 14%
2	Ed	163	73% 11% 16%
2	FD	163	78% 7% 14%
2	Fd	163	72% 10% 16%
2	GD	163	75% 10% 14%
2	Gd	163	75% 9% 16%
2	HD	163	74% 12% 14%
2	Hd	163	77% 6% 16%

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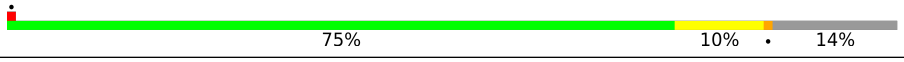
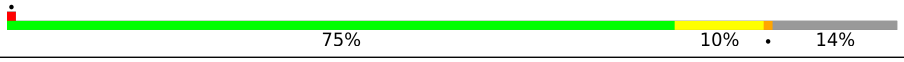
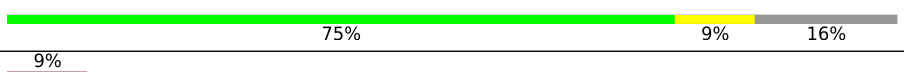
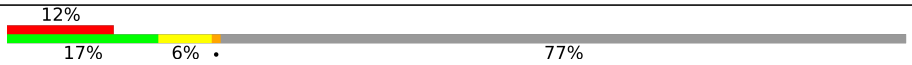
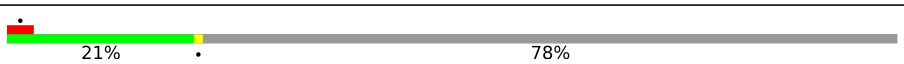
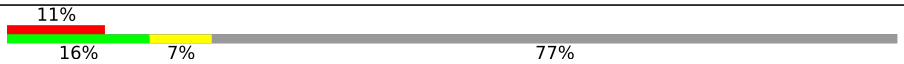
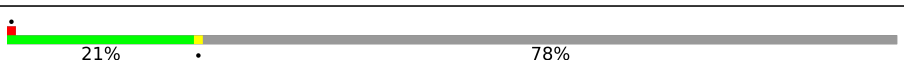
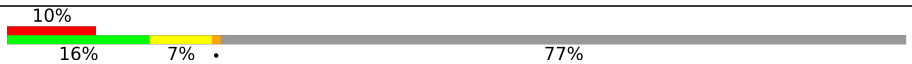
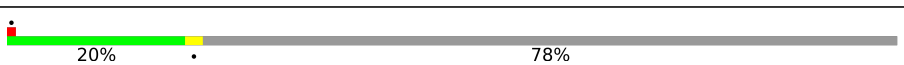
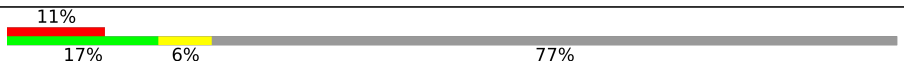
Mol	Chain	Length	Quality of chain
2	ID	163	
2	Id	163	
2	JD	163	
2	Jd	163	
2	KD	163	
2	Kd	163	
2	LD	163	
2	Ld	163	
2	MD	163	
2	Md	163	
3	AF	269	
3	Af	269	
3	BF	269	
3	Bf	269	
3	CF	269	
3	Cf	269	
3	DF	269	
3	Df	269	
3	EF	269	
3	Ef	269	
3	FF	269	
3	Ff	269	
3	GF	269	
3	Gf	269	
3	HF	269	

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Mol	Chain	Length	Quality of chain
3	Hf	269	17% 78%
3	IF	269	10% 17% 6% 77%
3	If	269	20% 78%
3	JF	269	9% 16% 7% 77%
3	Jf	269	19% 78%
3	KF	269	6% 17% 6% 77%
3	Kf	269	18% 78%
3	LF	269	10% 17% 6% 77%
3	Lf	269	19% 78%
3	MF	269	10% 17% 6% 77%
3	Mf	269	20% 78%
3	VF	269	7% 17% 6% 77%
3	WF	269	10% 16% 6% 77%
3	XF	269	9% 16% 6% 77%
3	YF	269	9% 17% 6% 77%
3	ZF	269	10% 16% 6% 77%
4	AG	1048	97%
4	BG	1048	97%
4	CG	1048	97%
4	DG	1048	97%
4	EG	1048	97%
4	FG	1048	97%
4	GG	1048	97%
4	HG	1048	97%
4	IG	1048	97%

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Mol	Chain	Length	Quality of chain
4	JG	1048	97%
4	KG	1048	97%
4	LG	1048	97%
4	MG	1048	97%
4	VG	1048	97%
4	WG	1048	97%
4	XG	1048	97%
4	YG	1048	97%
4	ZG	1048	97%
5	AH	361	7% 59% 11% 29%
5	BH	361	5% 59% 11% 29%
5	CH	361	5% 58% 12% 29%
5	DH	361	5% 61% 9% 29%
5	EH	361	7% 55% 16% 29%
5	FH	361	7% 61% 9% 29%
5	GH	361	5% 58% 12% 29%
5	HH	361	5% 59% 11% 29%
5	IH	361	6% 58% 12% 29%
5	JH	361	5% 60% 11% 29%
5	KH	361	6% 61% 10% 29%
5	LH	361	7% 58% 11% 29%
5	MH	361	6% 59% 12% 29%
5	VH	361	14% 60% 7% 33%
5	WH	361	27% 55% 11% 33%
5	XH	361	14% 57% 9% 33%

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Mol	Chain	Length	Quality of chain
5	YH	361	20% 59% 8% 33%
5	ZH	361	25% 56% 9% 33%
6	AK	189	69% 8% 20%
6	BK	189	68% 10% 20%
6	CK	189	68% 10% 20%
6	DK	189	68% 9% 20%
6	EK	189	67% 11% 20%
6	FK	189	69% 10% 20%
6	GK	189	69% 9% 20%
6	HK	189	69% 9% 20%
6	IK	189	67% 11% 20%
6	JK	189	68% 10% 20%
6	KK	189	68% 10% 20%
6	LK	189	69% 10% 20%
6	MK	189	69% 10% 20%
7	AL	249	57% 10% 31%
7	BL	249	56% 11% 31%
7	CL	249	56% 11% 31%
7	DL	249	56% 11% 31%
7	EL	249	57% 10% 31%
7	FL	249	56% 10% 31%
7	GL	249	56% 10% 31%
7	HL	249	56% 10% 31%
7	IL	249	56% 10% 31%
7	JL	249	56% 11% 31%

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Mol	Chain	Length	Quality of chain
7	KL	249	56% 10% 31%
7	LL	249	56% 11% 31%
7	ML	249	56% 10% 31%
8	AM	320	25% 71% 18% 9%
8	BM	320	24% 70% 19% 9%
8	CM	320	26% 70% 19% 9%
8	DM	320	25% 70% 19% 9%
8	EM	320	26% 70% 19% 9%
8	FM	320	25% 69% 20% 9%
8	GM	320	27% 72% 17% 9%
8	HM	320	26% 70% 19% 9%
8	IM	320	28% 69% 20% 9%
8	JM	320	27% 71% 18% 9%
8	KM	320	27% 69% 20% 9%
8	LM	320	22% 70% 19% 9%
8	MM	320	22% 70% 19% 9%
9	AN	124	5% 48% 11% 37%
9	BN	124	50% 10% 37%
9	CN	124	7% 52% 8% 37%
9	DN	124	5% 50% 10% 37%
9	EN	124	6% 51% 8% 37%
9	FN	124	7% 48% 11% 37%
9	GN	124	6% 50% 10% 37%
9	HN	124	5% 50% 10% 37%
9	IN	124	6% 52% 7% 37%

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Mol	Chain	Length	Quality of chain
9	JN	124	5% 51% 9% 37%
9	KN	124	5% 50% 10% 37%
9	LN	124	51% 10% 37%
9	MN	124	5% 51% 10% 37%
10	AU	9	100%
10	BU	9	100%
10	CU	9	100%
10	DU	9	100%
10	EU	9	100%
10	FU	9	100%
10	GU	9	100%
10	HU	9	100%
10	IU	9	100%
10	JU	9	100%
10	KU	9	100%
10	LU	9	100%
10	MU	9	100%
11	AX	48	85% 100%
11	BX	48	88% 100%
11	CX	48	85% 100%
11	DX	48	81% 100%
11	EX	48	79% 100%
11	FX	48	90% 100%
11	GX	48	90% 100%
11	HX	48	85% 100%

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Mol	Chain	Length	Quality of chain
11	IX	48	85% 100%
11	JX	48	90% 100%
11	KX	48	75% 98%
11	LX	48	79% 100%
11	MX	48	85% 100%
11	VX	48	81% 100%
11	WX	48	88% 100%
11	XX	48	83% 100%
11	YX	48	83% 100%
11	ZX	48	83% 100%

2 Entry composition (i)

There are 11 unique types of molecules in this entry. The entry contains 180380 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 DotC.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	AC	213	Total	C	N	O	S	0	0
			1685	1068	295	316	6		
1	BC	213	Total	C	N	O	S	0	0
			1685	1068	295	316	6		
1	CC	213	Total	C	N	O	S	0	0
			1685	1068	295	316	6		
1	DC	213	Total	C	N	O	S	0	0
			1685	1068	295	316	6		
1	EC	213	Total	C	N	O	S	0	0
			1685	1068	295	316	6		
1	FC	213	Total	C	N	O	S	0	0
			1685	1068	295	316	6		
1	GC	213	Total	C	N	O	S	0	0
			1685	1068	295	316	6		
1	HC	213	Total	C	N	O	S	0	0
			1685	1068	295	316	6		
1	IC	213	Total	C	N	O	S	0	0
			1685	1068	295	316	6		
1	JC	213	Total	C	N	O	S	0	0
			1685	1068	295	316	6		
1	KC	213	Total	C	N	O	S	0	0
			1685	1068	295	316	6		
1	LC	213	Total	C	N	O	S	0	0
			1685	1068	295	316	6		
1	MC	213	Total	C	N	O	S	0	0
			1685	1068	295	316	6		

- Molecule 2 is a protein called DotD.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	AD	140	Total	C	N	O	S	0	0
			1086	692	185	206	3		
2	Ad	137	Total	C	N	O	S	0	0
			1058	672	182	202	2		

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Mol	Chain	Residues	Atoms					AltConf	Trace
2	BD	140	Total	C	N	O	S	0	0
			1086	692	185	206	3		
2	Bd	137	Total	C	N	O	S	0	0
			1058	672	182	202	2		
2	CD	140	Total	C	N	O	S	0	0
			1086	692	185	206	3		
2	Cd	137	Total	C	N	O	S	0	0
			1058	672	182	202	2		
2	DD	140	Total	C	N	O	S	0	0
			1086	692	185	206	3		
2	Dd	137	Total	C	N	O	S	0	0
			1058	672	182	202	2		
2	ED	140	Total	C	N	O	S	0	0
			1086	692	185	206	3		
2	Ed	137	Total	C	N	O	S	0	0
			1058	672	182	202	2		
2	FD	140	Total	C	N	O	S	0	0
			1086	692	185	206	3		
2	Fd	137	Total	C	N	O	S	0	0
			1058	672	182	202	2		
2	GD	140	Total	C	N	O	S	0	0
			1086	692	185	206	3		
2	Gd	137	Total	C	N	O	S	0	0
			1058	672	182	202	2		
2	HD	140	Total	C	N	O	S	0	0
			1086	692	185	206	3		
2	Hd	137	Total	C	N	O	S	0	0
			1058	672	182	202	2		
2	ID	140	Total	C	N	O	S	0	0
			1086	692	185	206	3		
2	Id	137	Total	C	N	O	S	0	0
			1058	672	182	202	2		
2	JD	140	Total	C	N	O	S	0	0
			1086	692	185	206	3		
2	Jd	137	Total	C	N	O	S	0	0
			1058	672	182	202	2		
2	KD	140	Total	C	N	O	S	0	0
			1086	692	185	206	3		
2	Kd	137	Total	C	N	O	S	0	0
			1058	672	182	202	2		
2	LD	140	Total	C	N	O	S	0	0
			1086	692	185	206	3		

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Mol	Chain	Residues	Atoms					AltConf	Trace
2	Ld	137	Total	C	N	O	S	0	0
			1058	672	182	202	2		
2	MD	140	Total	C	N	O	S	0	0
			1086	692	185	206	3		
2	Md	137	Total	C	N	O	S	0	0
			1058	672	182	202	2		

- Molecule 3 is a protein called DotF.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	AF	63	Total	C	N	O	S	0	0
			483	308	84	90	1		
3	Af	59	Total	C	N	O	S	0	0
			449	290	77	81	1		
3	BF	63	Total	C	N	O	S	0	0
			483	308	84	90	1		
3	Bf	59	Total	C	N	O	S	0	0
			449	290	77	81	1		
3	CF	63	Total	C	N	O	S	0	0
			483	308	84	90	1		
3	Cf	59	Total	C	N	O	S	0	0
			449	290	77	81	1		
3	DF	63	Total	C	N	O	S	0	0
			483	308	84	90	1		
3	Df	59	Total	C	N	O	S	0	0
			449	290	77	81	1		
3	EF	63	Total	C	N	O	S	0	0
			483	308	84	90	1		
3	Ef	59	Total	C	N	O	S	0	0
			449	290	77	81	1		
3	FF	63	Total	C	N	O	S	0	0
			483	308	84	90	1		
3	Ff	59	Total	C	N	O	S	0	0
			449	290	77	81	1		
3	GF	63	Total	C	N	O	S	0	0
			483	308	84	90	1		
3	Gf	59	Total	C	N	O	S	0	0
			449	290	77	81	1		
3	HF	63	Total	C	N	O	S	0	0
			483	308	84	90	1		
3	Hf	59	Total	C	N	O	S	0	0
			449	290	77	81	1		

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Mol	Chain	Residues	Atoms					AltConf	Trace
3	IF	63	Total	C	N	O	S	0	0
			483	308	84	90	1		
3	If	59	Total	C	N	O	S	0	0
			449	290	77	81	1		
3	JF	63	Total	C	N	O	S	0	0
			483	308	84	90	1		
3	Jf	59	Total	C	N	O	S	0	0
			449	290	77	81	1		
3	KF	63	Total	C	N	O	S	0	0
			483	308	84	90	1		
3	Kf	59	Total	C	N	O	S	0	0
			449	290	77	81	1		
3	LF	63	Total	C	N	O	S	0	0
			483	308	84	90	1		
3	Lf	59	Total	C	N	O	S	0	0
			449	290	77	81	1		
3	MF	63	Total	C	N	O	S	0	0
			483	308	84	90	1		
3	Mf	59	Total	C	N	O	S	0	0
			449	290	77	81	1		
3	VF	63	Total	C	N	O	S	0	0
			483	308	84	90	1		
3	WF	63	Total	C	N	O	S	0	0
			483	308	84	90	1		
3	XF	63	Total	C	N	O	S	0	0
			483	308	84	90	1		
3	YF	63	Total	C	N	O	S	0	0
			483	308	84	90	1		
3	ZF	63	Total	C	N	O	S	0	0
			483	308	84	90	1		

- Molecule 4 is a protein called IcmE protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	AG	34	Total	C	N	O	S	0	0
			276	168	47	60	1		
4	BG	34	Total	C	N	O	S	0	0
			276	168	47	60	1		
4	CG	34	Total	C	N	O	S	0	0
			276	168	47	60	1		
4	DG	34	Total	C	N	O	S	0	0
			276	168	47	60	1		

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Mol	Chain	Residues	Atoms					AltConf	Trace
4	EG	34	Total	C	N	O	S	0	0
			276	168	47	60	1		
4	FG	34	Total	C	N	O	S	0	0
			276	168	47	60	1		
4	GG	34	Total	C	N	O	S	0	0
			276	168	47	60	1		
4	HG	34	Total	C	N	O	S	0	0
			276	168	47	60	1		
4	IG	34	Total	C	N	O	S	0	0
			276	168	47	60	1		
4	JG	34	Total	C	N	O	S	0	0
			276	168	47	60	1		
4	KG	34	Total	C	N	O	S	0	0
			276	168	47	60	1		
4	LG	34	Total	C	N	O	S	0	0
			276	168	47	60	1		
4	MG	34	Total	C	N	O	S	0	0
			276	168	47	60	1		
4	VG	34	Total	C	N	O	S	0	0
			276	168	47	60	1		
4	WG	34	Total	C	N	O	S	0	0
			276	168	47	60	1		
4	XG	34	Total	C	N	O	S	0	0
			276	168	47	60	1		
4	YG	34	Total	C	N	O	S	0	0
			276	168	47	60	1		
4	ZG	34	Total	C	N	O	S	0	0
			276	168	47	60	1		

- Molecule 5 is a protein called Type IV secretion protein IcmK.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	AH	258	Total	C	N	O	S	0	0
			1983	1268	336	371	8		
5	BH	258	Total	C	N	O	S	0	0
			1983	1268	336	371	8		
5	CH	258	Total	C	N	O	S	0	0
			1983	1268	336	371	8		
5	DH	258	Total	C	N	O	S	0	0
			1983	1268	336	371	8		
5	EH	258	Total	C	N	O	S	0	0
			1983	1268	336	371	8		

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

- Molecule 6 is a protein called Inner membrane lipoprotein YiaD.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	AK	151	Total	C	N	O	S	0	0
			1175	747	209	215	4		
6	BK	151	Total	C	N	O	S	0	0
			1175	747	209	215	4		
6	CK	151	Total	C	N	O	S	0	0
			1175	747	209	215	4		
6	DK	151	Total	C	N	O	S	0	0
			1175	747	209	215	4		
6	EK	151	Total	C	N	O	S	0	0
			1175	747	209	215	4		
6	FK	151	Total	C	N	O	S	0	0
			1175	747	209	215	4		

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Mol	Chain	Residues	Atoms					AltConf	Trace
6	GK	151	Total	C	N	O	S	0	0
			1175	747	209	215	4		
6	HK	151	Total	C	N	O	S	0	0
			1175	747	209	215	4		
6	IK	151	Total	C	N	O	S	0	0
			1175	747	209	215	4		
6	JK	151	Total	C	N	O	S	0	0
			1175	747	209	215	4		
6	KK	151	Total	C	N	O	S	0	0
			1175	747	209	215	4		
6	LK	151	Total	C	N	O	S	0	0
			1175	747	209	215	4		
6	MK	151	Total	C	N	O	S	0	0
			1175	747	209	215	4		

- Molecule 7 is a protein called Outer membrane protein, OmpA family protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	AL	173	Total	C	N	O	S	0	0
			1388	877	253	253	5		
7	BL	173	Total	C	N	O	S	0	0
			1388	877	253	253	5		
7	CL	173	Total	C	N	O	S	0	0
			1388	877	253	253	5		
7	DL	173	Total	C	N	O	S	0	0
			1388	877	253	253	5		
7	EL	173	Total	C	N	O	S	0	0
			1388	877	253	253	5		
7	FL	173	Total	C	N	O	S	0	0
			1388	877	253	253	5		
7	GL	173	Total	C	N	O	S	0	0
			1388	877	253	253	5		
7	HL	173	Total	C	N	O	S	0	0
			1388	877	253	253	5		
7	IL	173	Total	C	N	O	S	0	0
			1388	877	253	253	5		
7	JL	173	Total	C	N	O	S	0	0
			1388	877	253	253	5		
7	KL	173	Total	C	N	O	S	0	0
			1388	877	253	253	5		
7	LL	173	Total	C	N	O	S	0	0
			1388	877	253	253	5		

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Mol	Chain	Residues	Atoms					AltConf	Trace
7	ML	173	Total	C	N	O	S	0	0
			1388	877	253	253	5		

- Molecule 8 is a protein called DUF2807 domain-containing protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	AM	292	Total	C	N	O	S	0	0
			2320	1467	416	433	4		
8	BM	292	Total	C	N	O	S	0	0
			2320	1467	416	433	4		
8	CM	292	Total	C	N	O	S	0	0
			2320	1467	416	433	4		
8	DM	292	Total	C	N	O	S	0	0
			2320	1467	416	433	4		
8	EM	292	Total	C	N	O	S	0	0
			2320	1467	416	433	4		
8	FM	292	Total	C	N	O	S	0	0
			2320	1467	416	433	4		
8	GM	292	Total	C	N	O	S	0	0
			2320	1467	416	433	4		
8	HM	292	Total	C	N	O	S	0	0
			2320	1467	416	433	4		
8	IM	292	Total	C	N	O	S	0	0
			2320	1467	416	433	4		
8	JM	292	Total	C	N	O	S	0	0
			2320	1467	416	433	4		
8	KM	292	Total	C	N	O	S	0	0
			2320	1467	416	433	4		
8	LM	292	Total	C	N	O	S	0	0
			2320	1467	416	433	4		
8	MM	292	Total	C	N	O	S	0	0
			2320	1467	416	433	4		

- Molecule 9 is a protein called Neurogenic locus notch.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	AN	78	Total	C	N	O	S	0	0
			582	357	99	113	13		
9	BN	78	Total	C	N	O	S	0	0
			582	357	99	113	13		
9	CN	78	Total	C	N	O	S	0	0
			582	357	99	113	13		

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

- Molecule 10 is a protein called Unknown protein fragment.

Mol	Chain	Residues	Atoms				AltConf	Trace
10	AU	9	Total	C	N	O	0	0
			45	27	9	9		
10	BU	9	Total	C	N	O	0	0
			45	27	9	9		
10	CU	9	Total	C	N	O	0	0
			45	27	9	9		
10	DU	9	Total	C	N	O	0	0
			45	27	9	9		
10	EU	9	Total	C	N	O	0	0
			45	27	9	9		
10	FU	9	Total	C	N	O	0	0
			45	27	9	9		
10	GU	9	Total	C	N	O	0	0
			45	27	9	9		
10	HU	9	Total	C	N	O	0	0
			45	27	9	9		
10	IU	9	Total	C	N	O	0	0
			45	27	9	9		

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Mol	Chain	Residues	Atoms				AltConf	Trace
10	JU	9	Total 45	C 27	N 9	O 9	0	0
10	KU	9	Total 45	C 27	N 9	O 9	0	0
10	LU	9	Total 45	C 27	N 9	O 9	0	0
10	MU	9	Total 45	C 27	N 9	O 9	0	0

- Molecule 11 is a protein called Unknown protein fragment.

Mol	Chain	Residues	Atoms				AltConf	Trace
11	AX	48	Total 240	C 144	N 48	O 48	0	0
11	BX	48	Total 240	C 144	N 48	O 48	0	0
11	CX	48	Total 240	C 144	N 48	O 48	0	0
11	DX	48	Total 240	C 144	N 48	O 48	0	0
11	EX	48	Total 240	C 144	N 48	O 48	0	0
11	FX	48	Total 240	C 144	N 48	O 48	0	0
11	GX	48	Total 240	C 144	N 48	O 48	0	0
11	HX	48	Total 240	C 144	N 48	O 48	0	0
11	IX	48	Total 240	C 144	N 48	O 48	0	0
11	JX	48	Total 240	C 144	N 48	O 48	0	0
11	KX	48	Total 240	C 144	N 48	O 48	0	0
11	LX	48	Total 240	C 144	N 48	O 48	0	0
11	MX	48	Total 240	C 144	N 48	O 48	0	0
11	VX	48	Total 240	C 144	N 48	O 48	0	0
11	WX	48	Total 240	C 144	N 48	O 48	0	0

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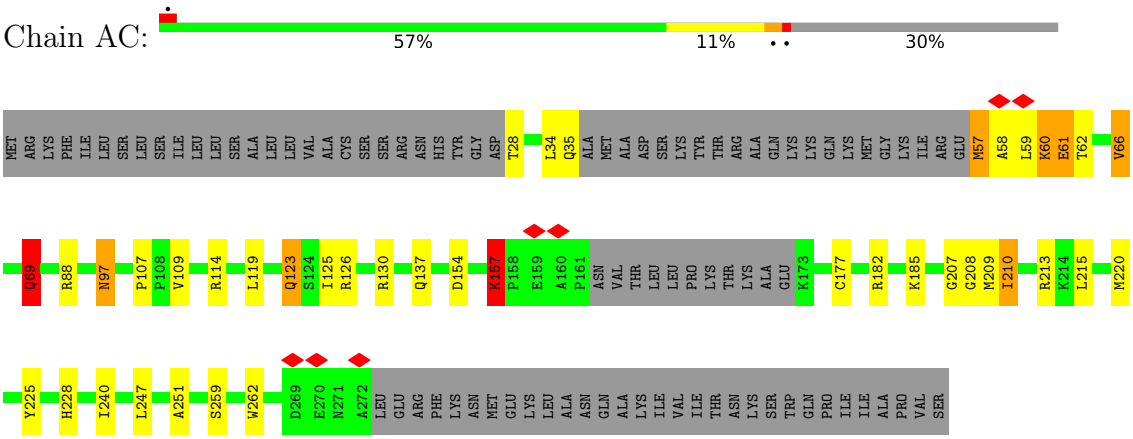
Mol	Chain	Residues	Atoms				AltConf	Trace
11	XX	48	Total	C	N	O	0	0
			240	144	48	48		
11	YX	48	Total	C	N	O	0	0
			240	144	48	48		
11	ZX	48	Total	C	N	O	0	0
			240	144	48	48		

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.

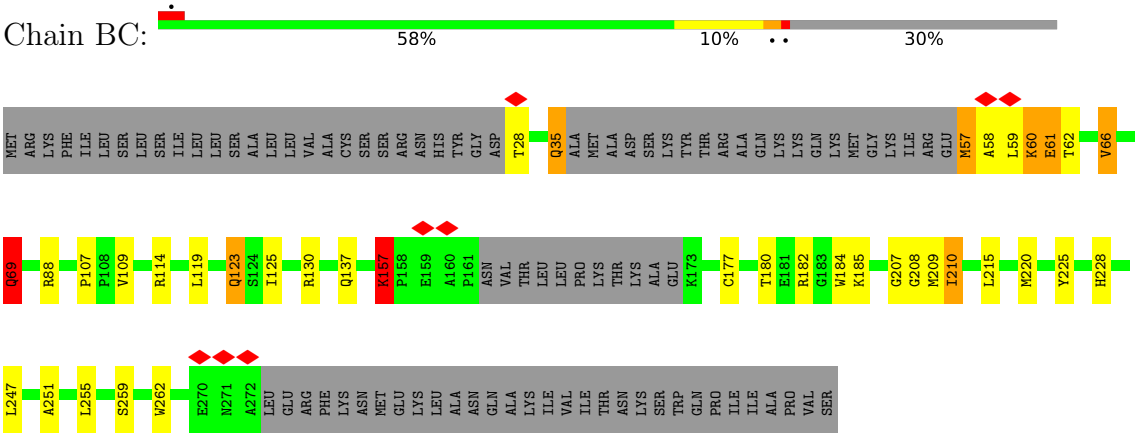
• Molecule 1: DotC

Chain AC:



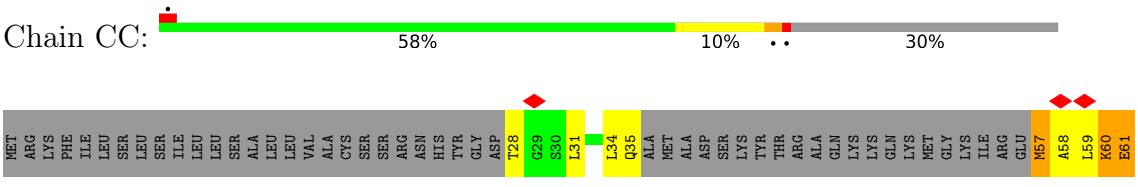
• Molecule 1: DotC

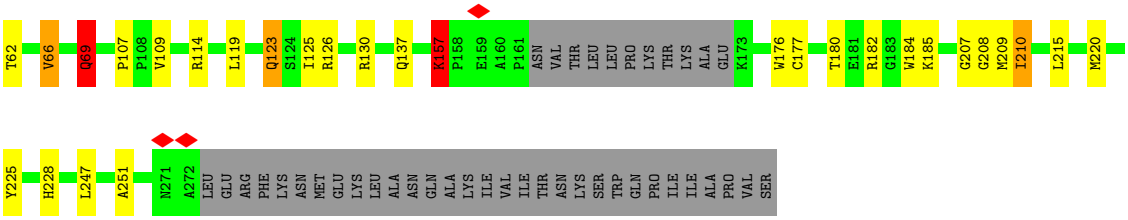
Chain BC:



• Molecule 1: DotC

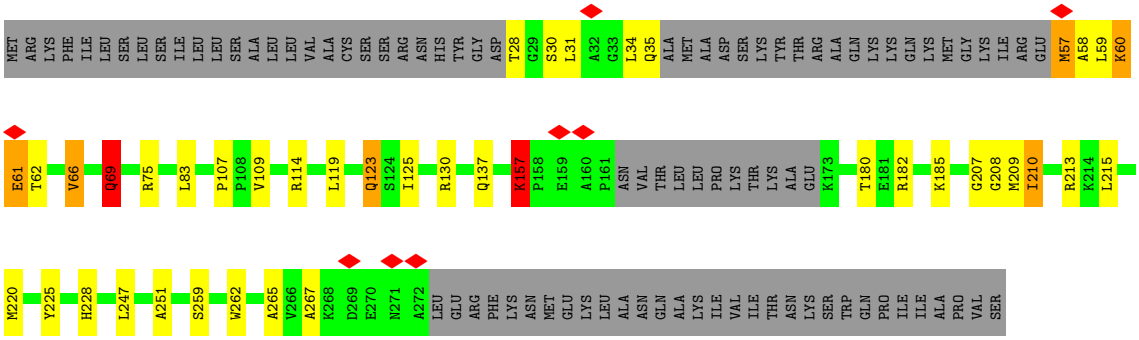
Chain CC:





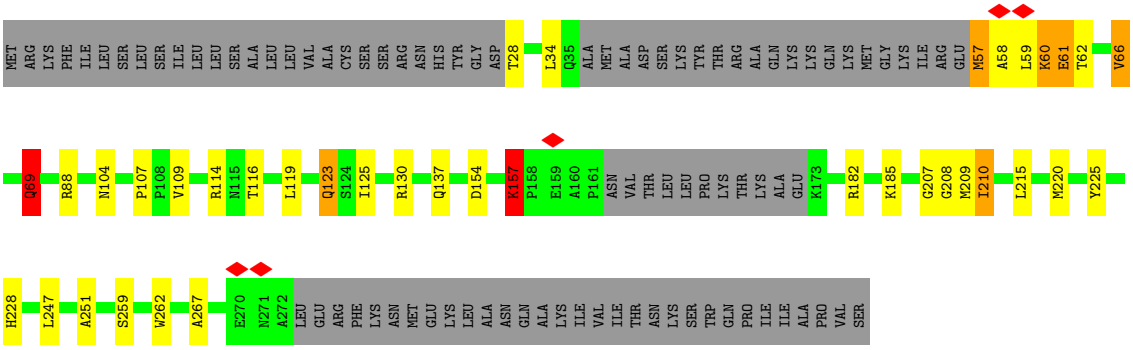
• Molecule 1: DotC

Chain DC:



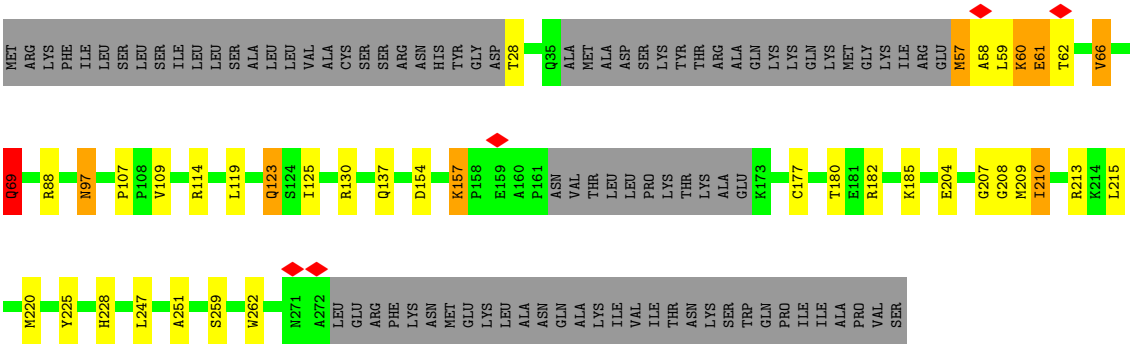
• Molecule 1: DotC

Chain EC:



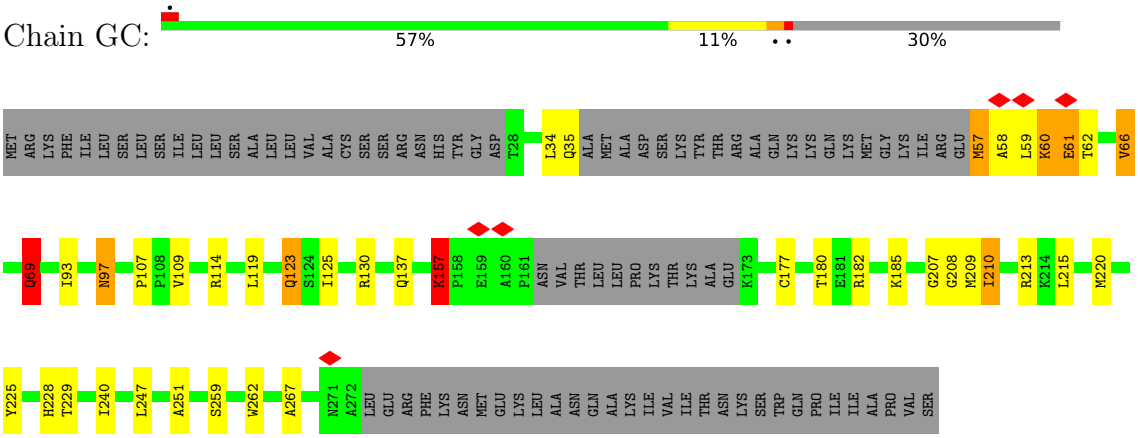
• Molecule 1: DotC

Chain FC:



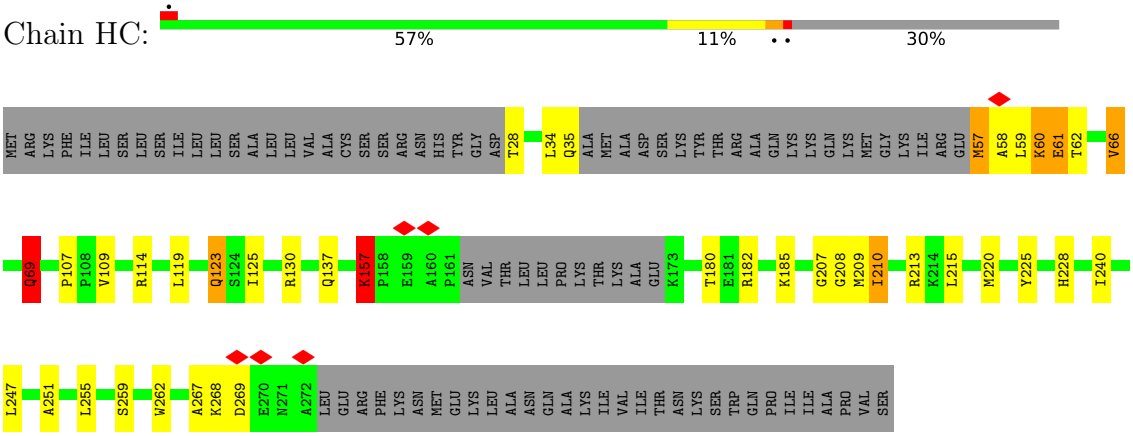
- Molecule 1: DotC

Chain GC:



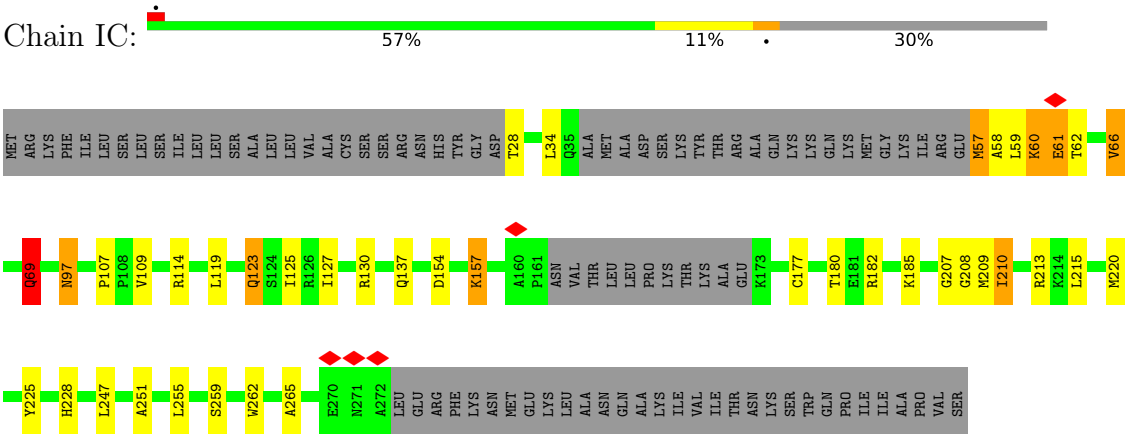
- Molecule 1: DotC

Chain HC:



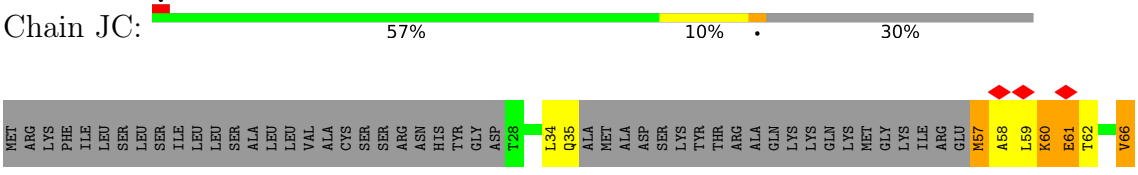
- Molecule 1: DotC

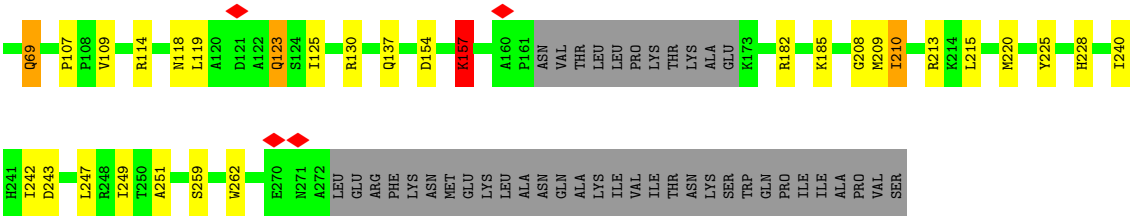
Chain IC:



- Molecule 1: DotC

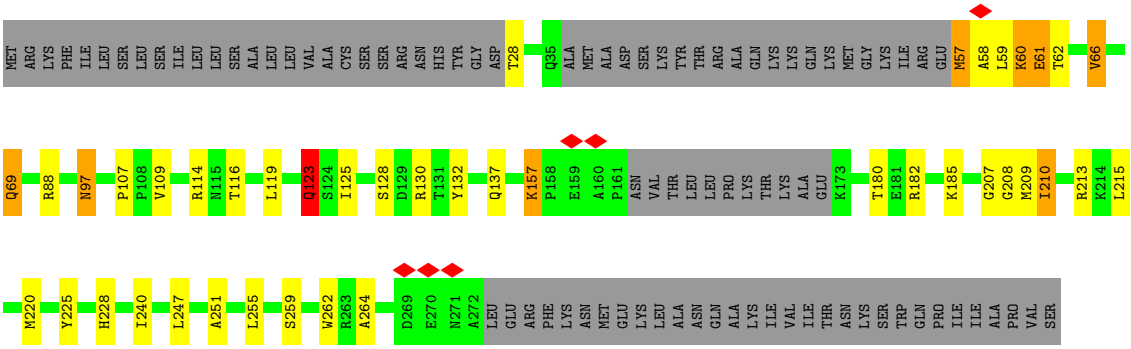
Chain JC:





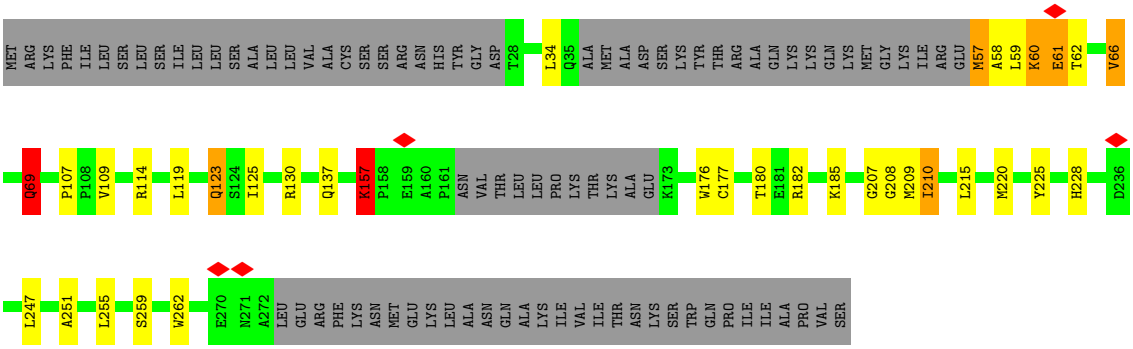
• Molecule 1: DotC

Chain KC:



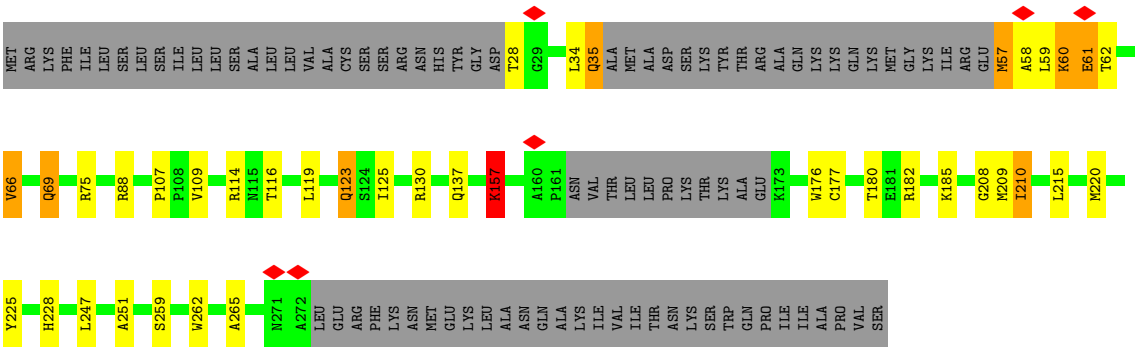
• Molecule 1: DotC

Chain LC:



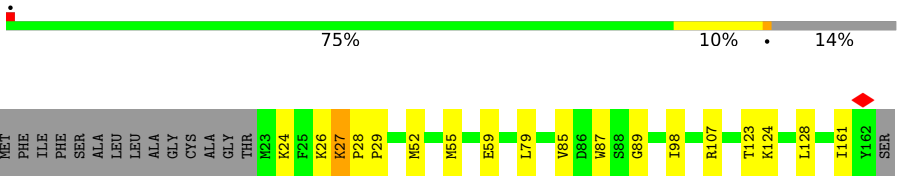
• Molecule 1: DotC

Chain MC:



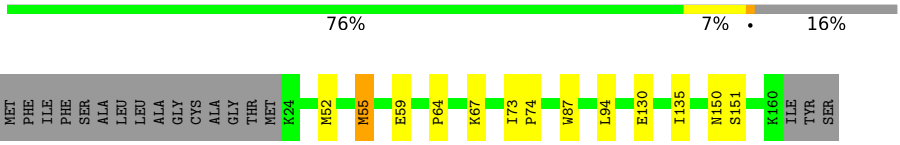
• Molecule 2: DotD

Chain AD:



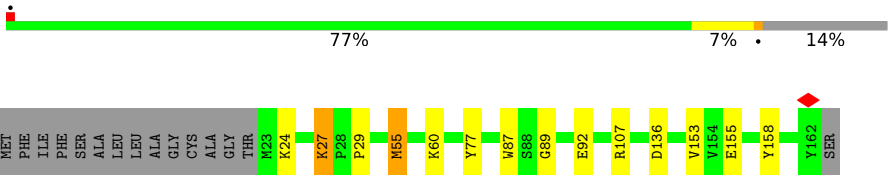
• Molecule 2: DotD

Chain Ad:



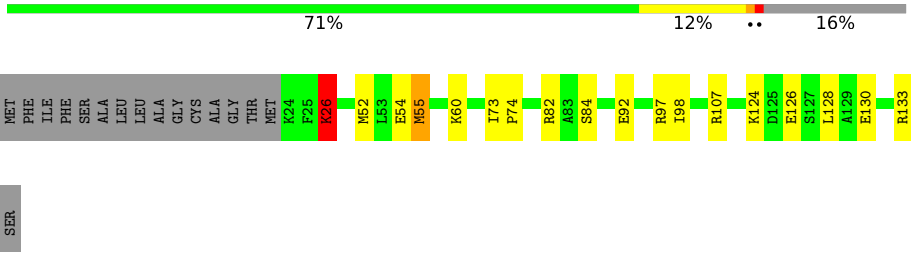
• Molecule 2: DotD

Chain BD:



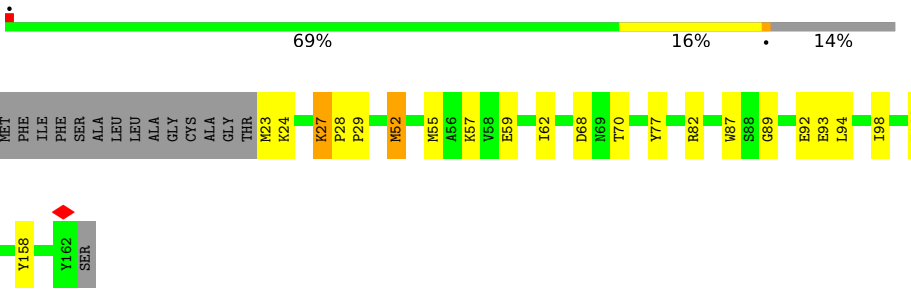
• Molecule 2: DotD

Chain Bd:



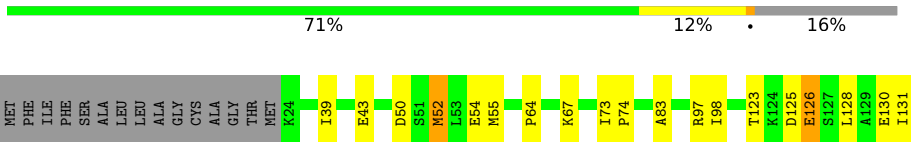
• Molecule 2: DotD

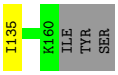
Chain CD:



• Molecule 2: DotD

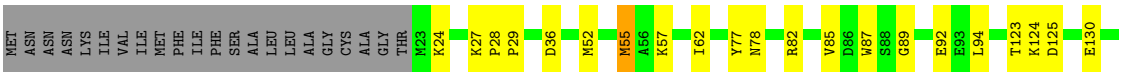
Chain Cd:

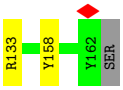




• Molecule 2: DotD

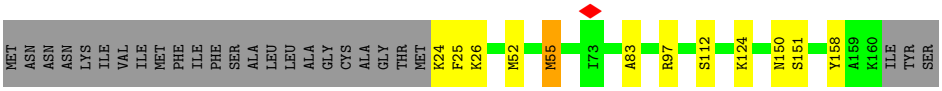
Chain DD:





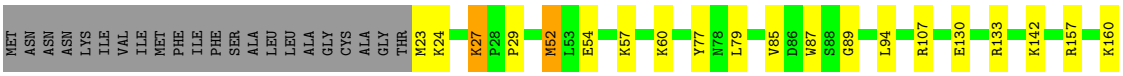
• Molecule 2: DotD

Chain Dd:



• Molecule 2: DotD

Chain ED:

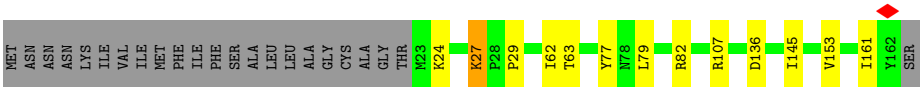


• Molecule 2: DotD

Chain Ed:

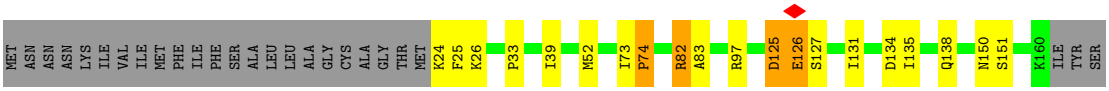
• Molecule 2: DotD

Chain FD:



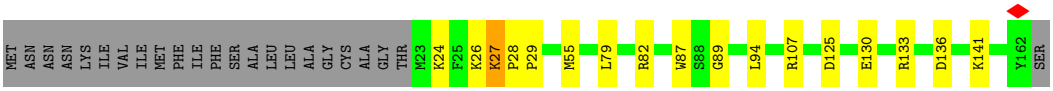
• Molecule 2: DotD

Chain Fd:



• Molecule 2: DotD

Chain GD:

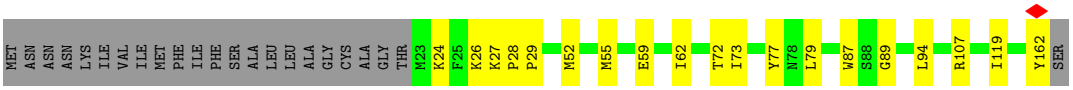


• Molecule 2: DotD

Chain Gd:

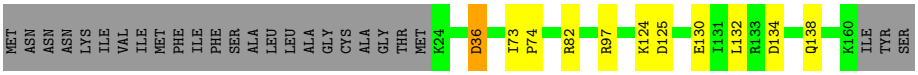
• Molecule 2: DotD

Chain HD:



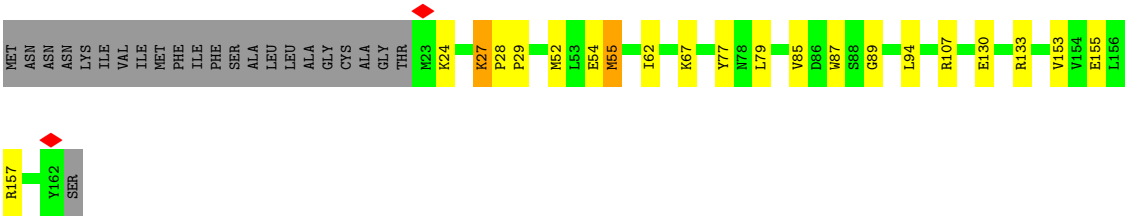
• Molecule 2: DotD

Chain Hd:



• Molecule 2: DotD

Chain ID:



• Molecule 2: DotD

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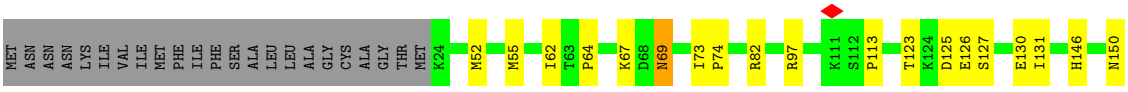
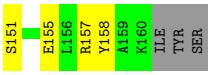
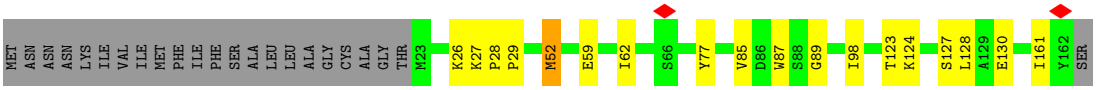




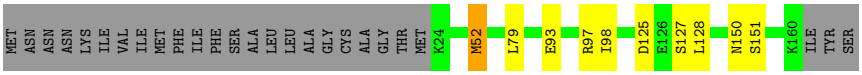
• Molecule 2: DotD

Chain JD:



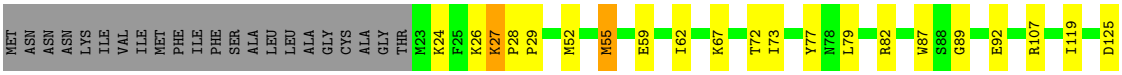
• Molecule 2: DotD

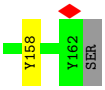
Chain Jd:



• Molecule 2: DotD

Chain KD:





• Molecule 2: DotD

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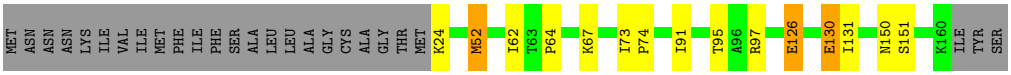

• Molecule 2: DotD

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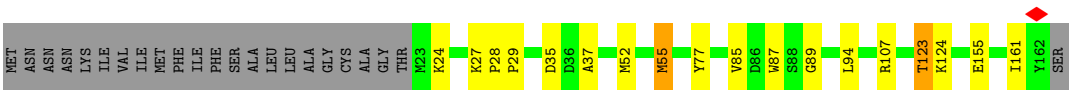
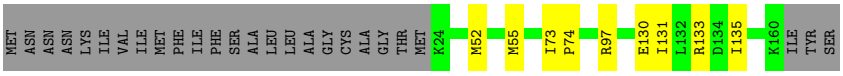


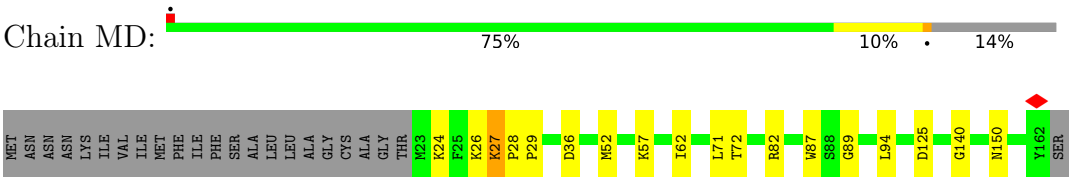
• Molecule 2: DotD

Chain Ld:



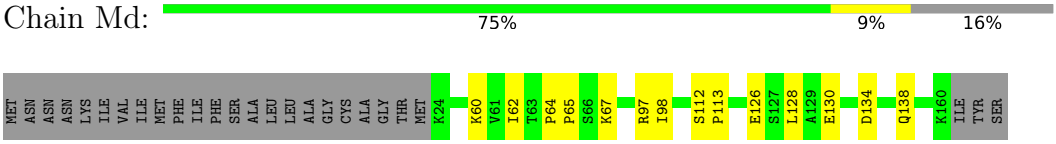
• Molecule 2: DotD

Chain MD:



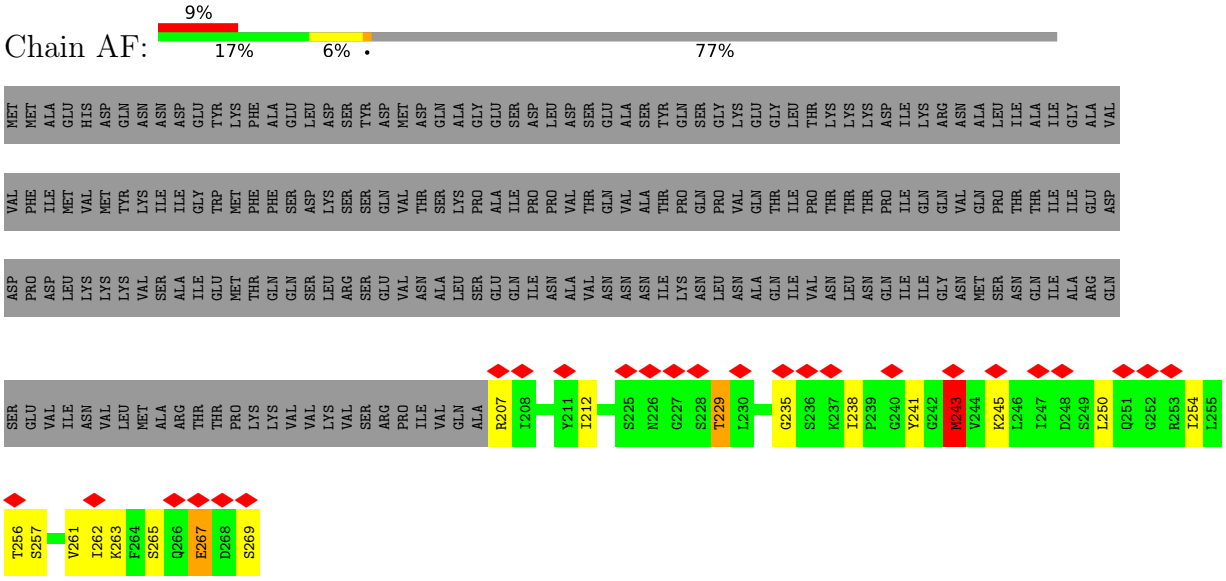
• Molecule 2: DotD

Chain Md:



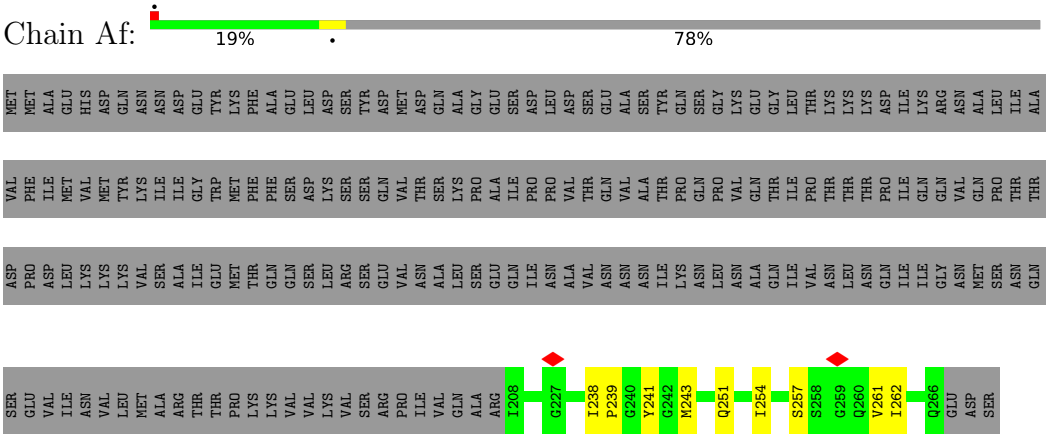
• Molecule 3: DotF

Chain AF:



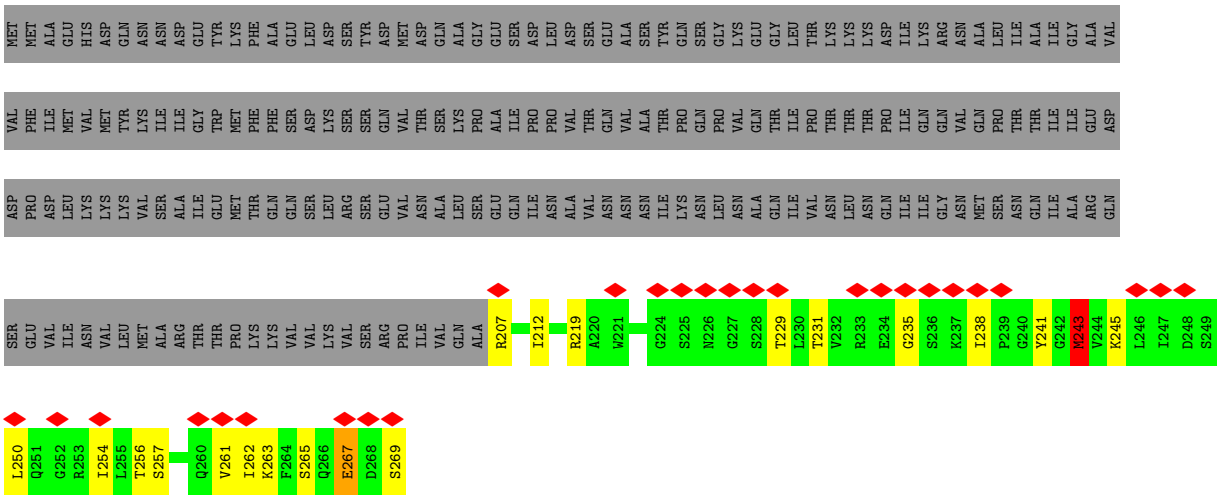
• Molecule 3: DotF

Chain Af:



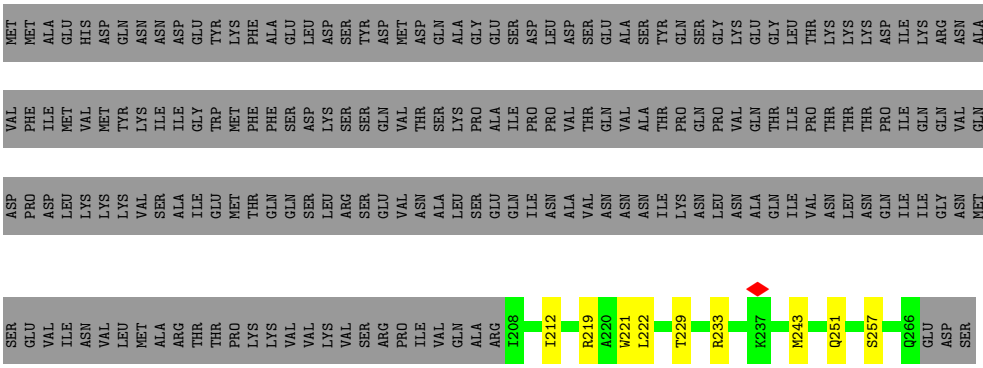
• Molecule 3: DotF

Chain BF:



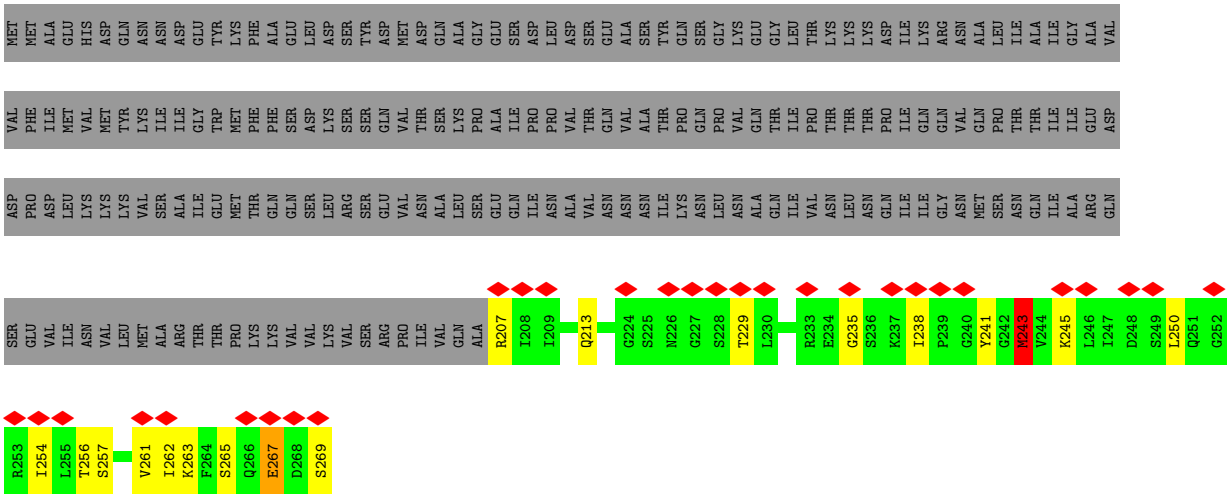
• Molecule 3: DotF

Chain Bf:



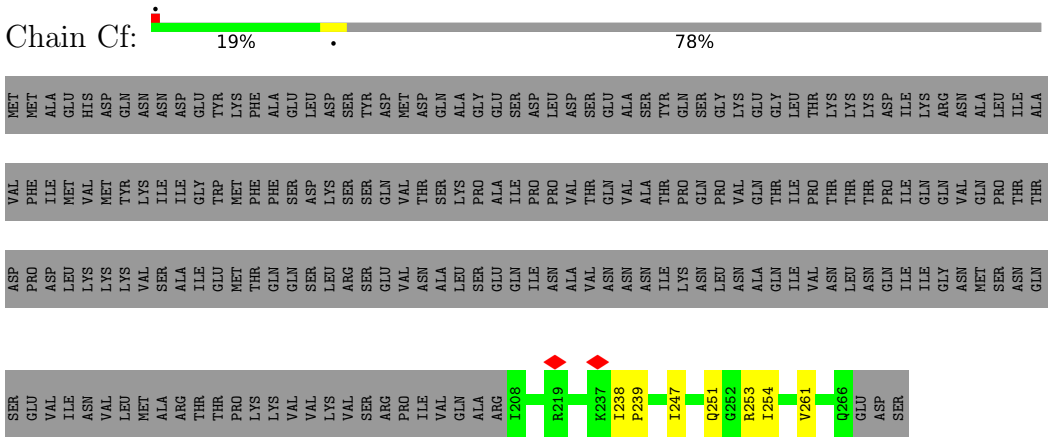
• Molecule 3: DotF

Chain CF:



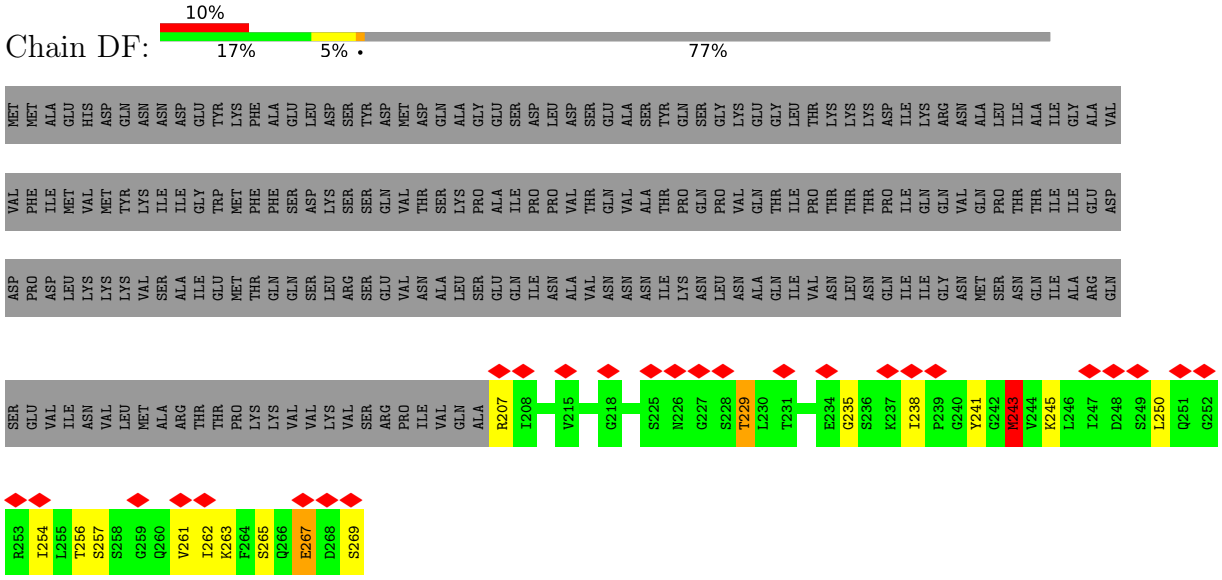
• Molecule 3: DotF

Chain Cf:



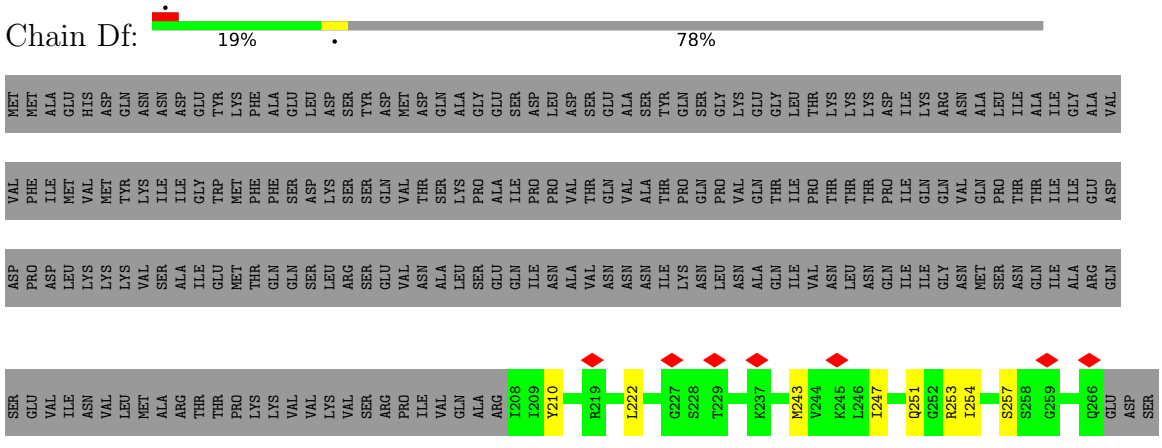
• Molecule 3: DotF

Chain DF:



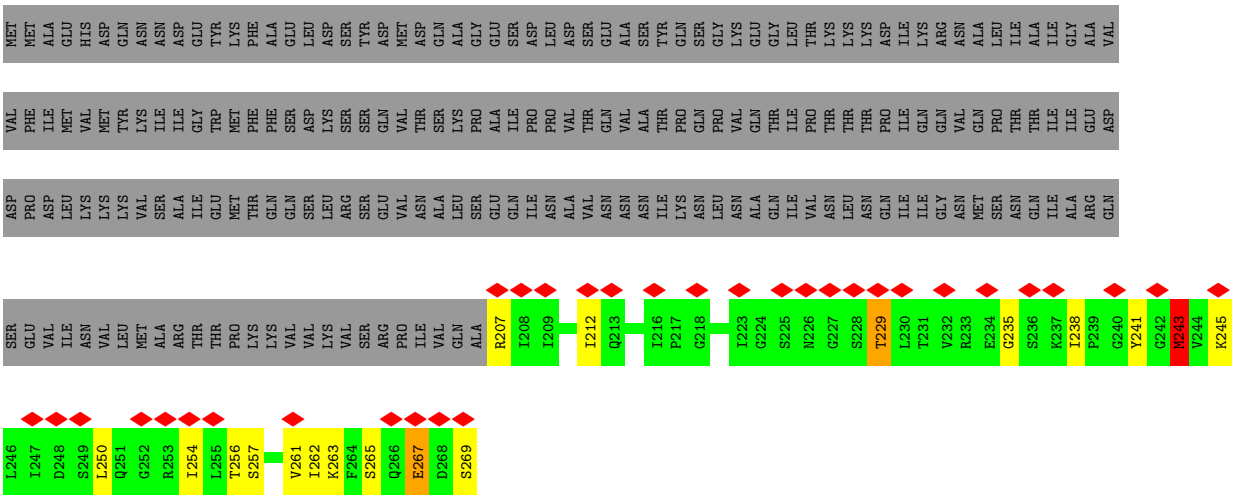
• Molecule 3: DotF

Chain Df:



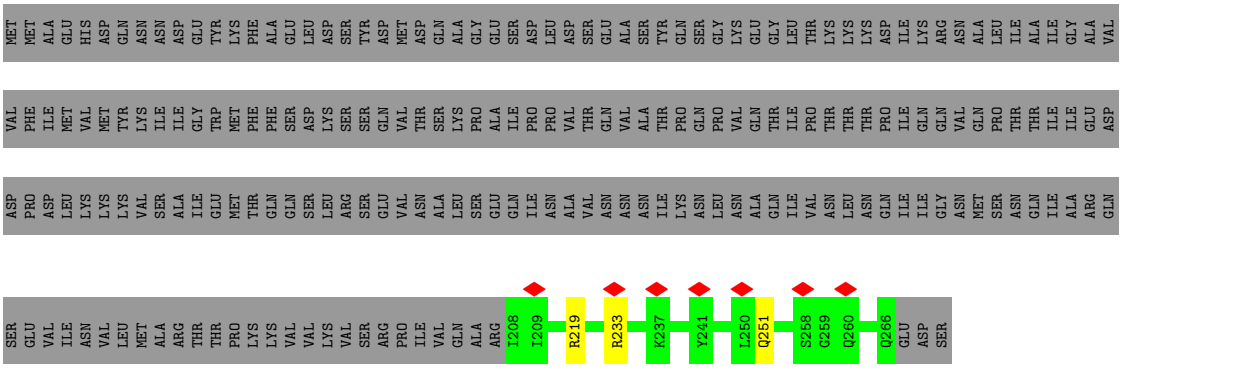
• Molecule 3: DotF

Chain EF:



- Molecule 3: DotF

Chain Ef:

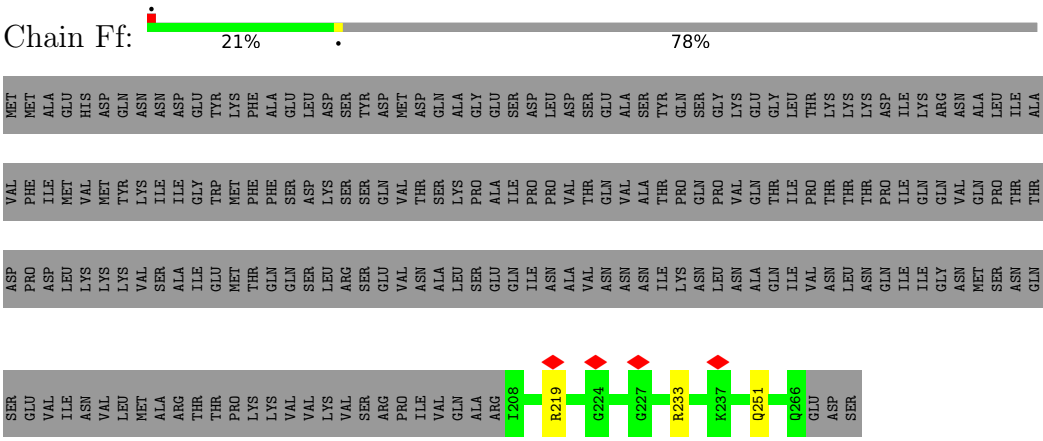


- Molecule 3: DotF

Chain FF:

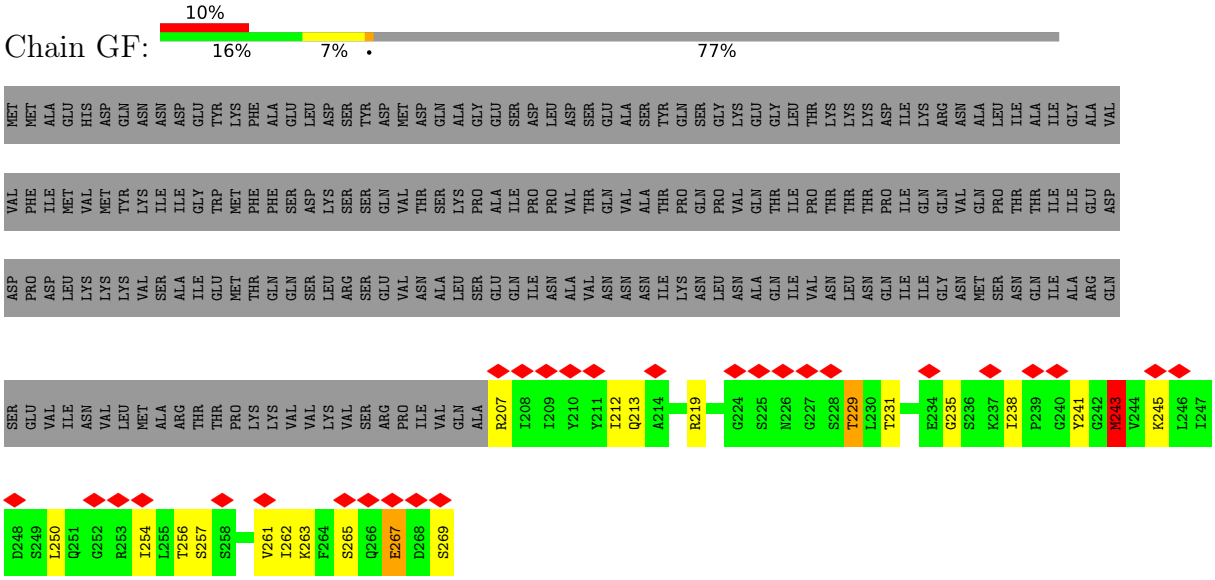
- Molecule 3: DotF

Chain Ff:



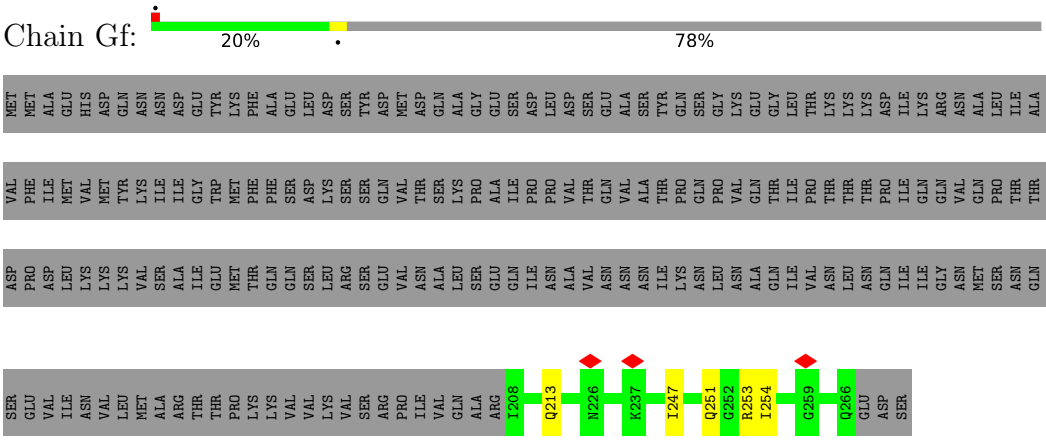
• Molecule 3: DotF

Chain GF:



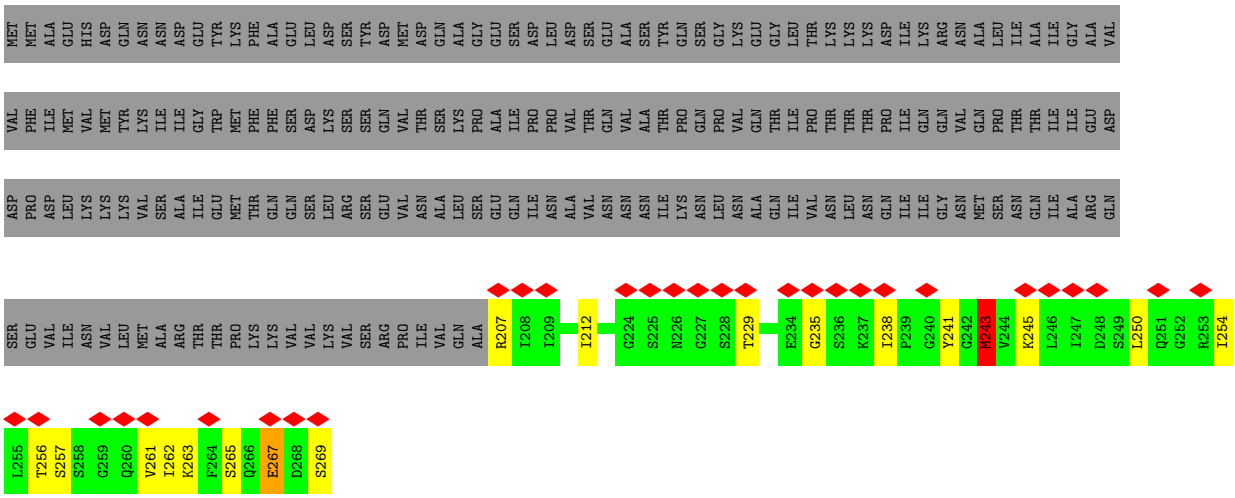
• Molecule 3: DotF

Chain Gf:



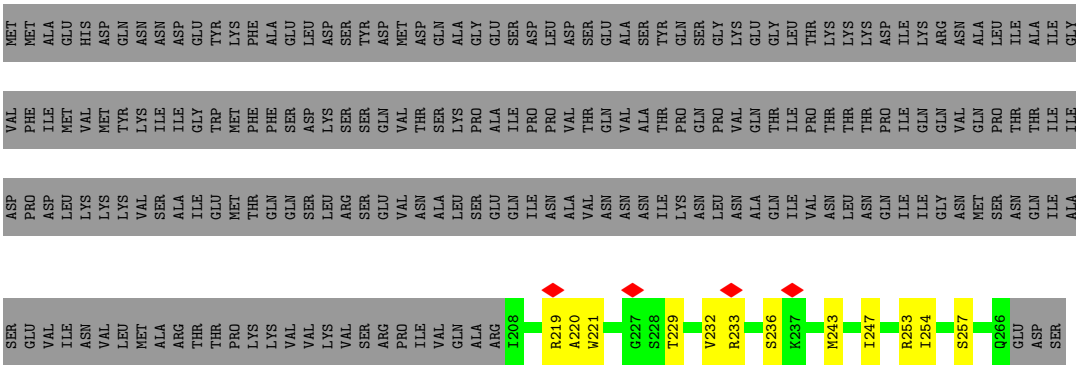
• Molecule 3: DotF

Chain HF:



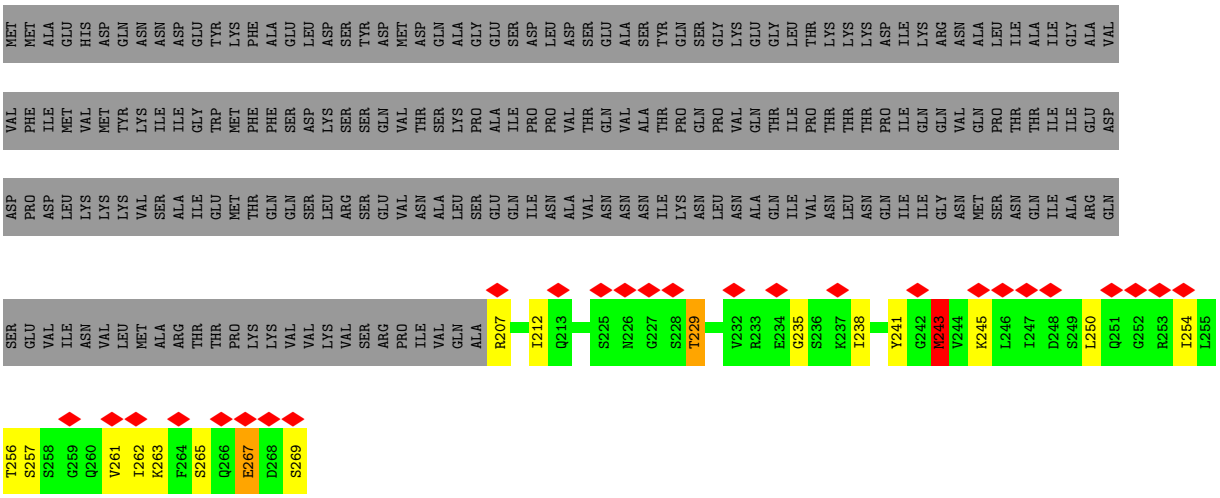
- Molecule 3: DotF

Chain Hf:



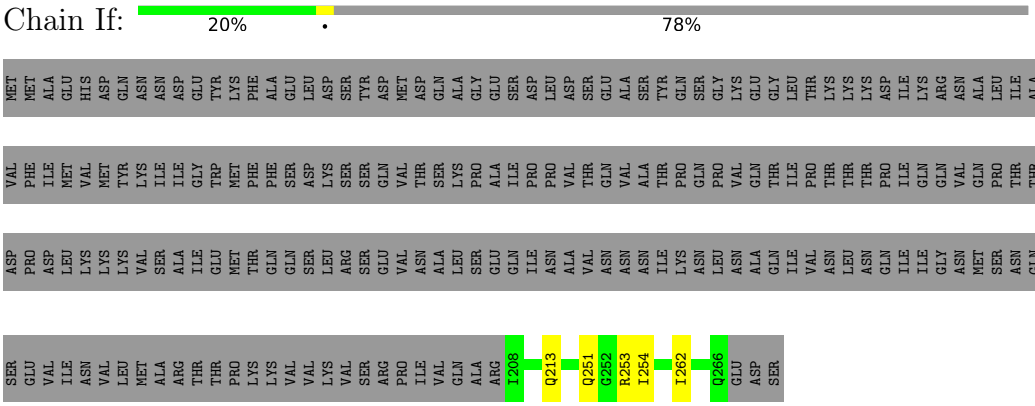
- Molecule 3: DotF

Chain IF:



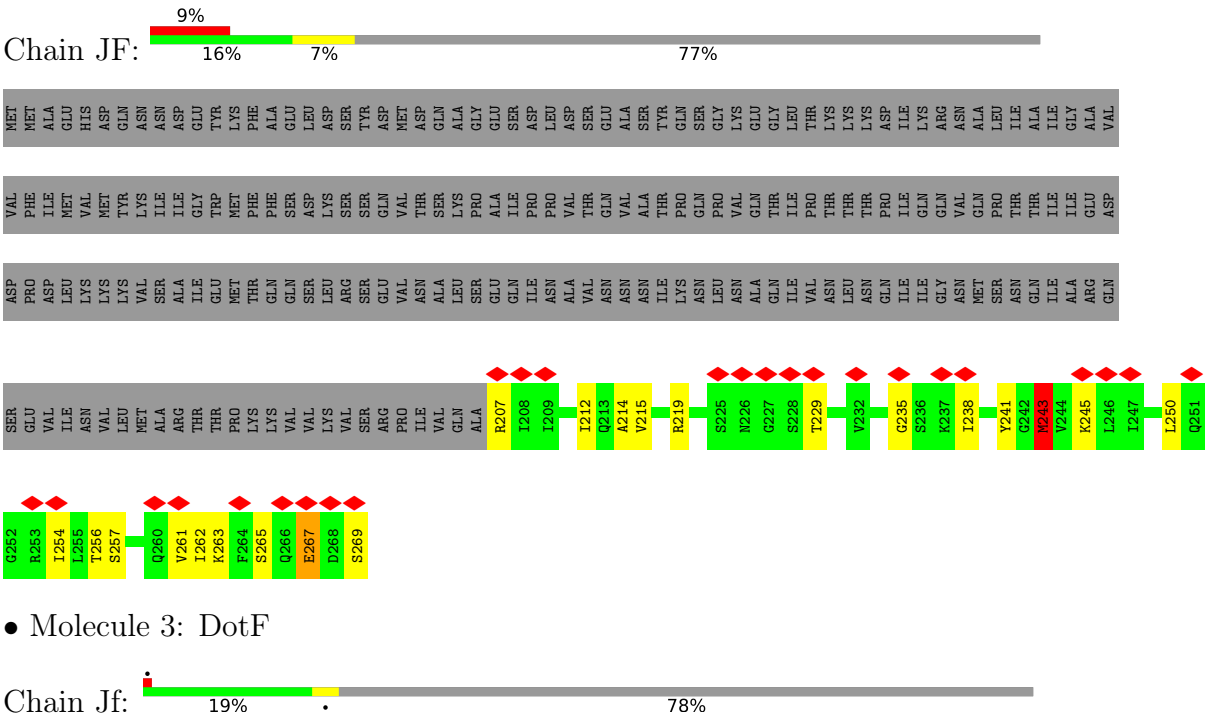
- Molecule 3: DotF

Chain If:



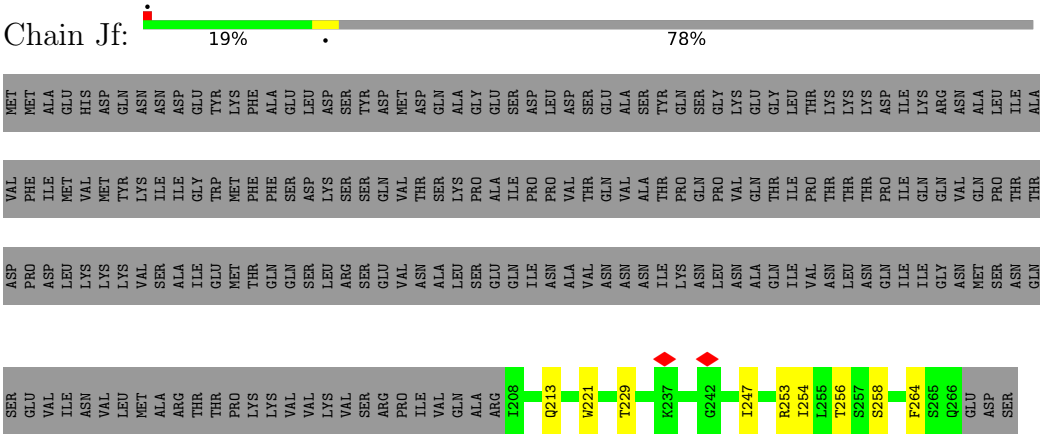
• Molecule 3: DotF

Chain JF:



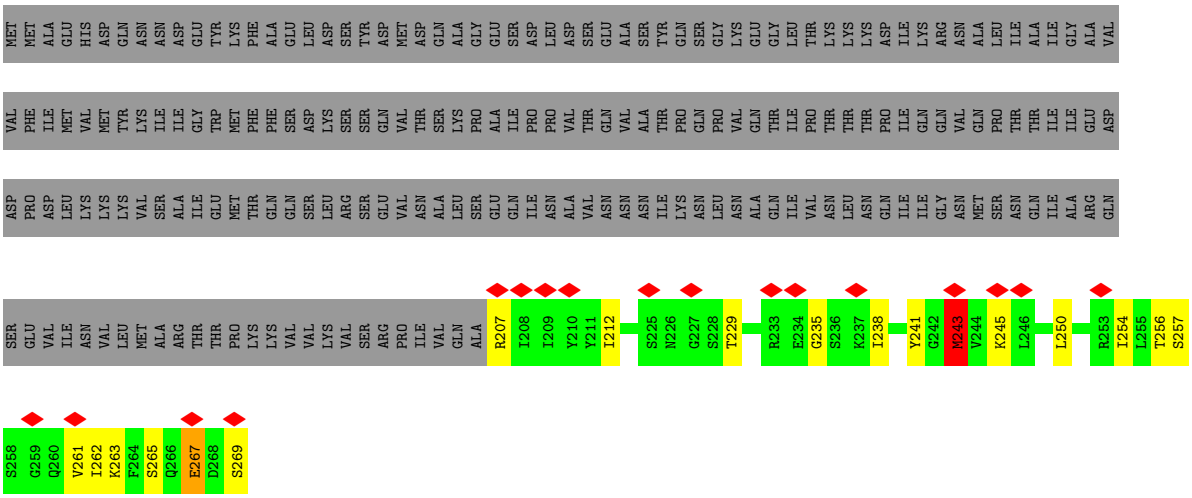
• Molecule 3: DotF

Chain Jf:



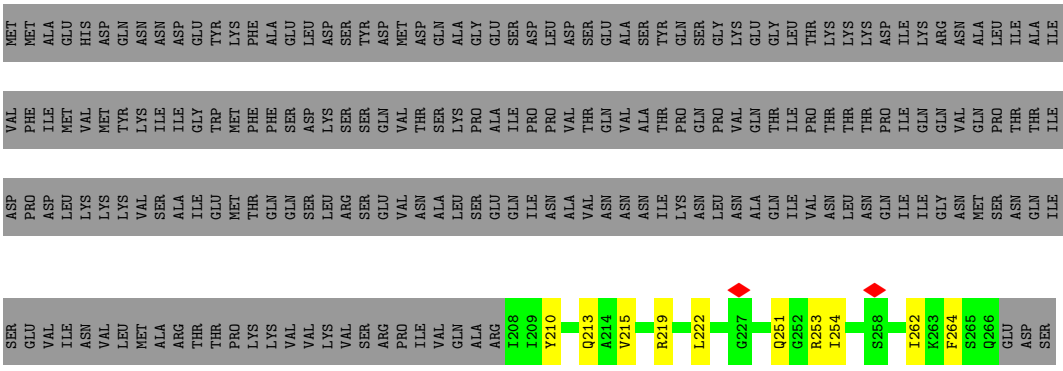
• Molecule 3: DotF

Chain KF:



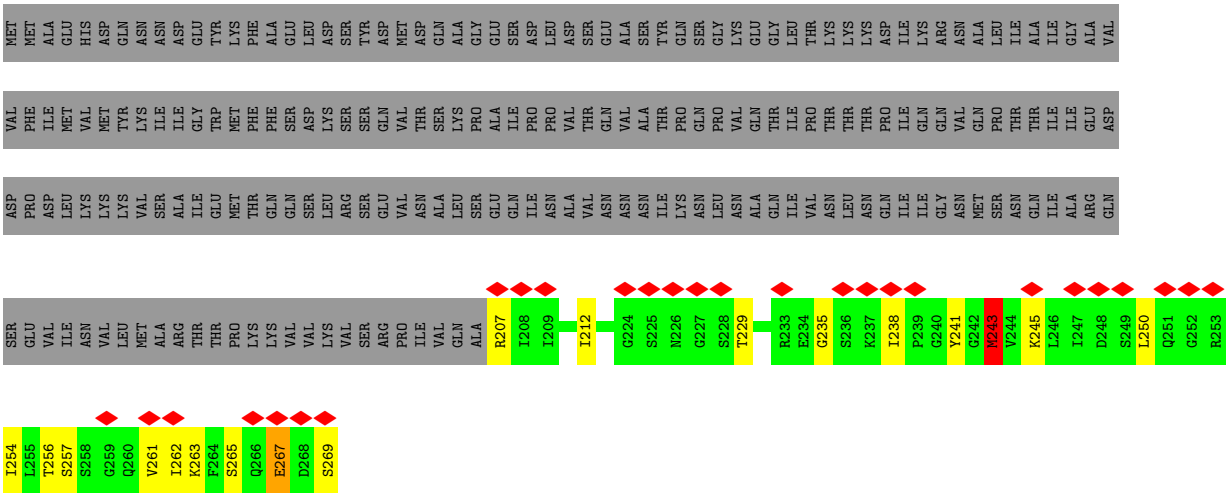
• Molecule 3: DotF

Chain Kf:



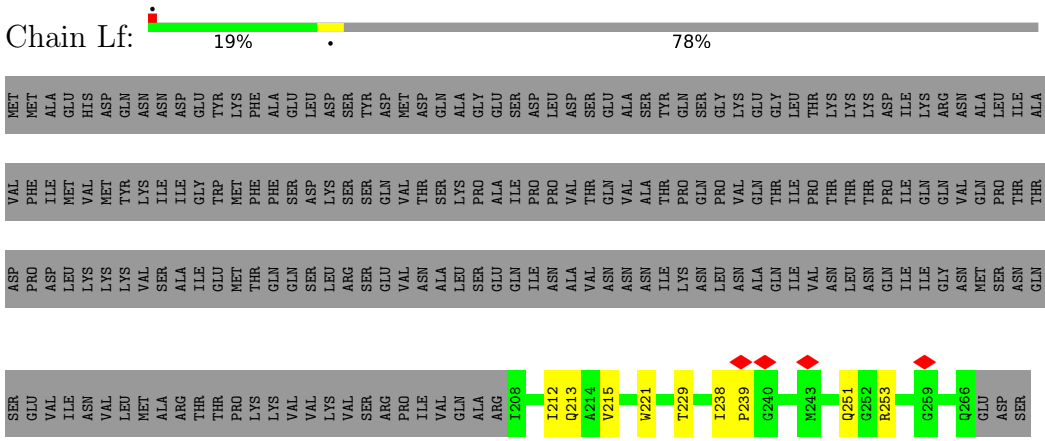
• Molecule 3: DotF

Chain LF:



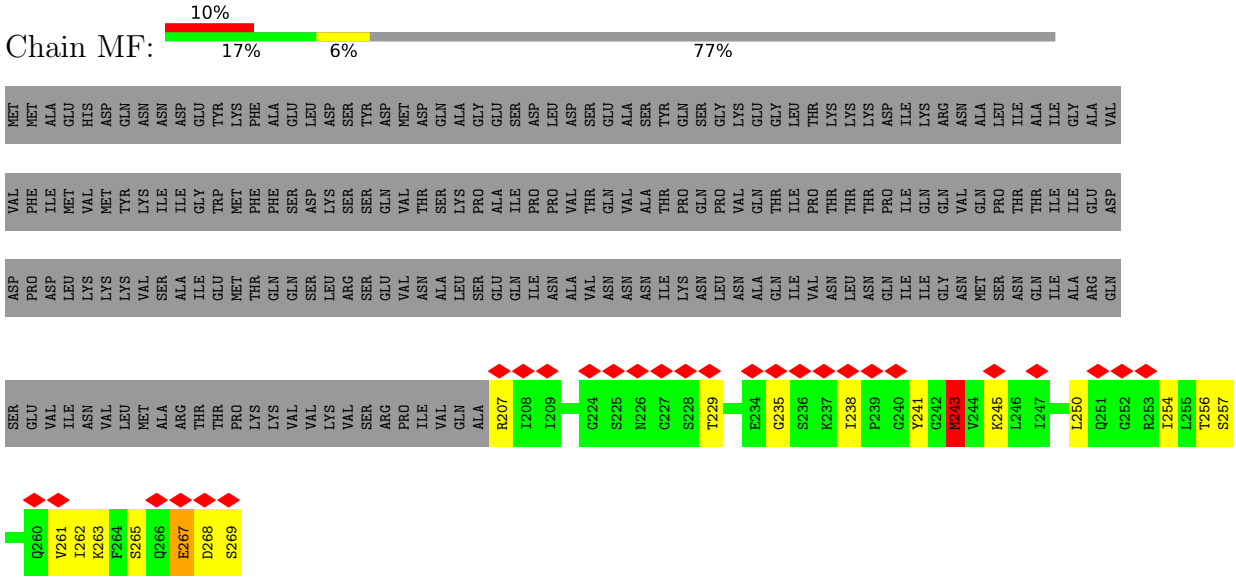
• Molecule 3: DotF

Chain Lf:



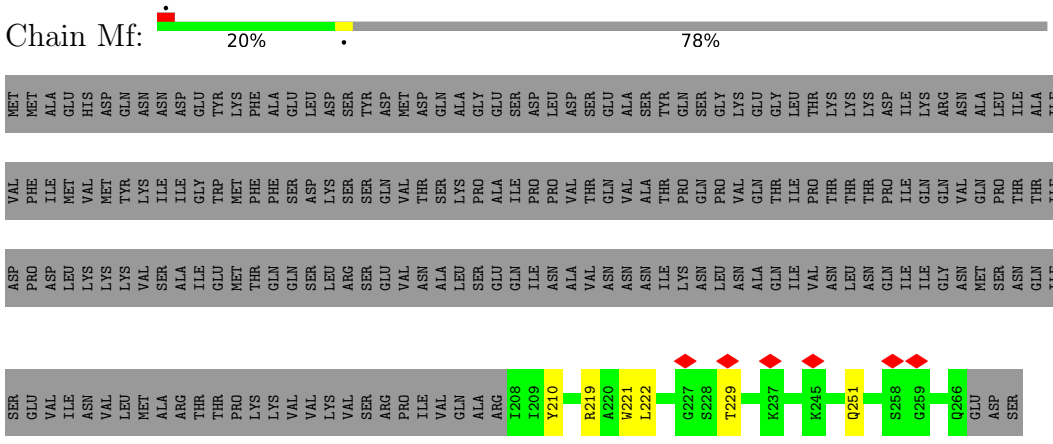
• Molecule 3: DotF

Chain MF:



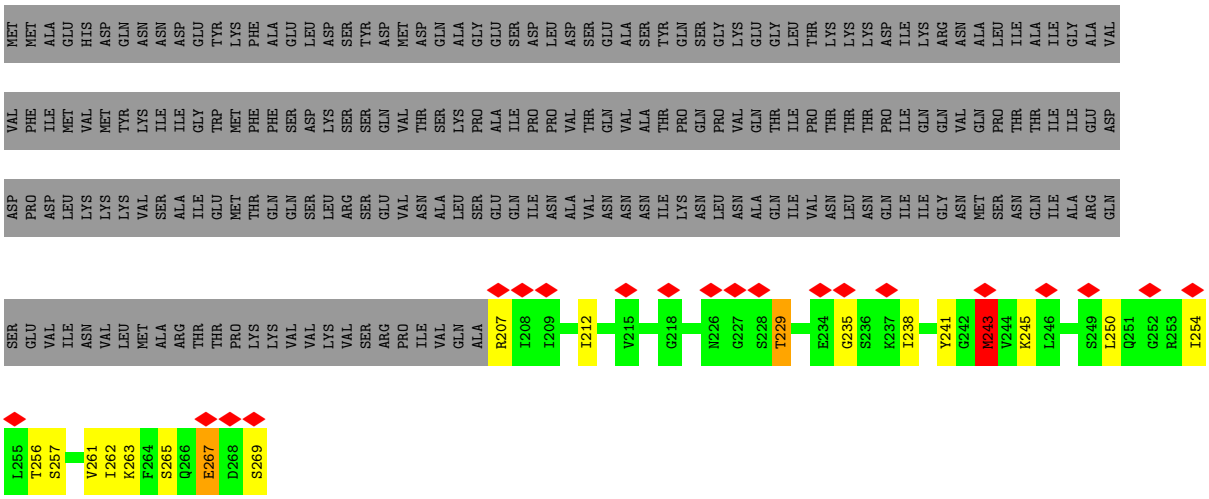
• Molecule 3: DotF

Chain Mf:



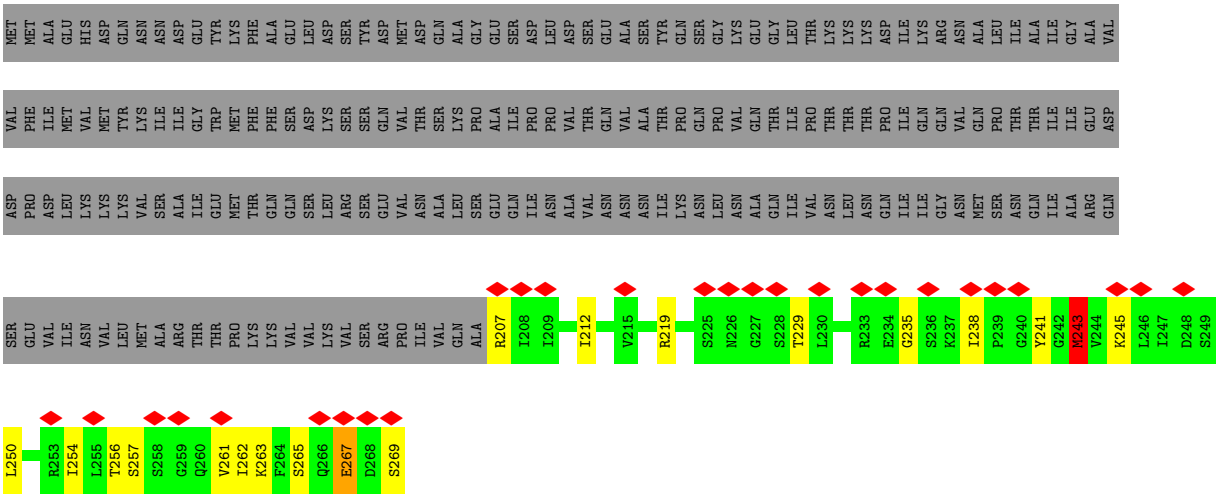
• Molecule 3: DotF

Chain VF:



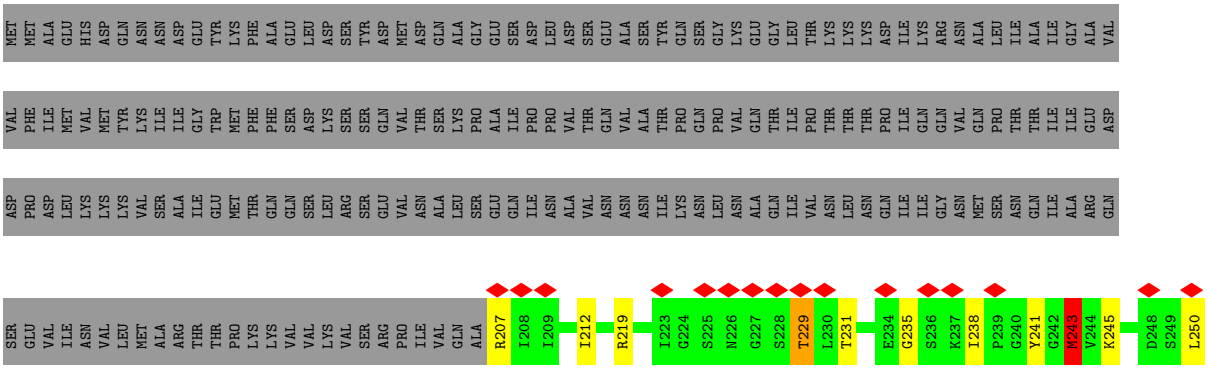
• Molecule 3: DotF

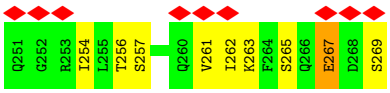
Chain WF:



• Molecule 3: DotF

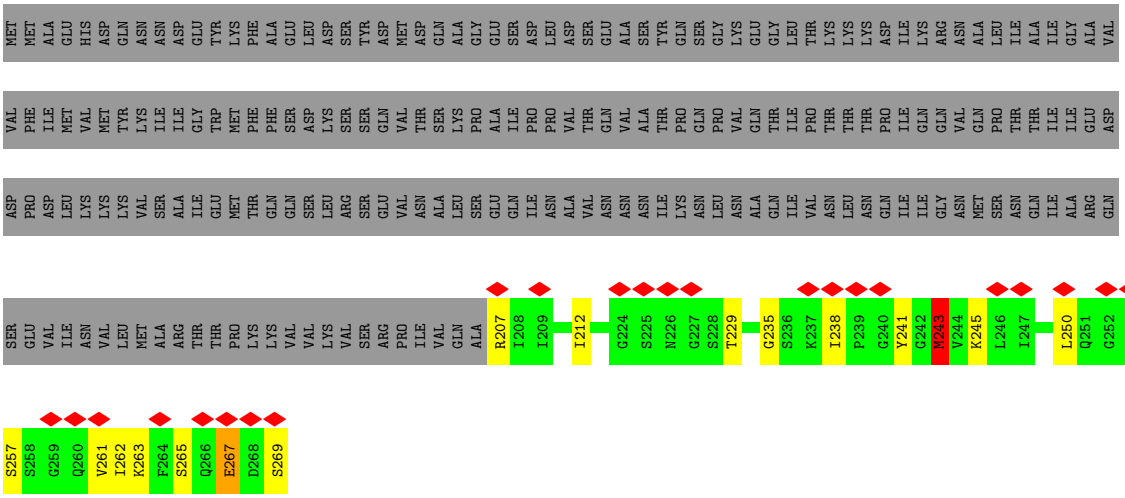
Chain XF:





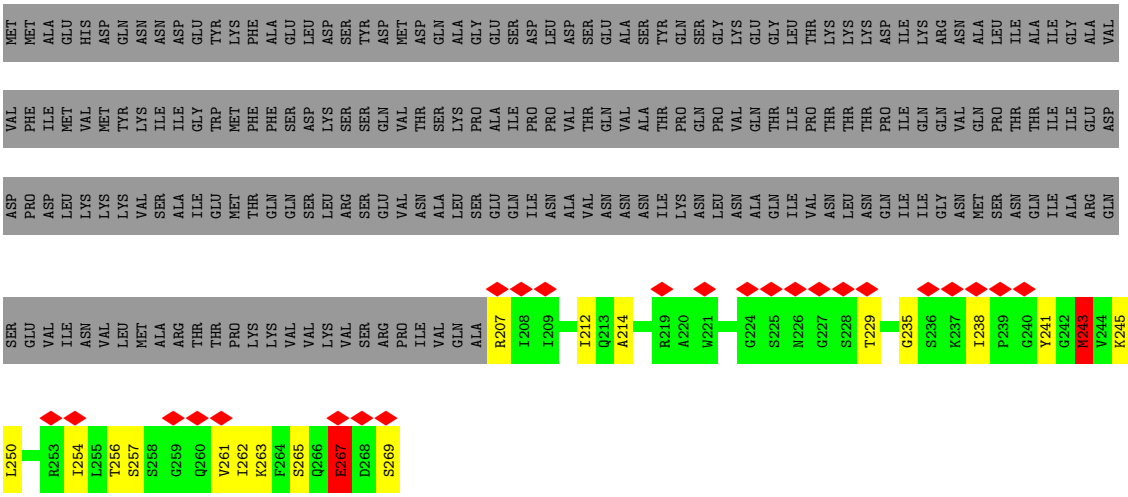
• Molecule 3: DotF

Chain YF:



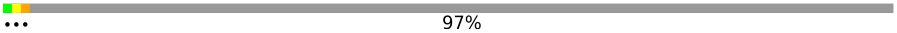
• Molecule 3: DotF

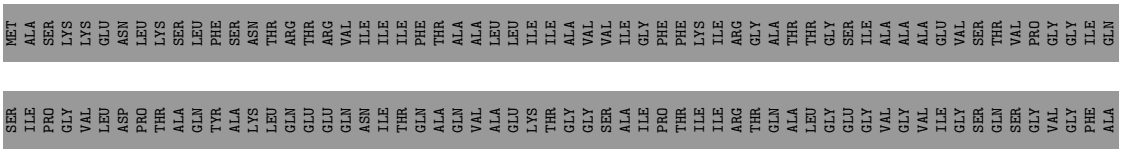
Chain ZF:

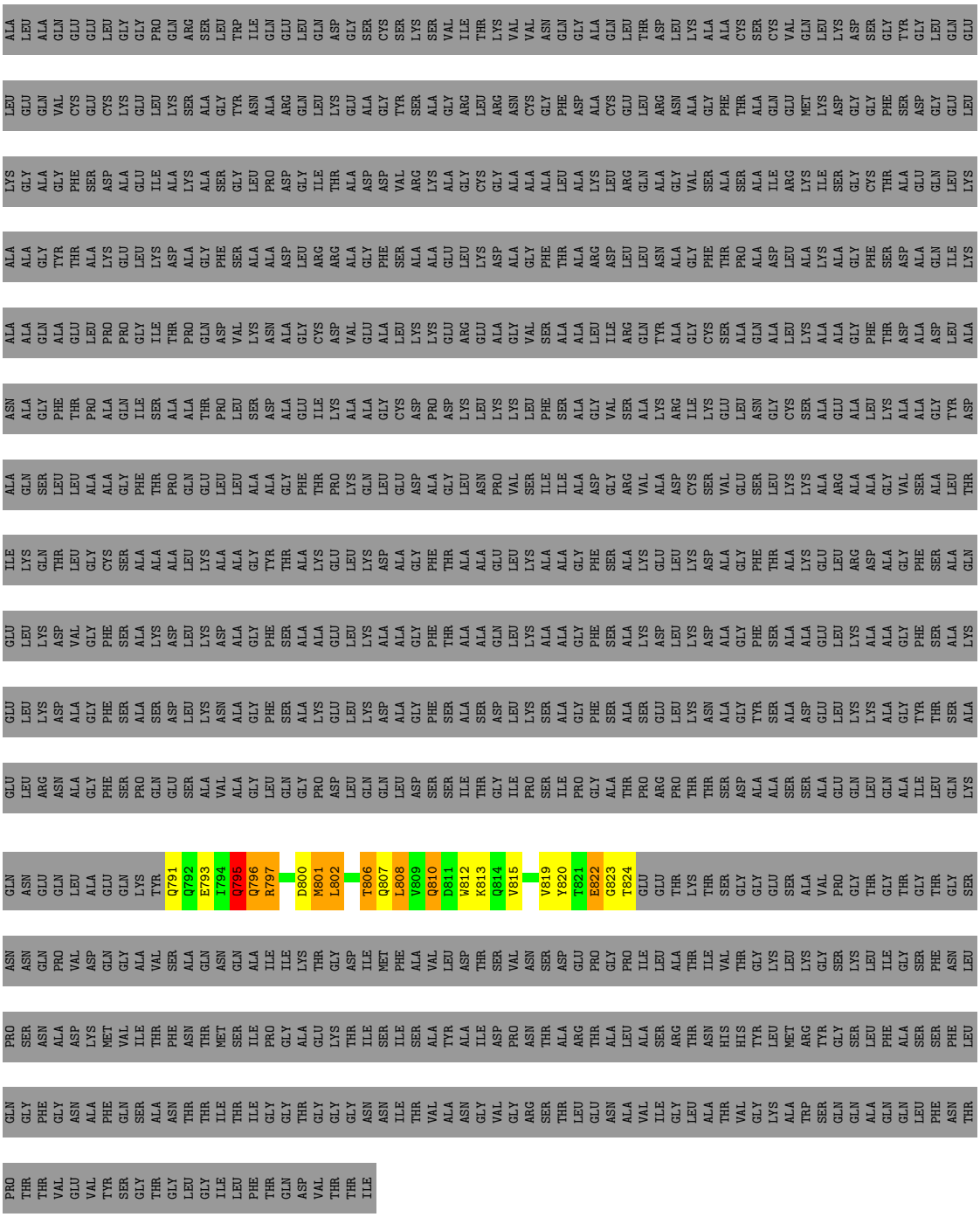


• Molecule 4: IcmE protein

Chain AG:

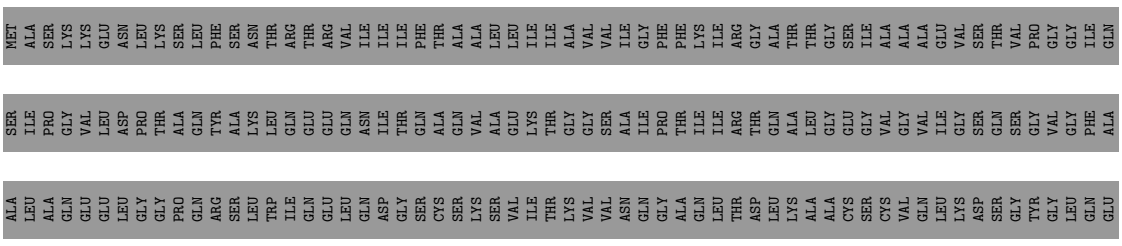


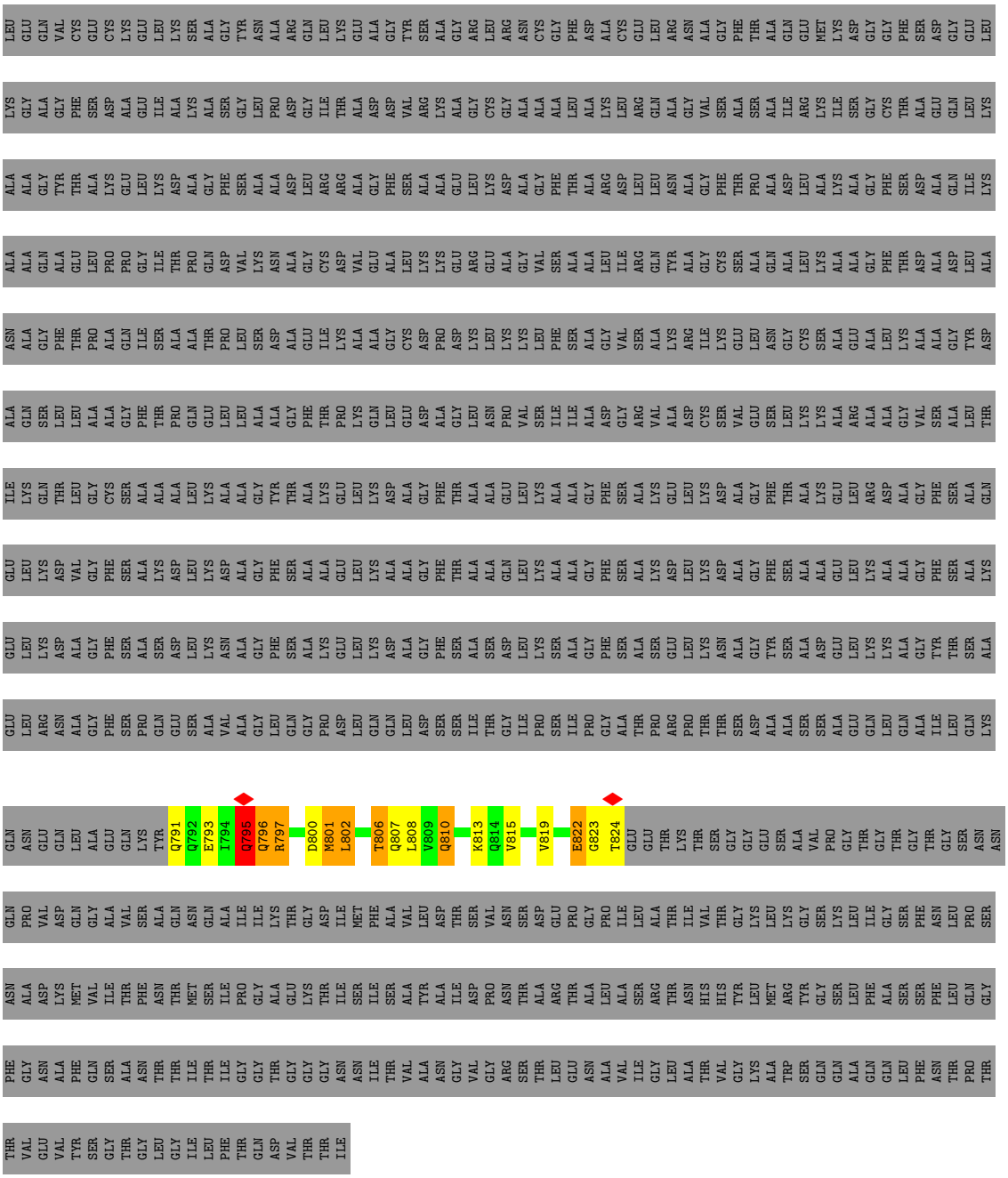




• Molecule 4: IcmE protein

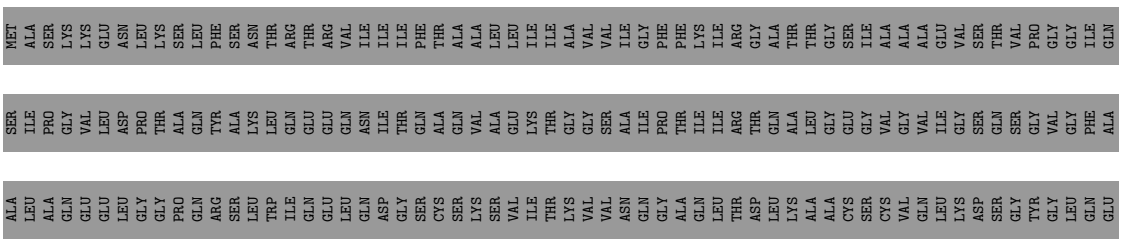
Chain BG: 97%

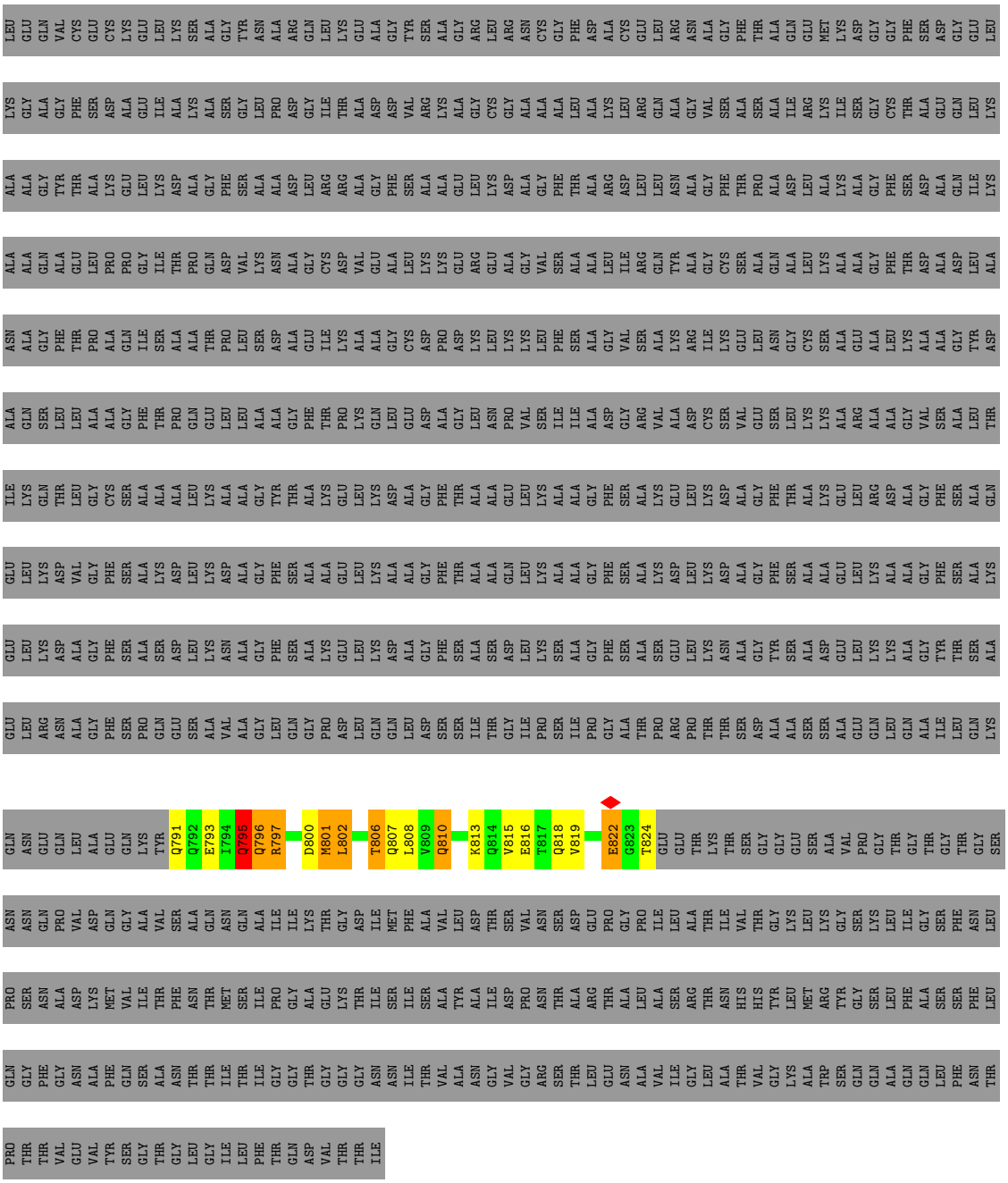




• Molecule 4: IcmE protein

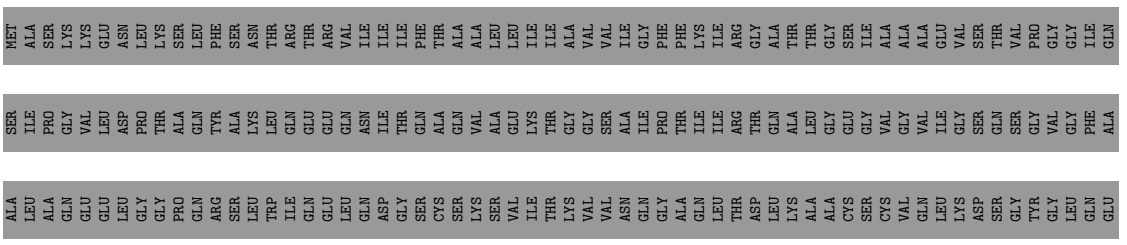
Chain CG: 97%

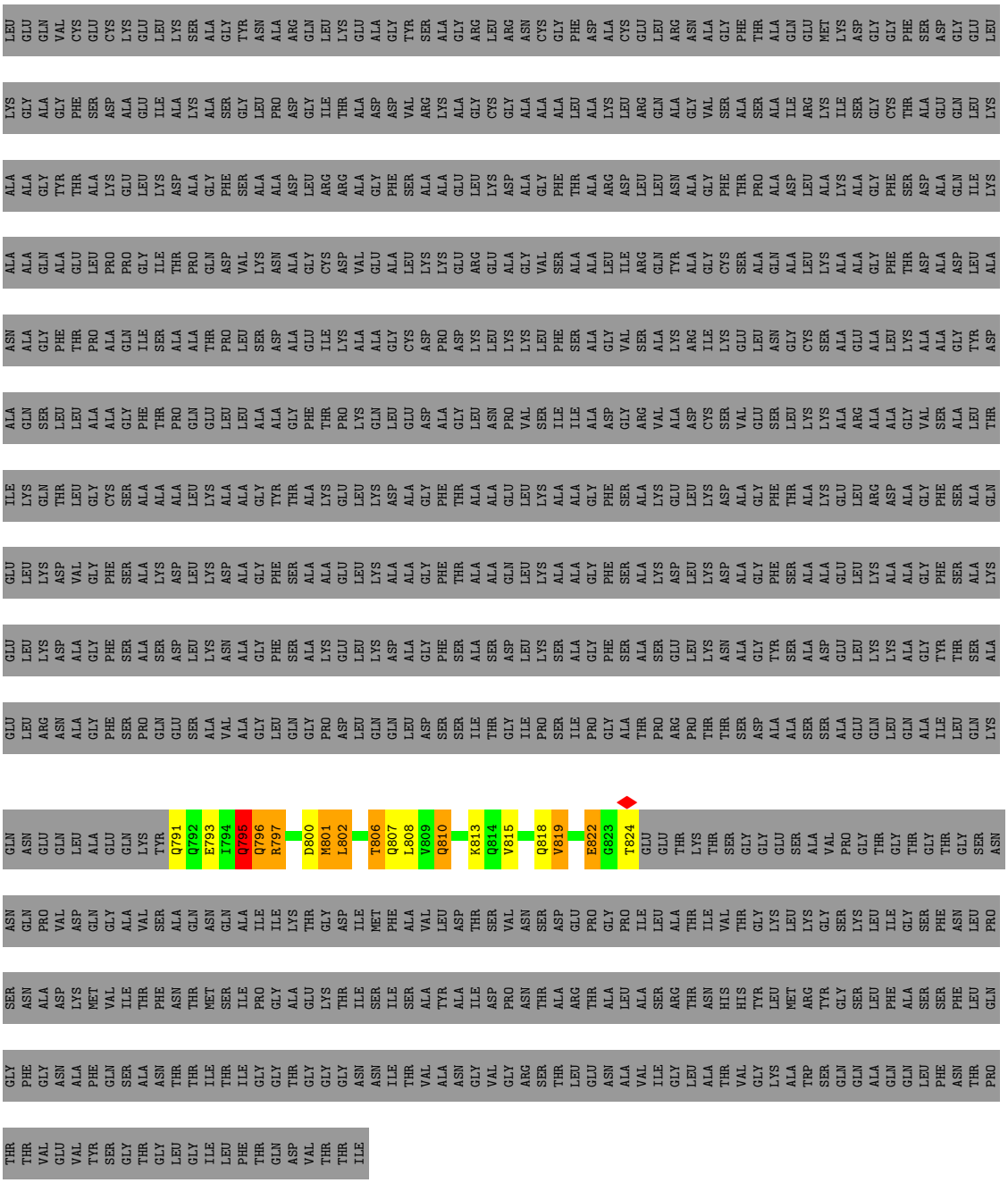




• Molecule 4: IcmE protein

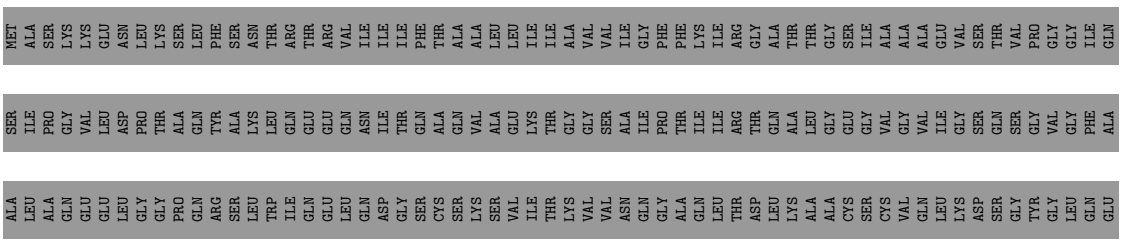
Chain DG: 97%

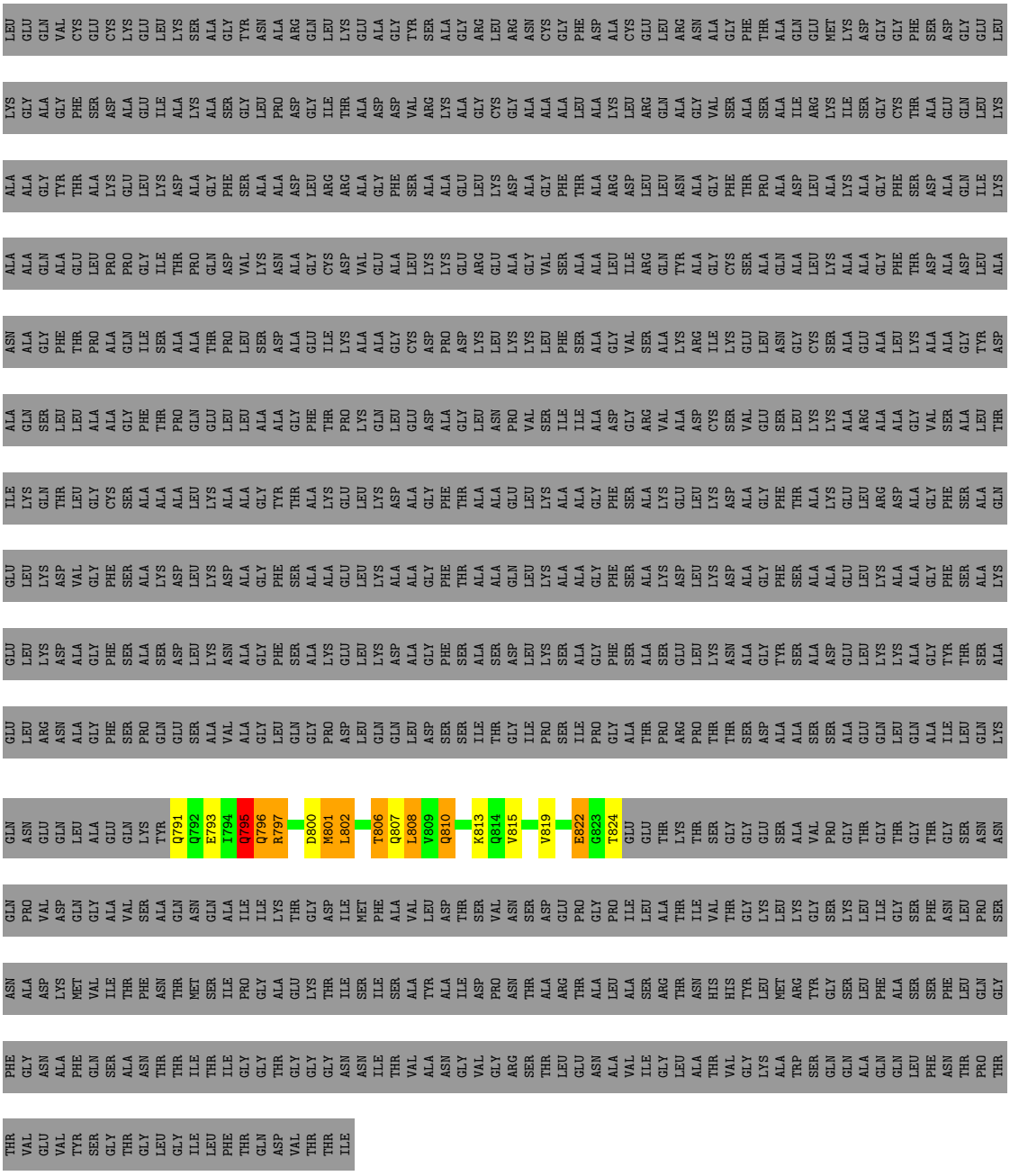




• Molecule 4: IcmE protein

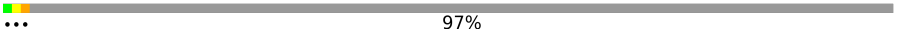
Chain EG: 97%

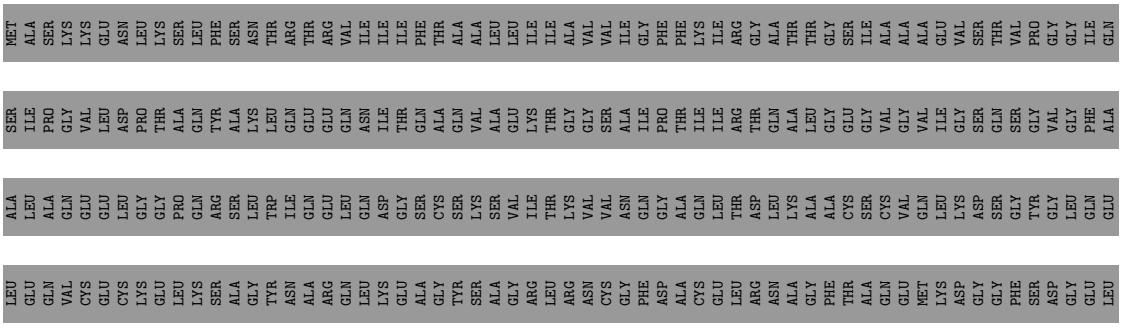


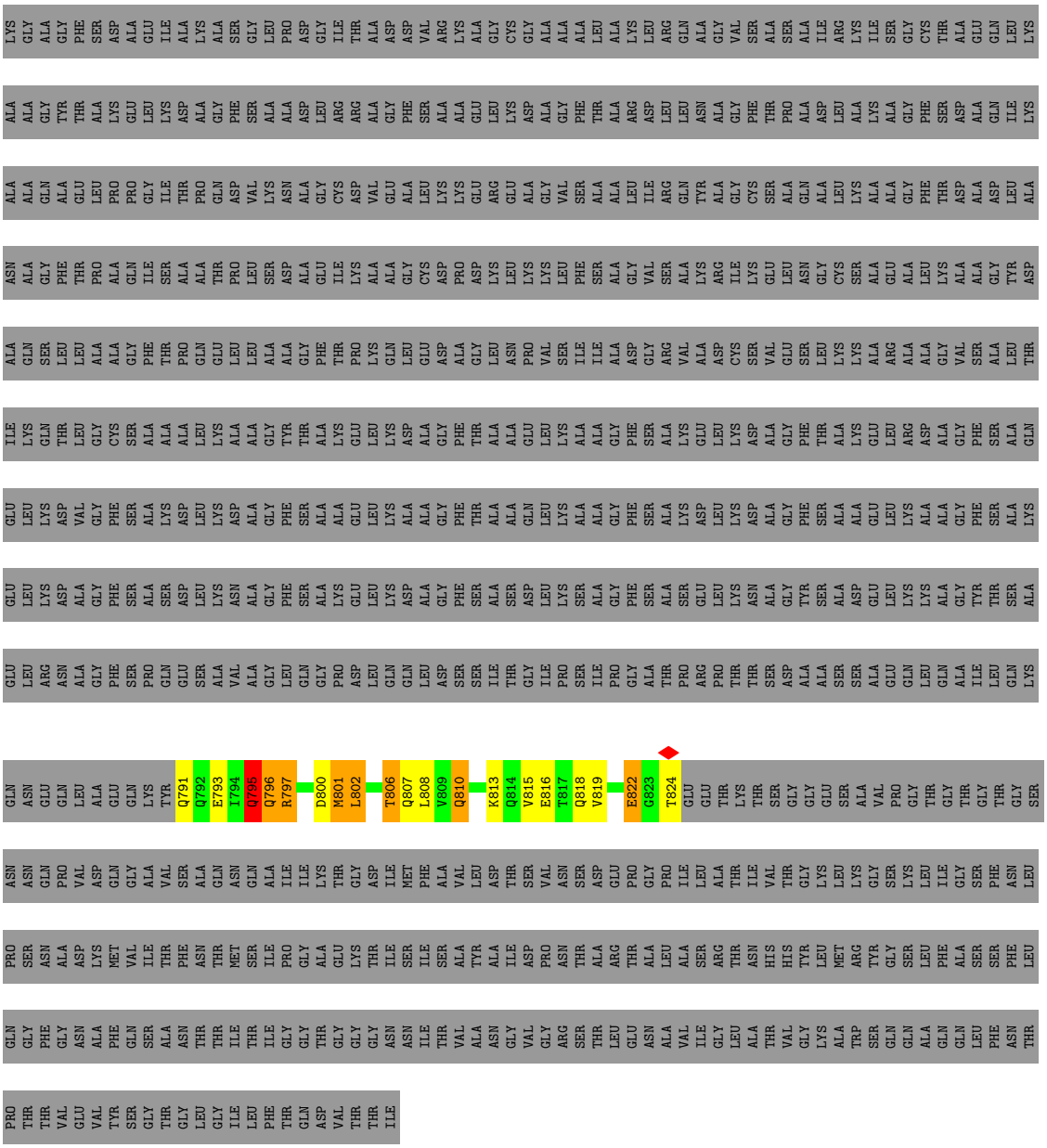


• Molecule 4: IcmE protein

Chain FG:

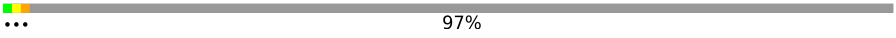


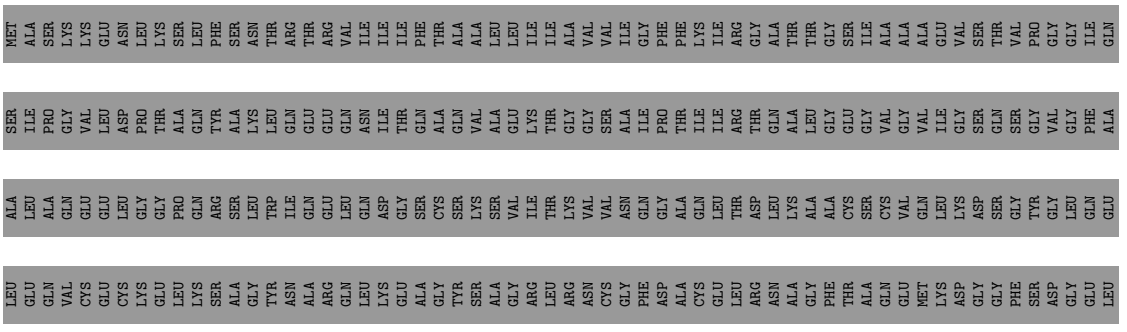


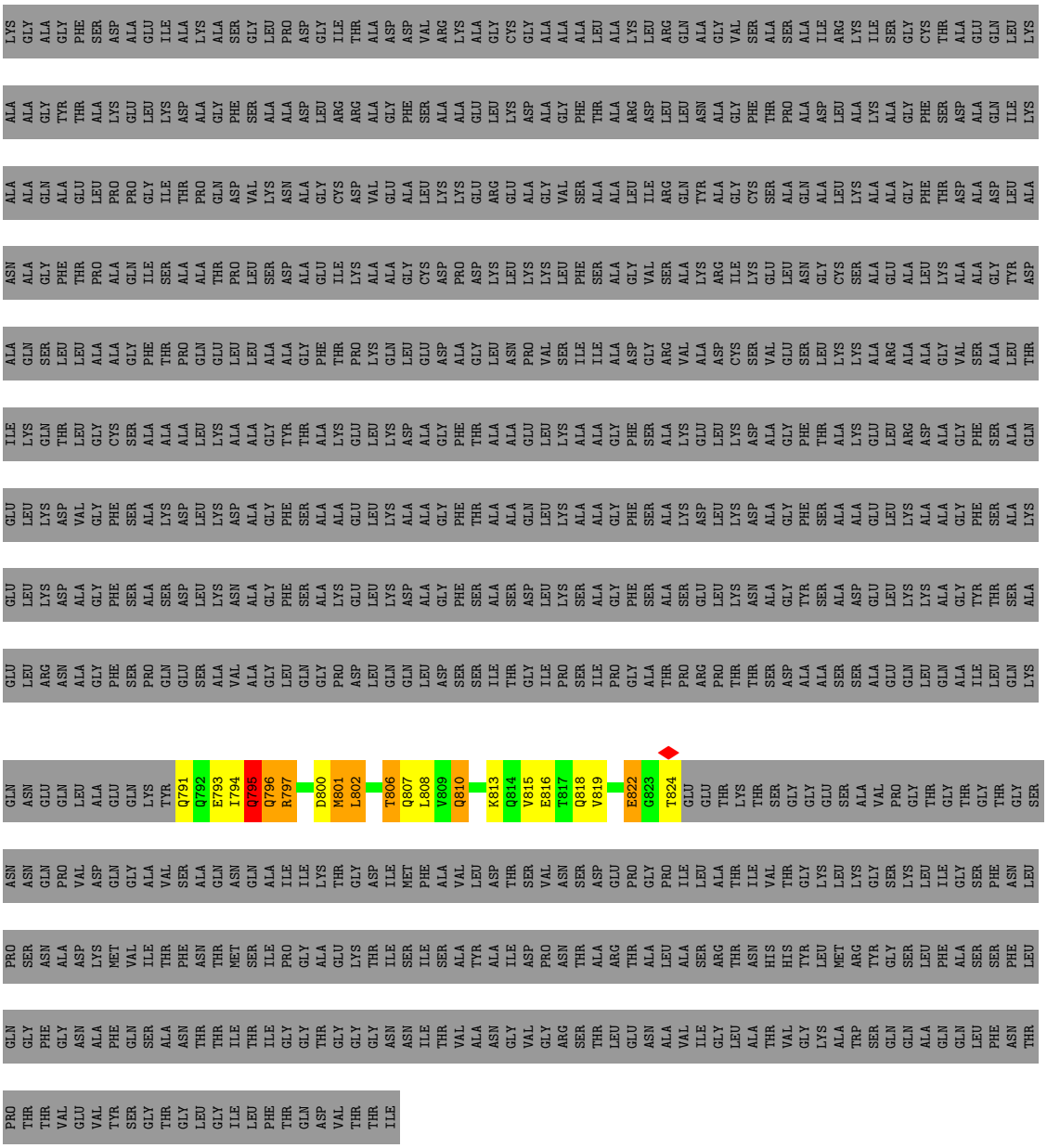


• Molecule 4: IcmE protein

Chain GG:

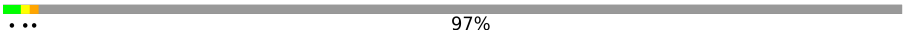


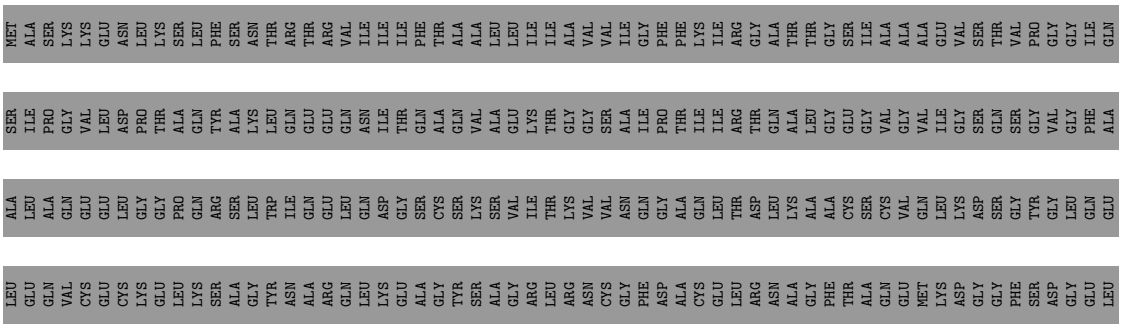


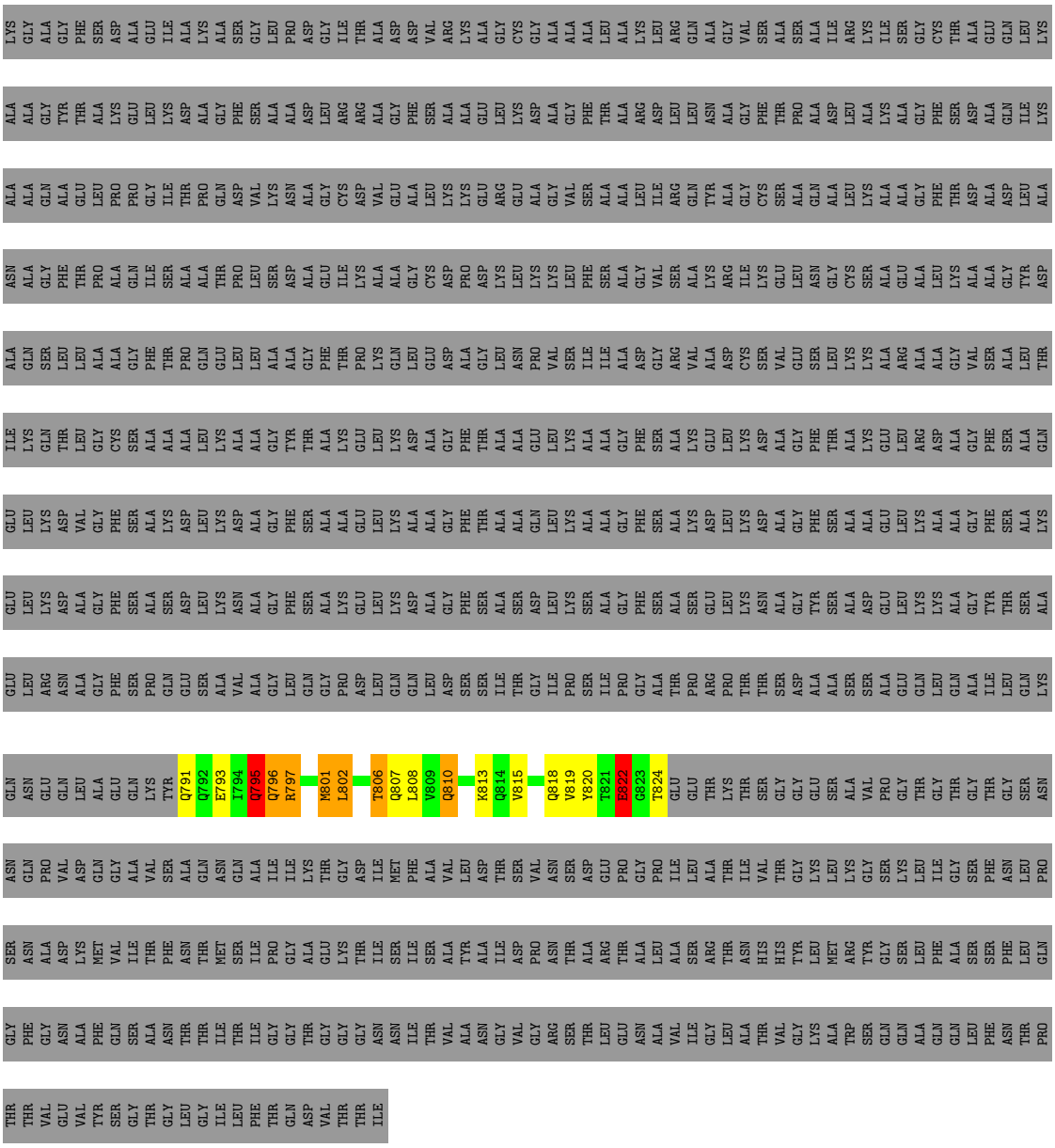


• Molecule 4: IcmE protein

Chain HG:

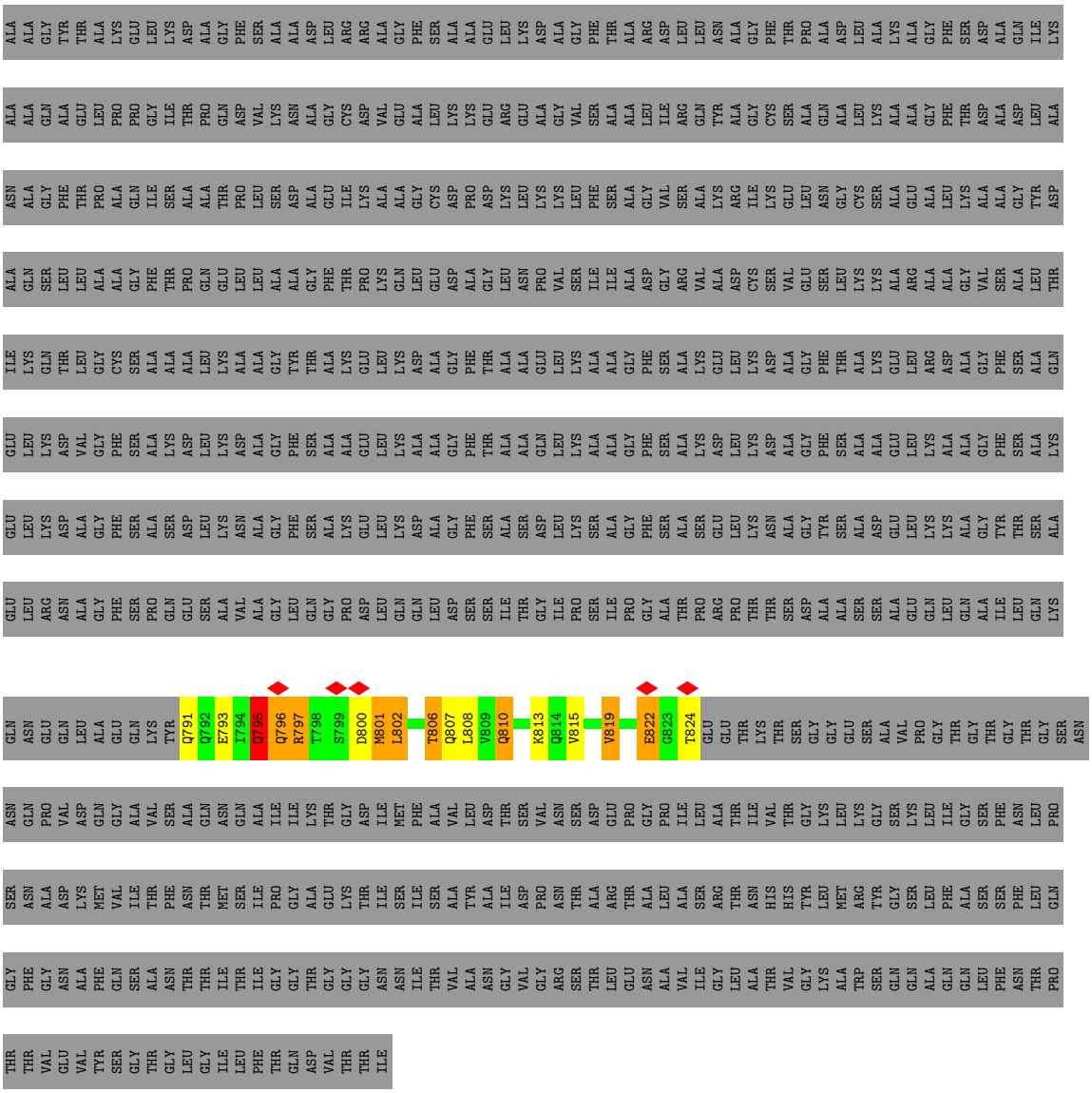






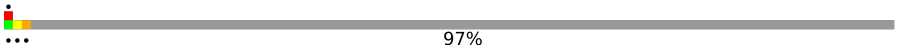
• Molecule 4: IcmE protein

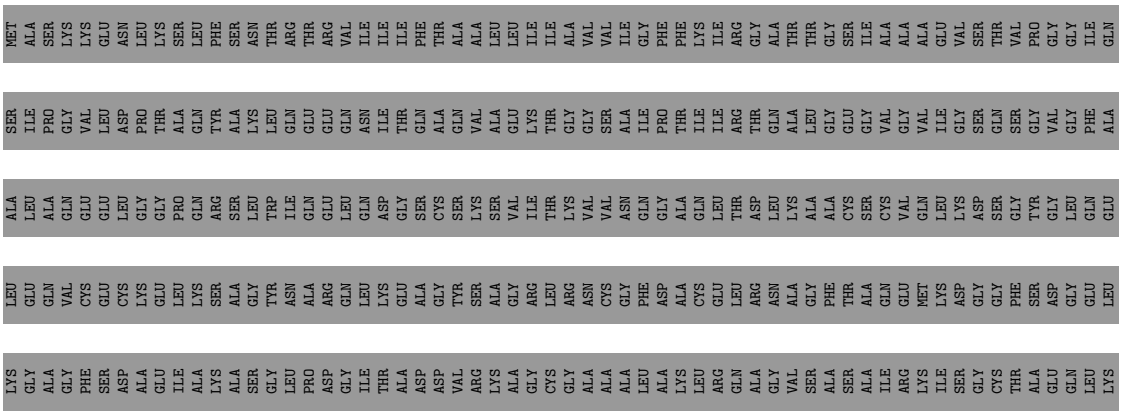
Chain IG: 97%

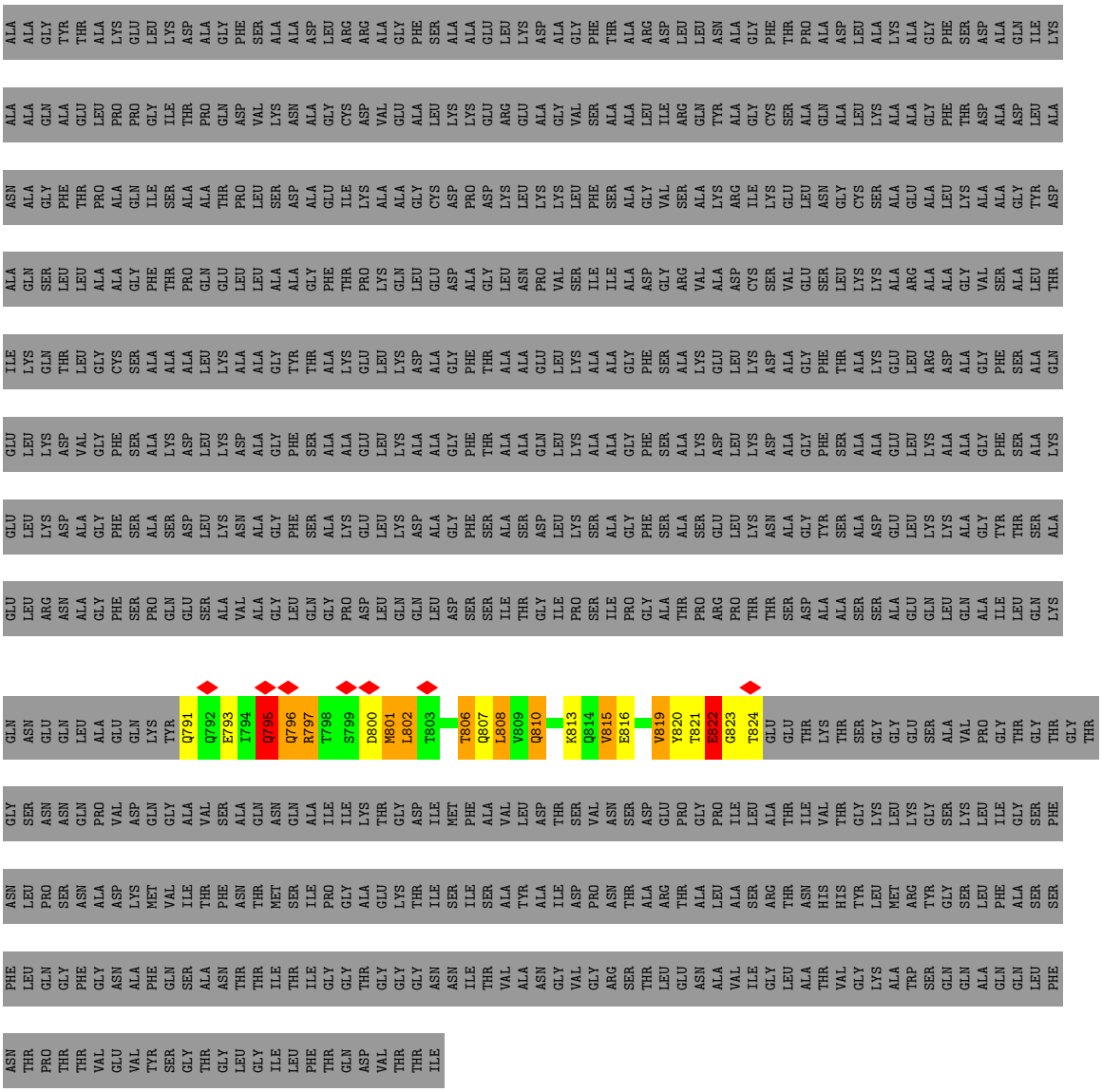


• Molecule 4: IcmE protein

Chain JG:

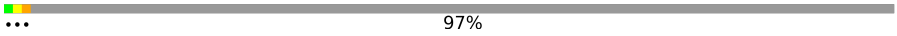


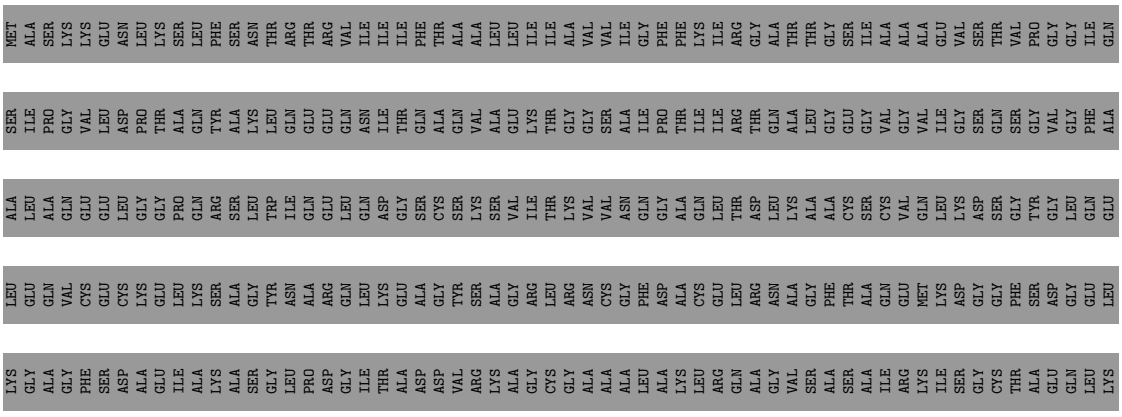


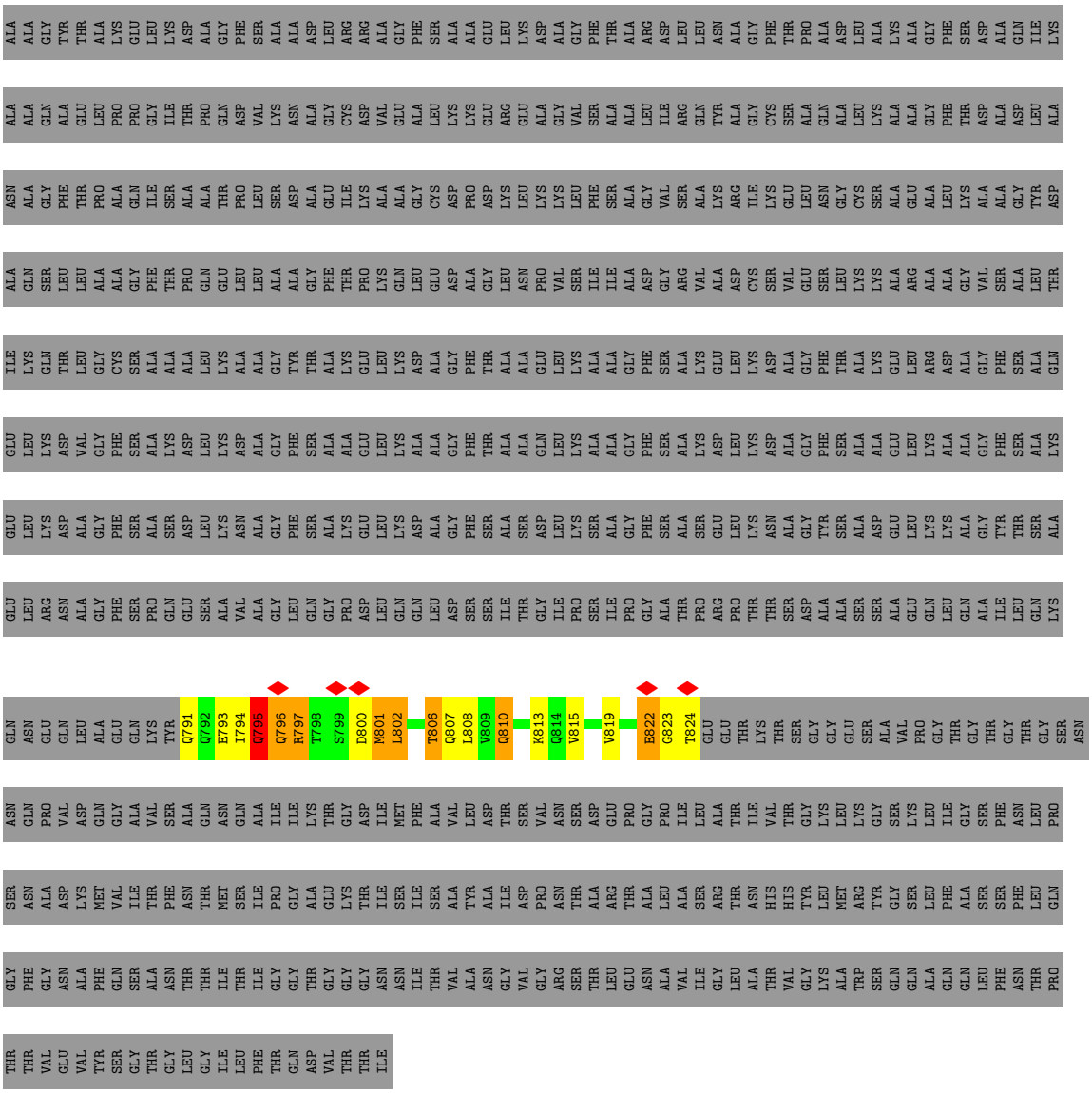


• Molecule 4: IcmE protein

Chain KG:

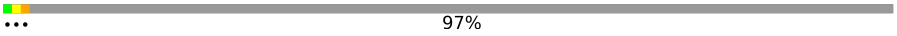


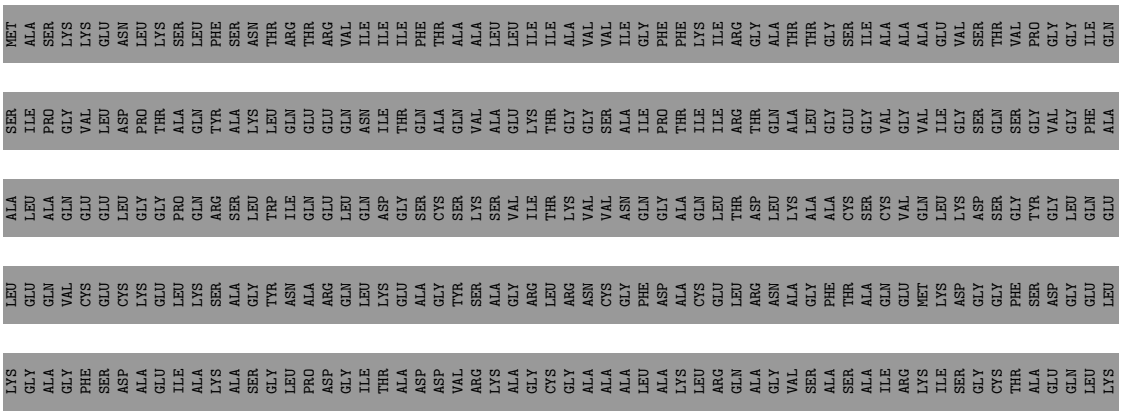


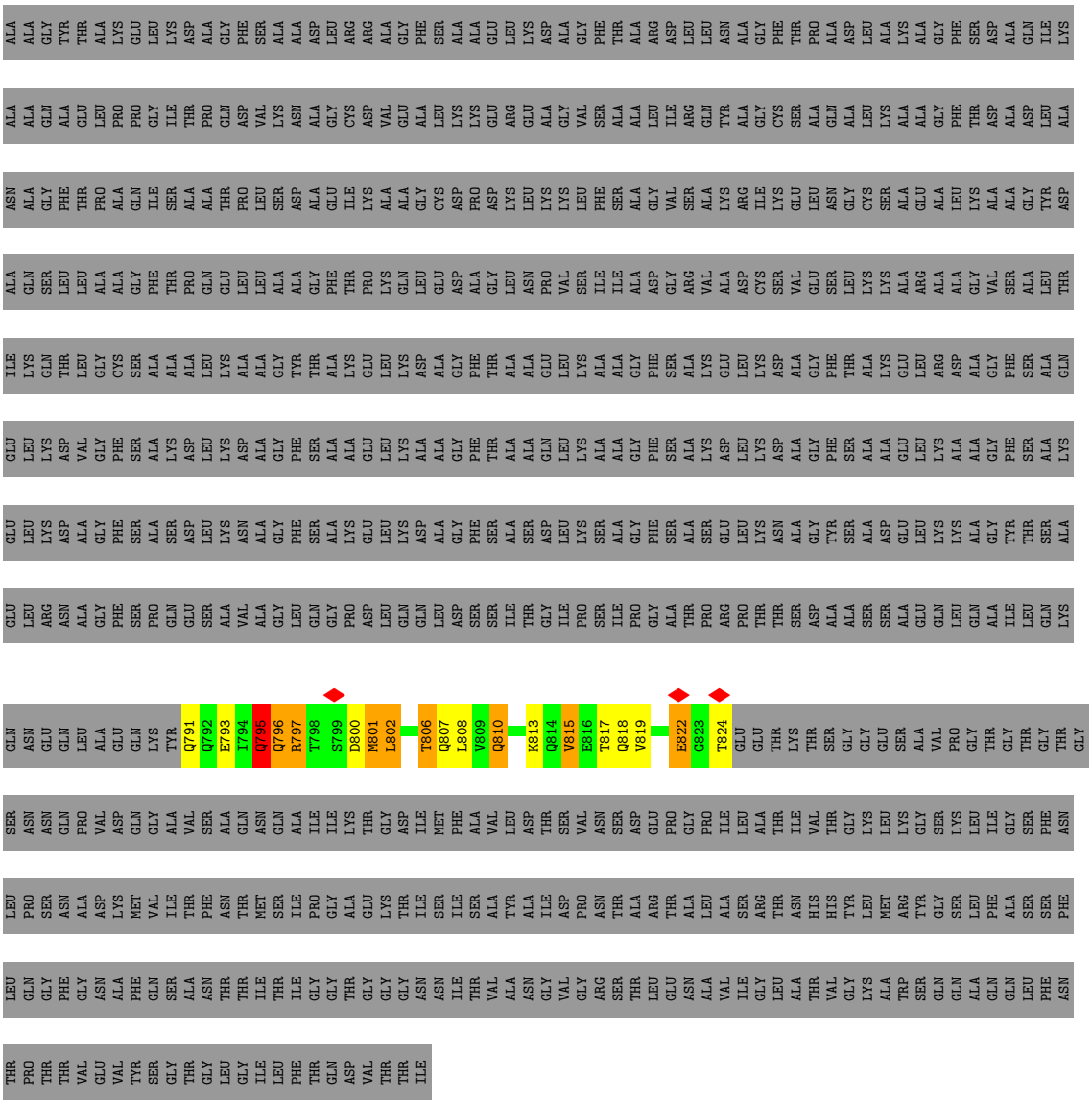


• Molecule 4: IcmE protein

Chain LG:

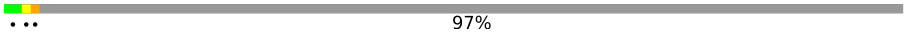


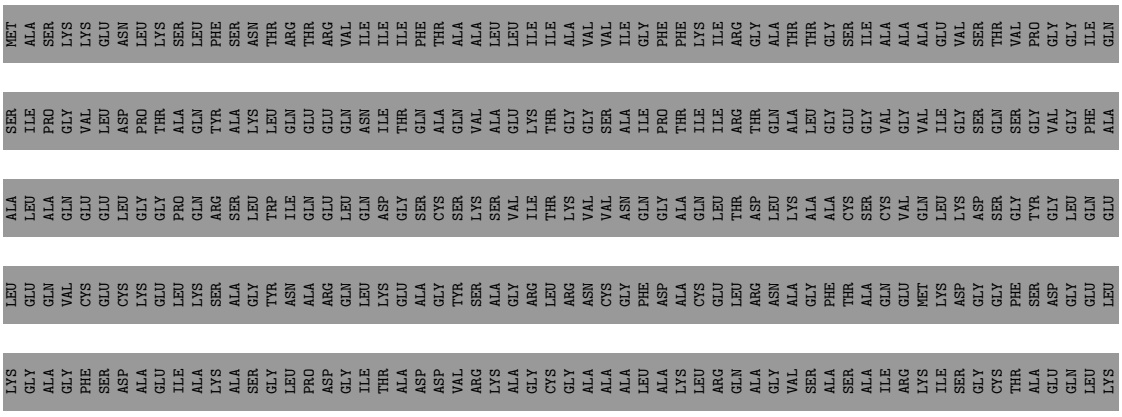


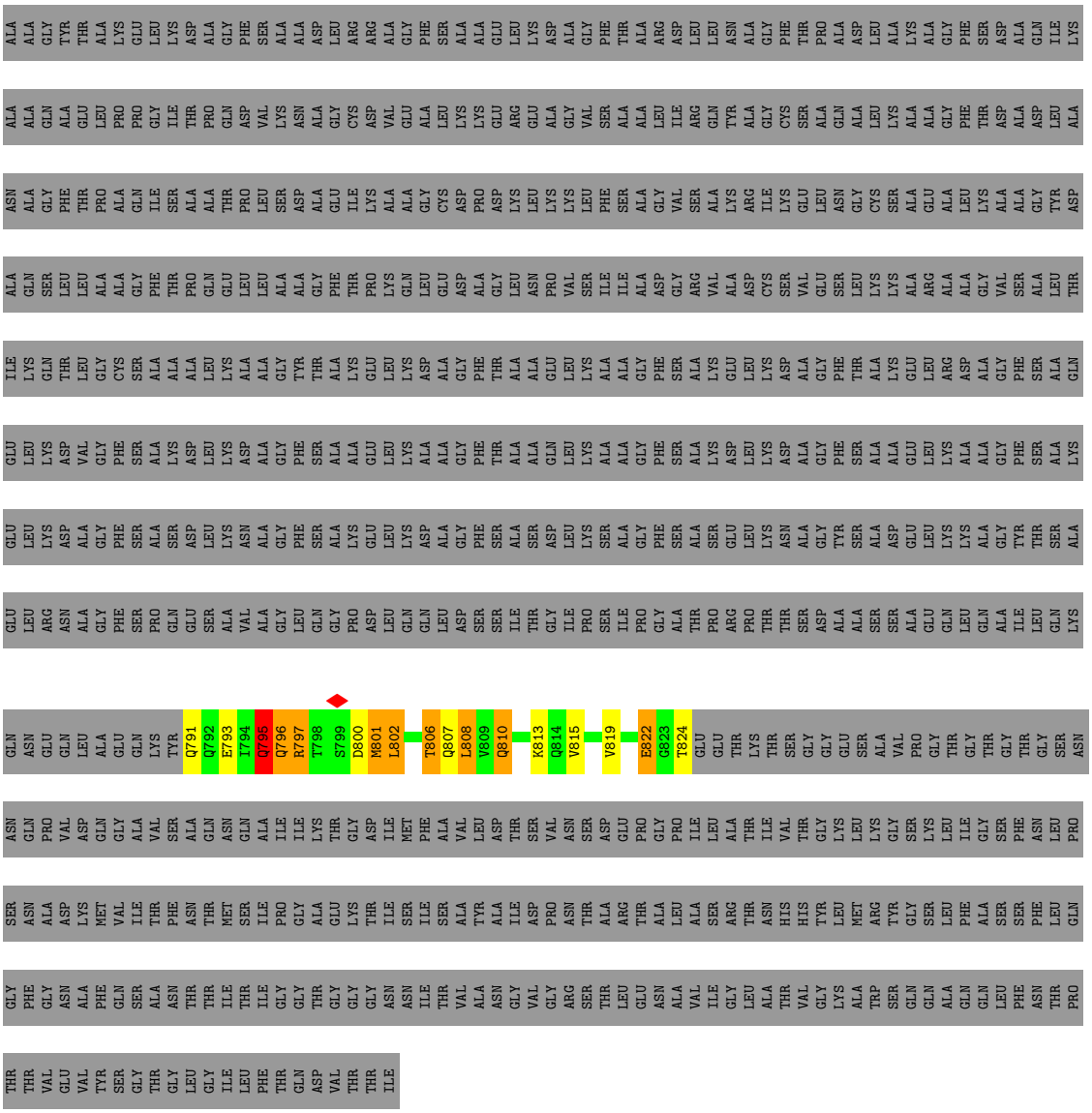


• Molecule 4: IcmE protein

Chain MG:

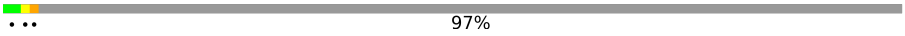


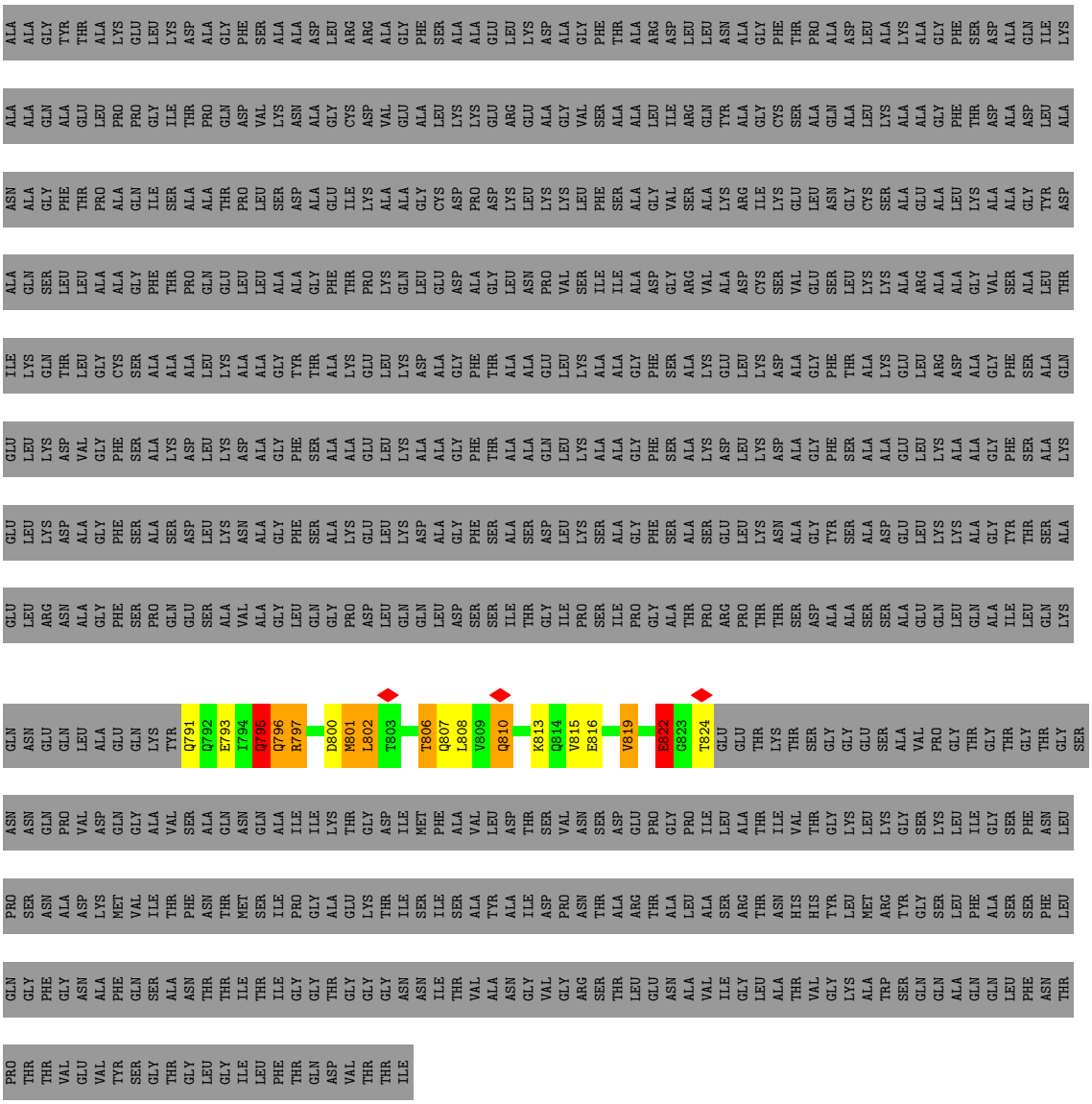




• Molecule 4: IcmE protein

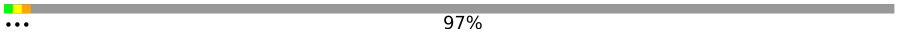
Chain VG:

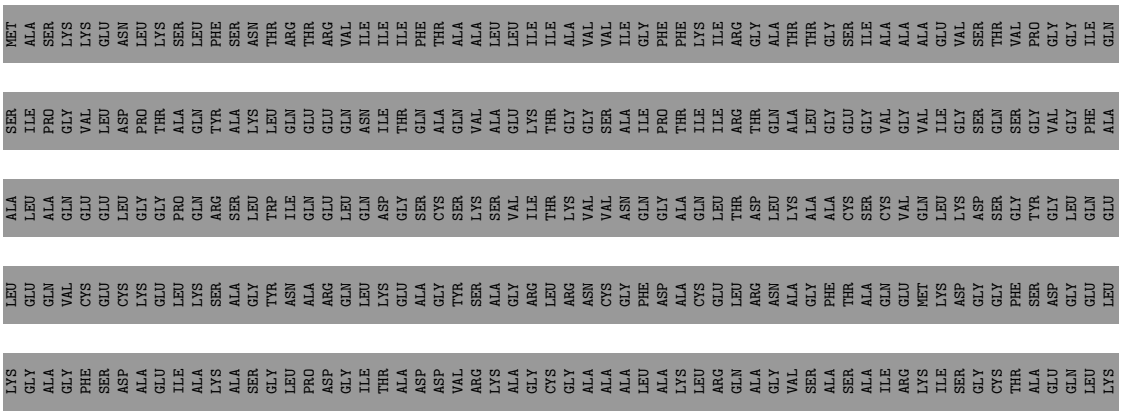


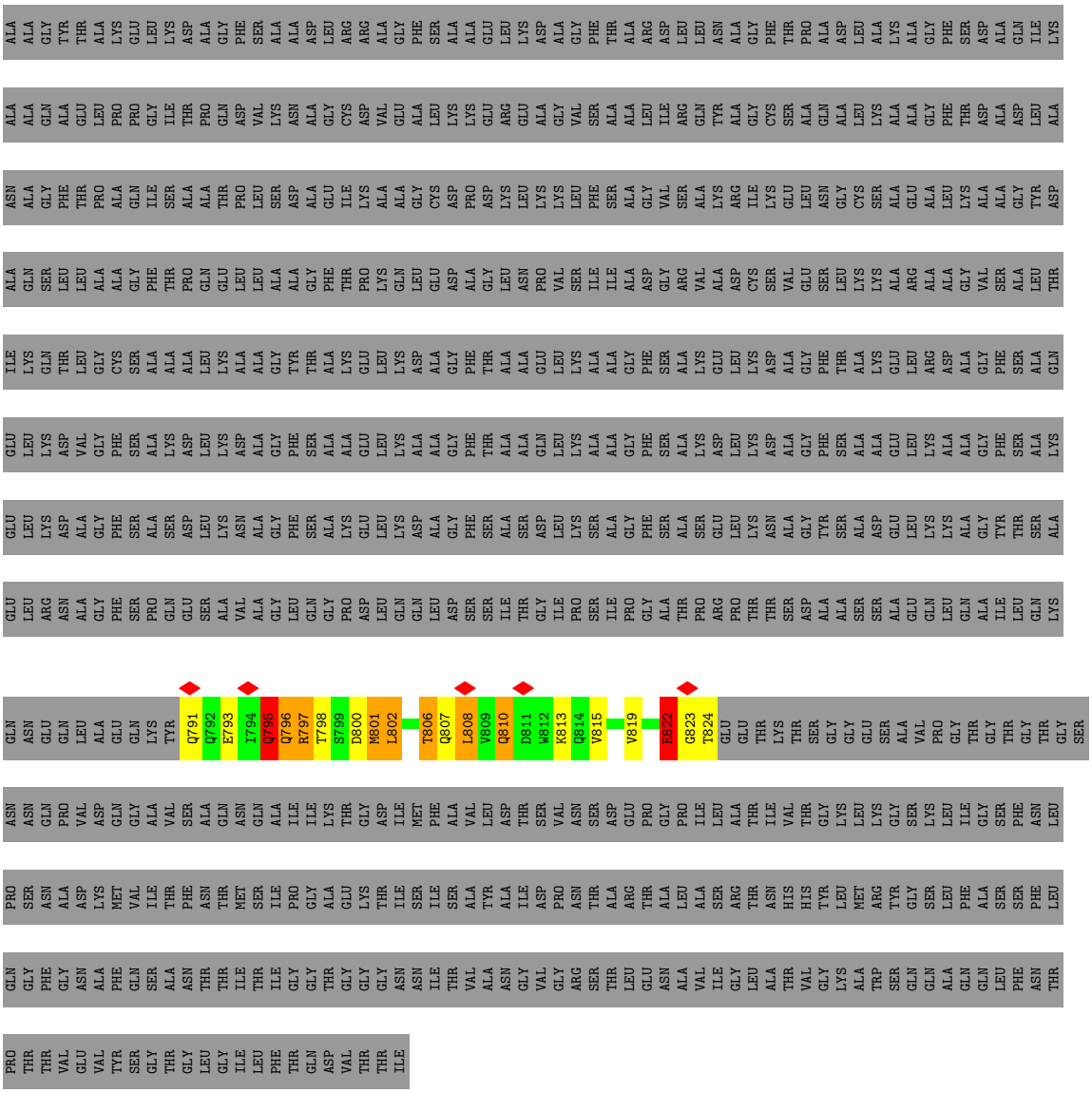


• Molecule 4: IcmE protein

Chain WG:

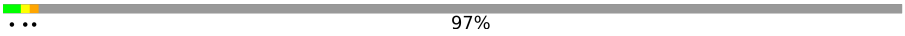


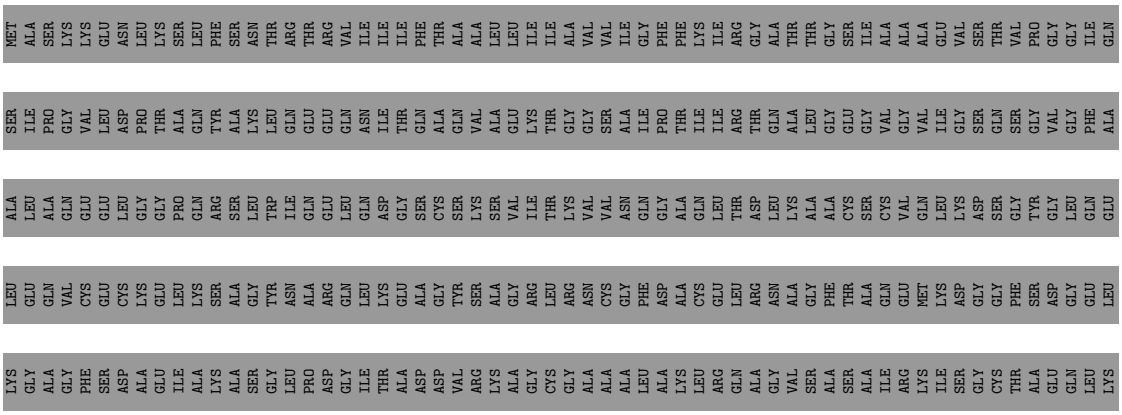


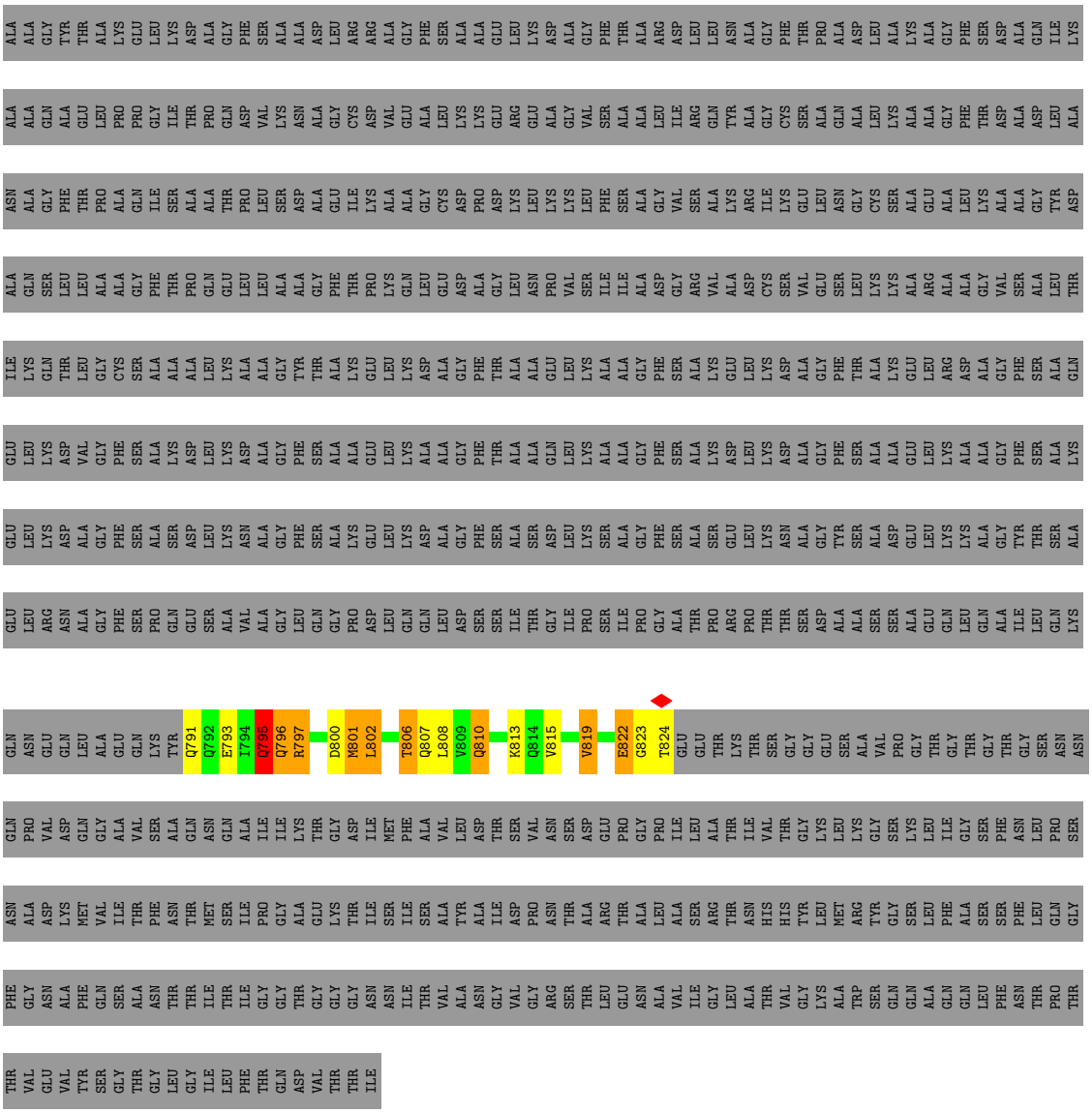


• Molecule 4: IcmE protein

Chain XG:

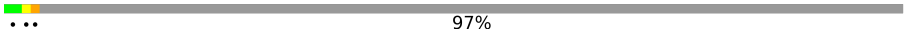


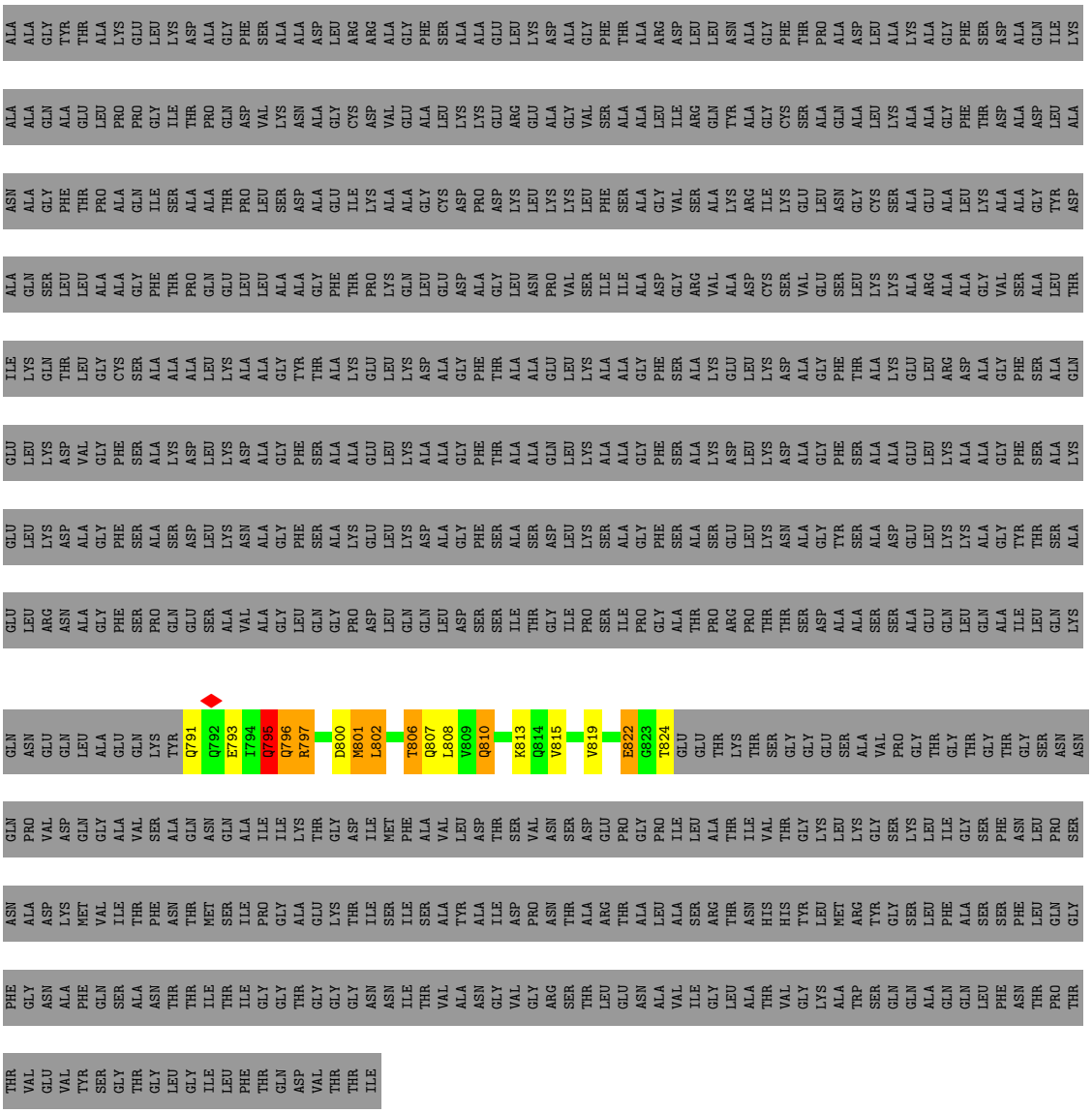




• Molecule 4: IcmE protein

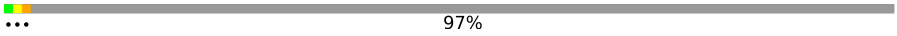
Chain YG:

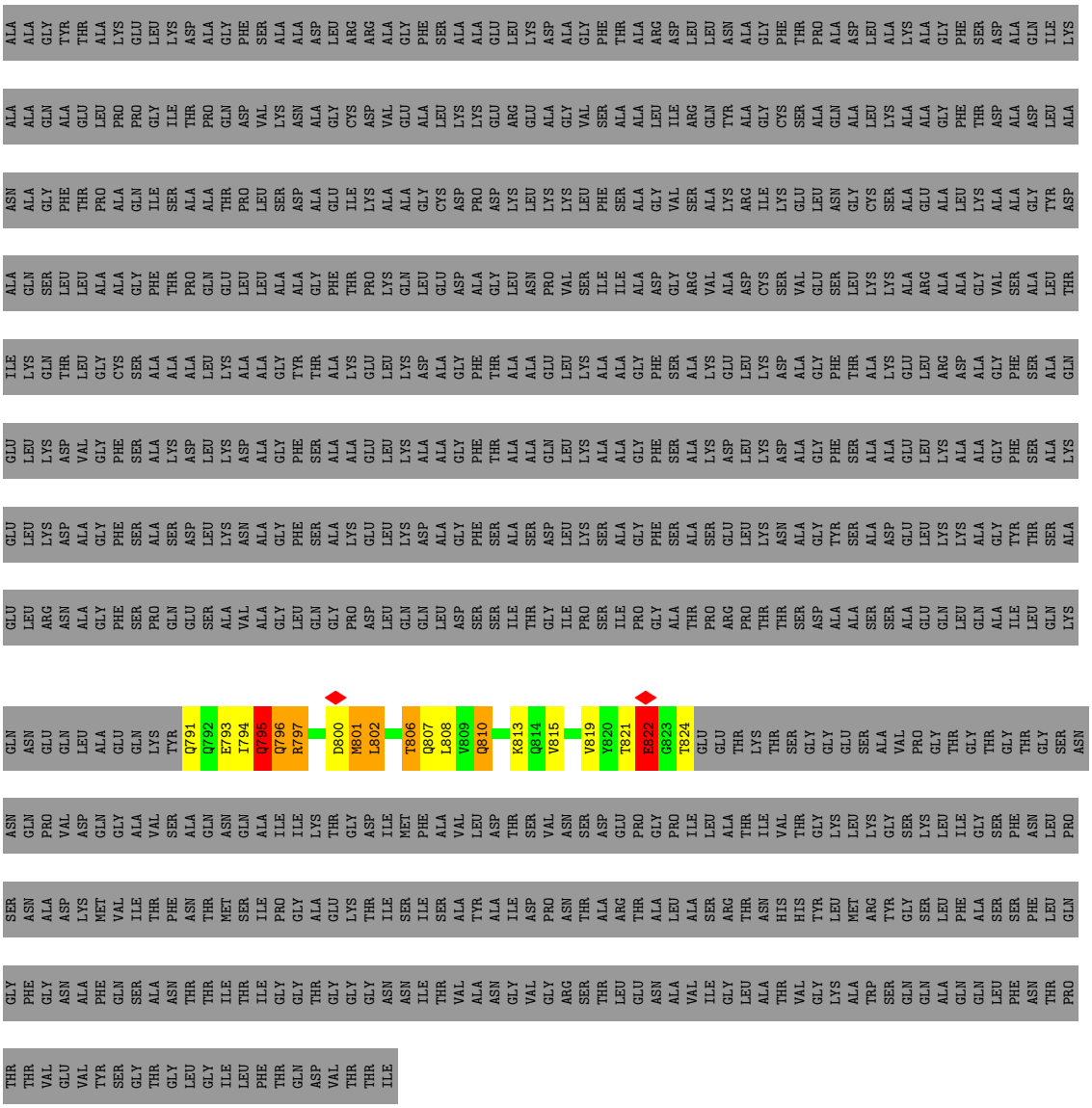




• Molecule 4: IcmE protein

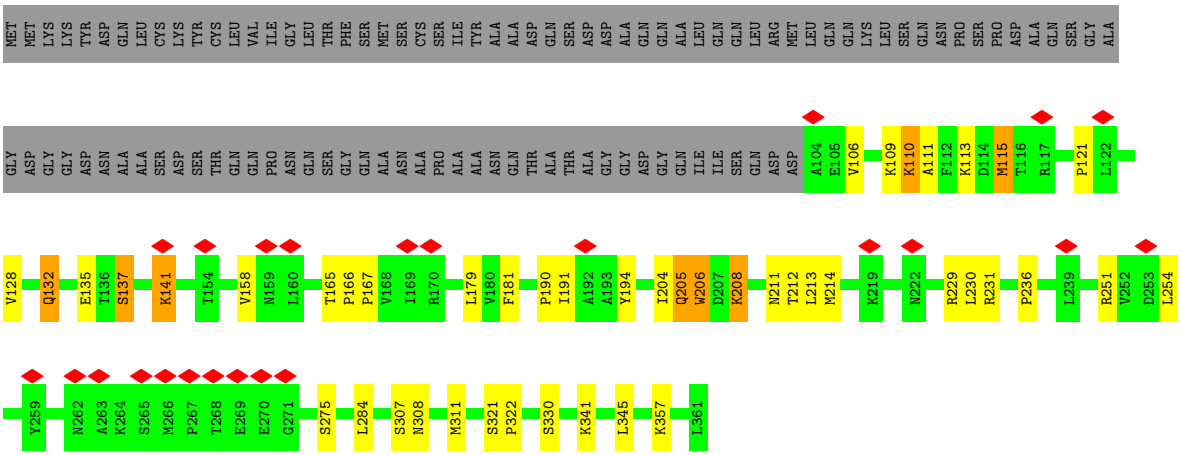
Chain ZG:





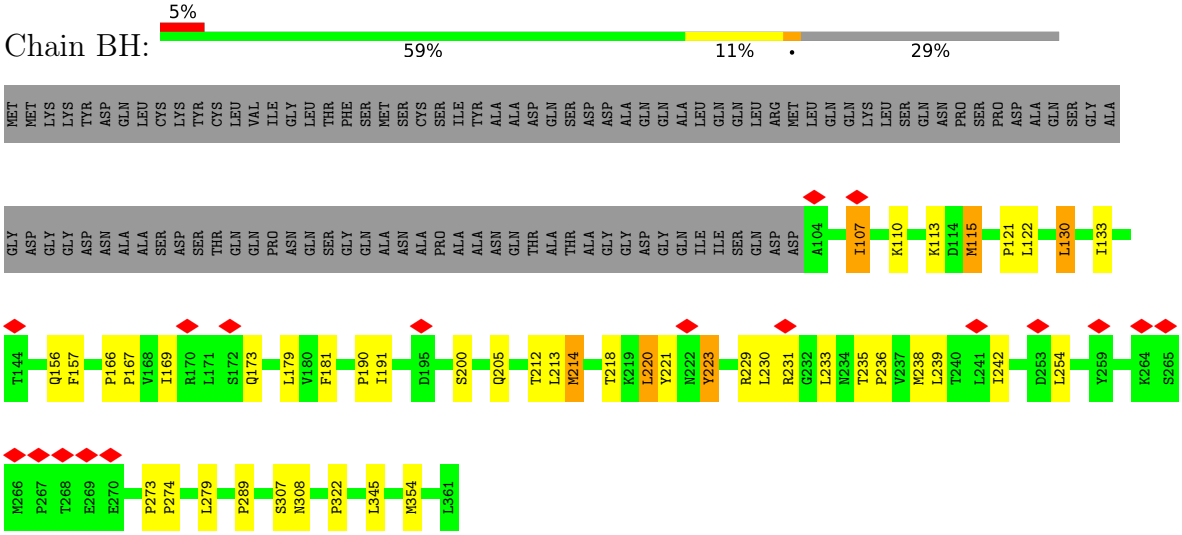
• Molecule 5: Type IV secretion protein IcmK

Chain AH:



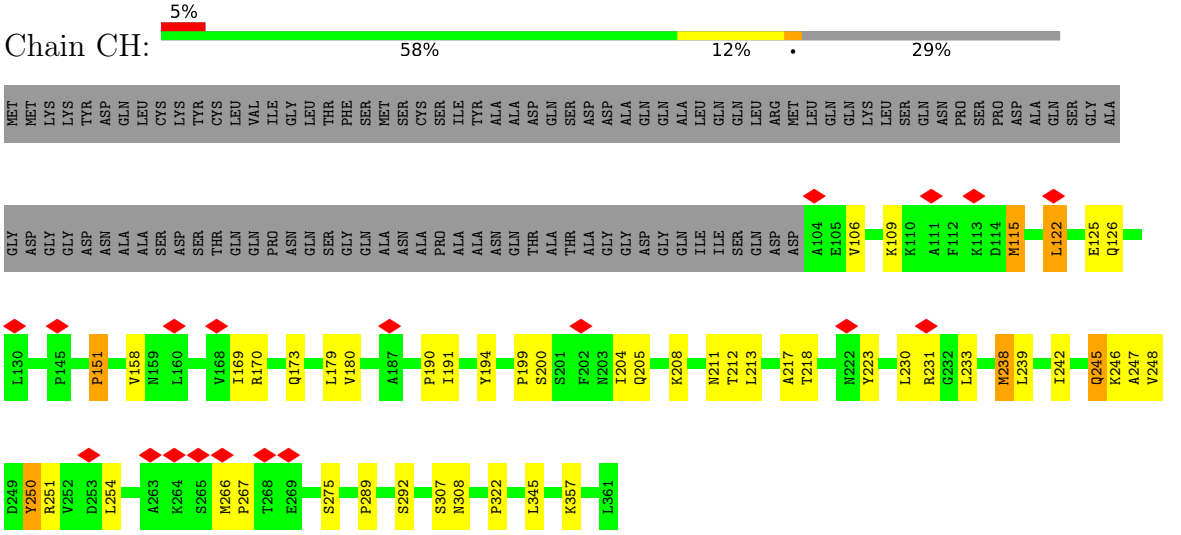
• Molecule 5: Type IV secretion protein IcmK

Chain BH:



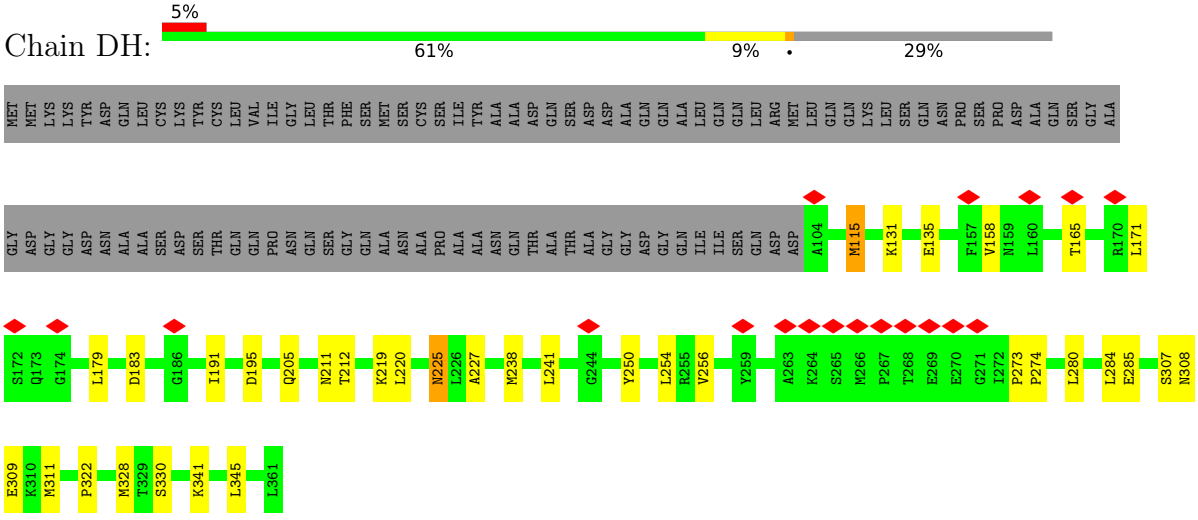
• Molecule 5: Type IV secretion protein IcmK

Chain CH:



• Molecule 5: Type IV secretion protein IcmK

Chain DH:



• Molecule 5: Type IV secretion protein IcmK

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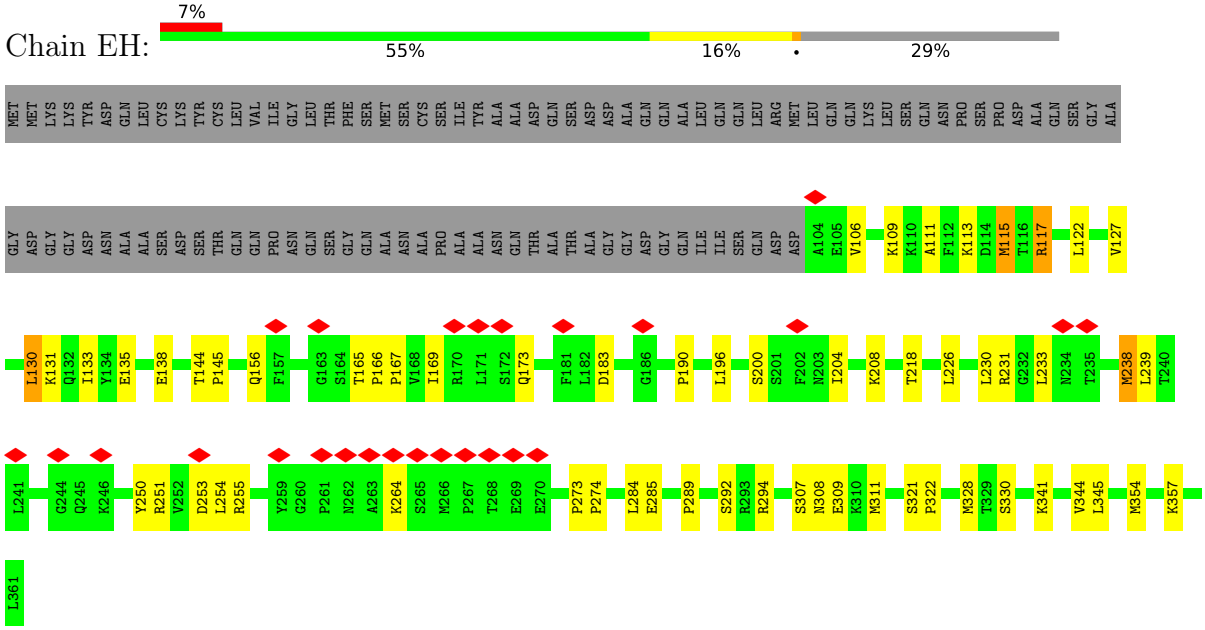
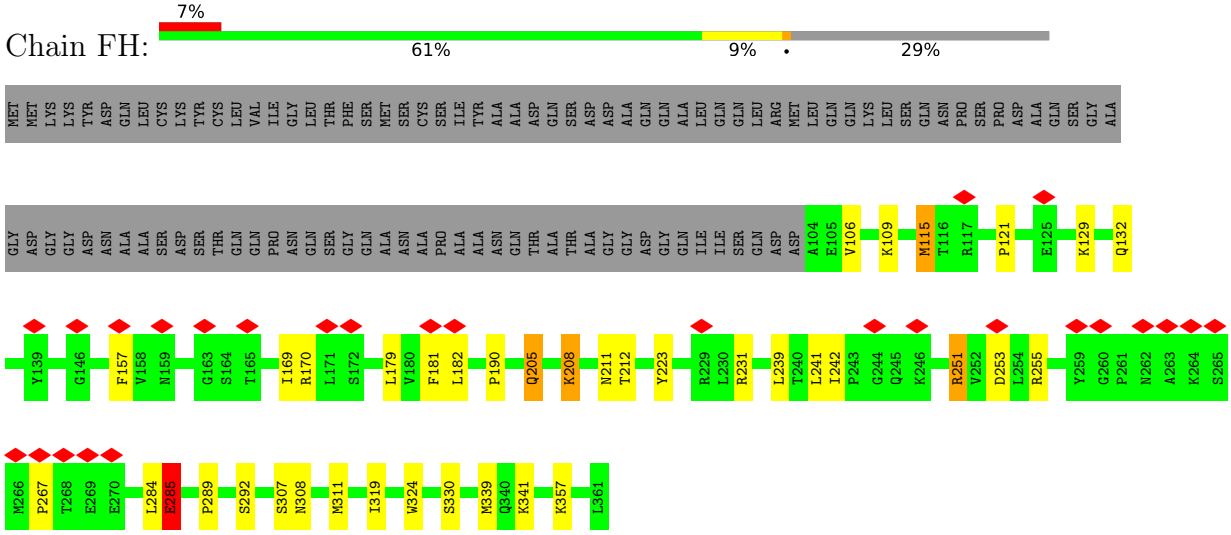


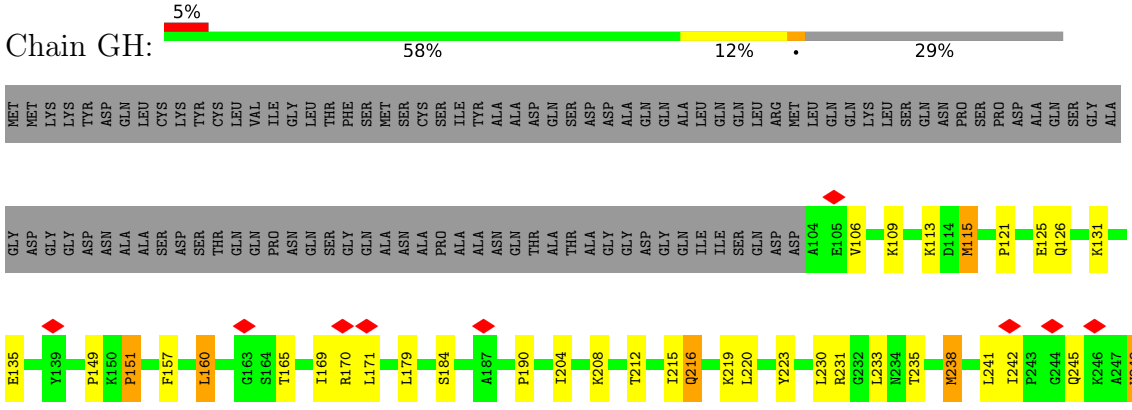
• Molecule 5: Type IV secretion protein IcmK

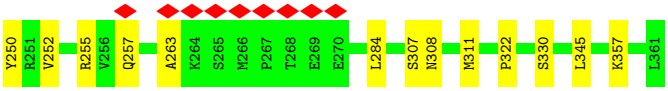
Chain FH:



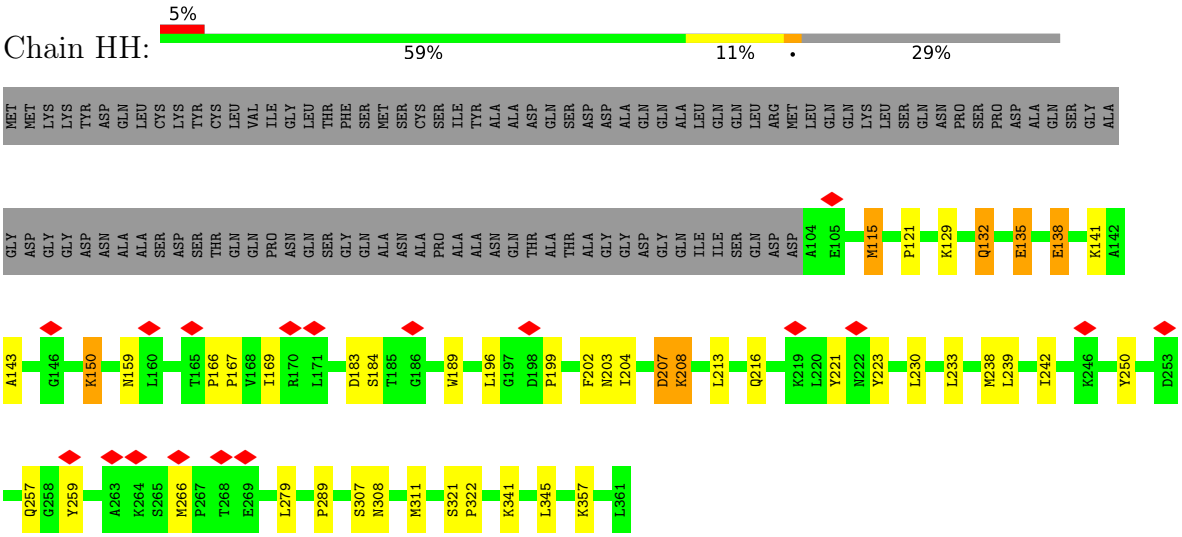
• Molecule 5: Type IV secretion protein IcmK

Chain GH:

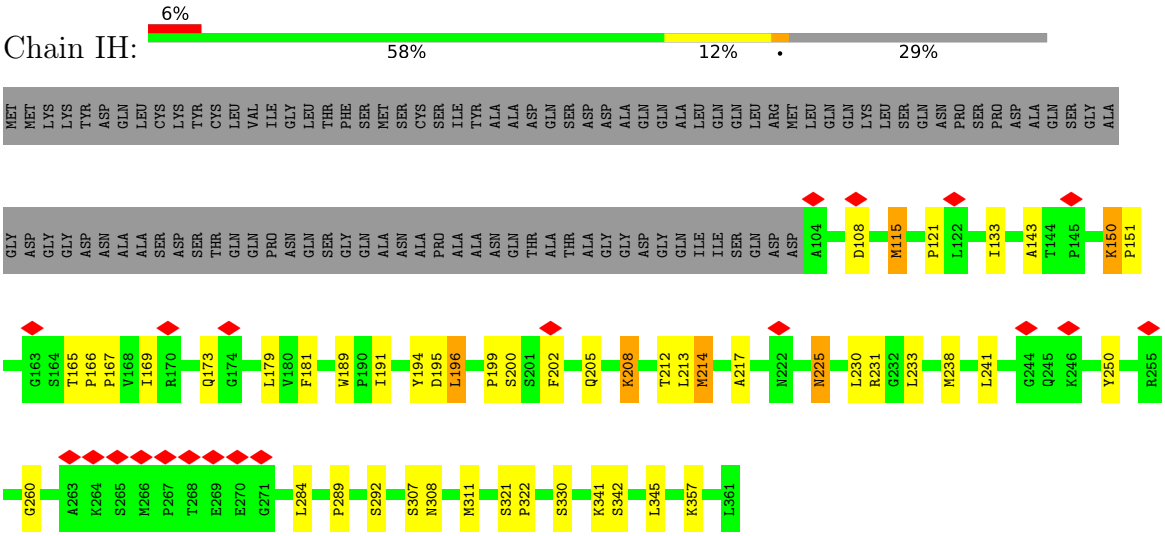




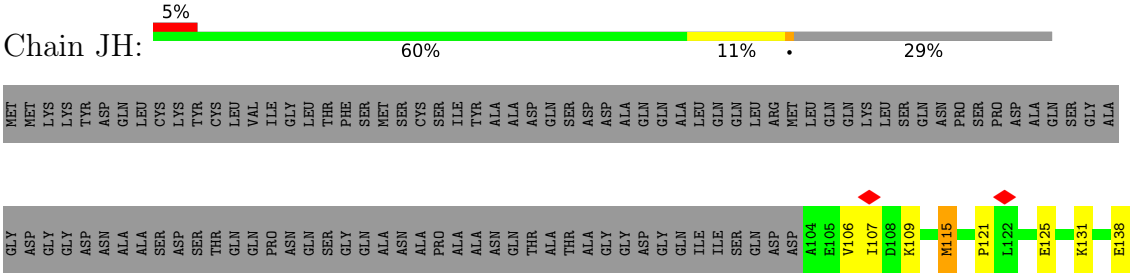
• Molecule 5: Type IV secretion protein IcmK

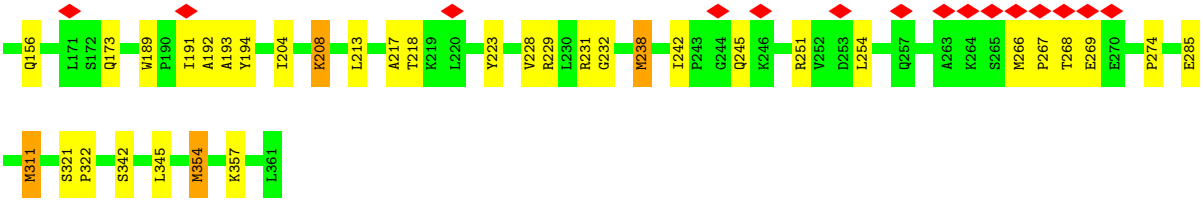


• Molecule 5: Type IV secretion protein IcmK

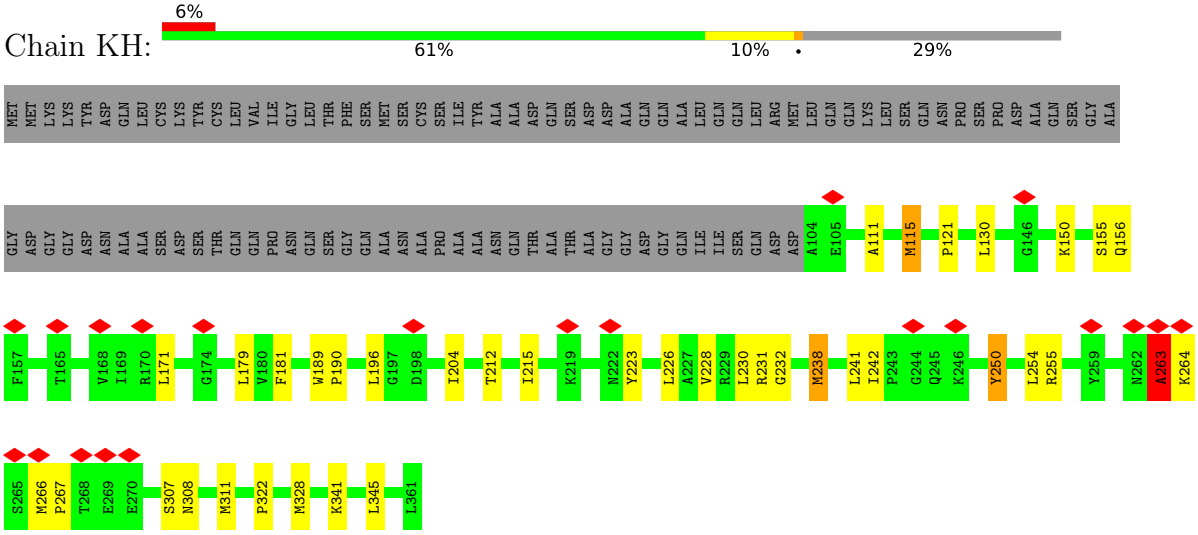


• Molecule 5: Type IV secretion protein IcmK

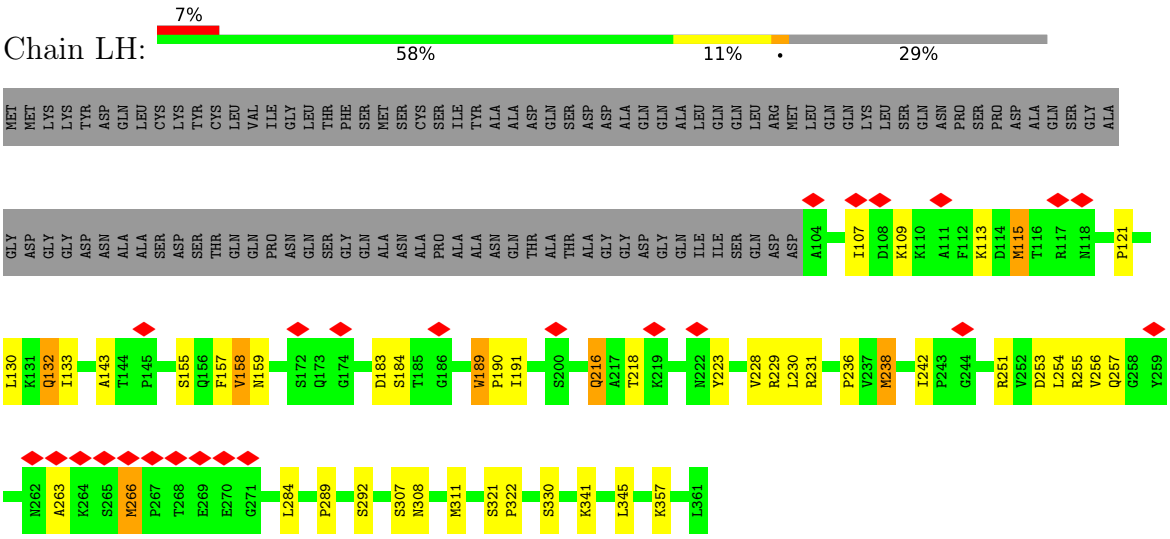




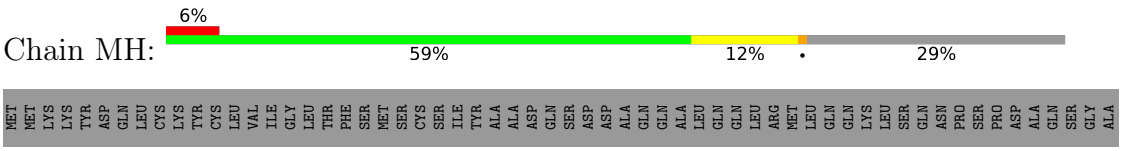
• Molecule 5: Type IV secretion protein IcmK

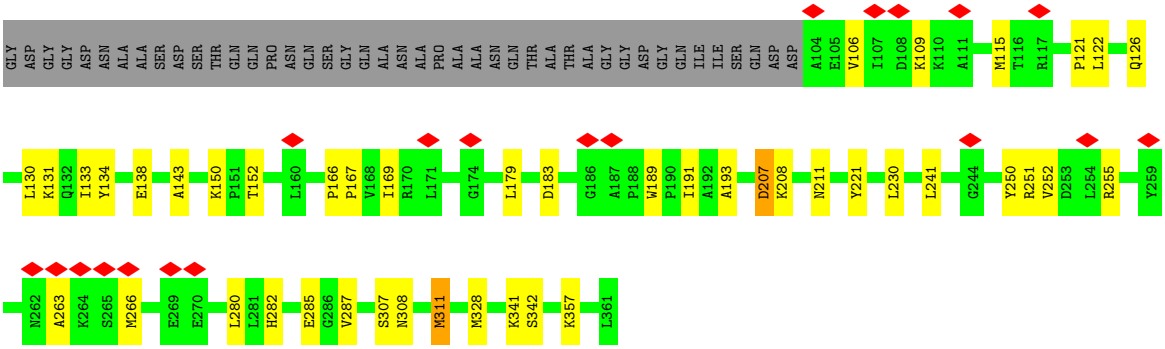


• Molecule 5: Type IV secretion protein IcmK



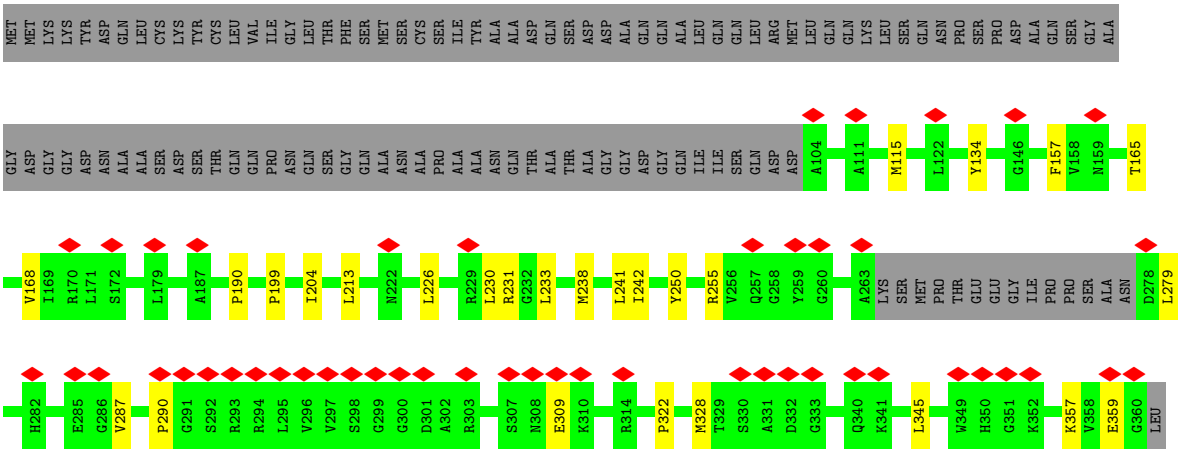
• Molecule 5: Type IV secretion protein IcmK





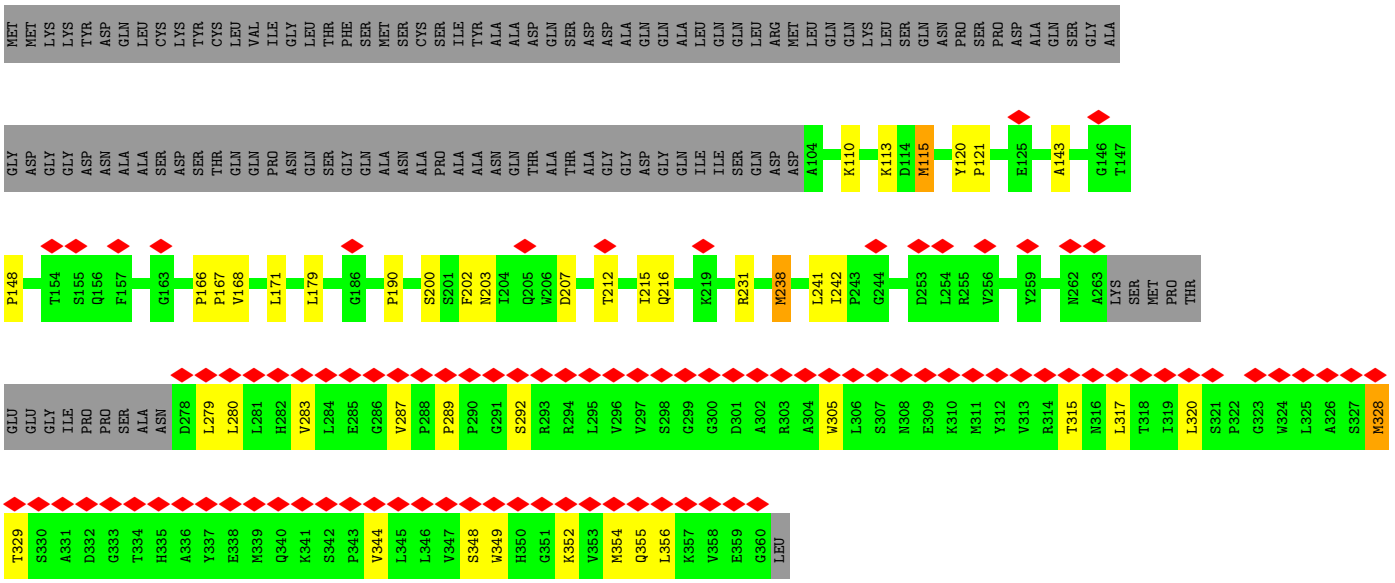
• Molecule 5: Type IV secretion protein IcmK

Chain VH:



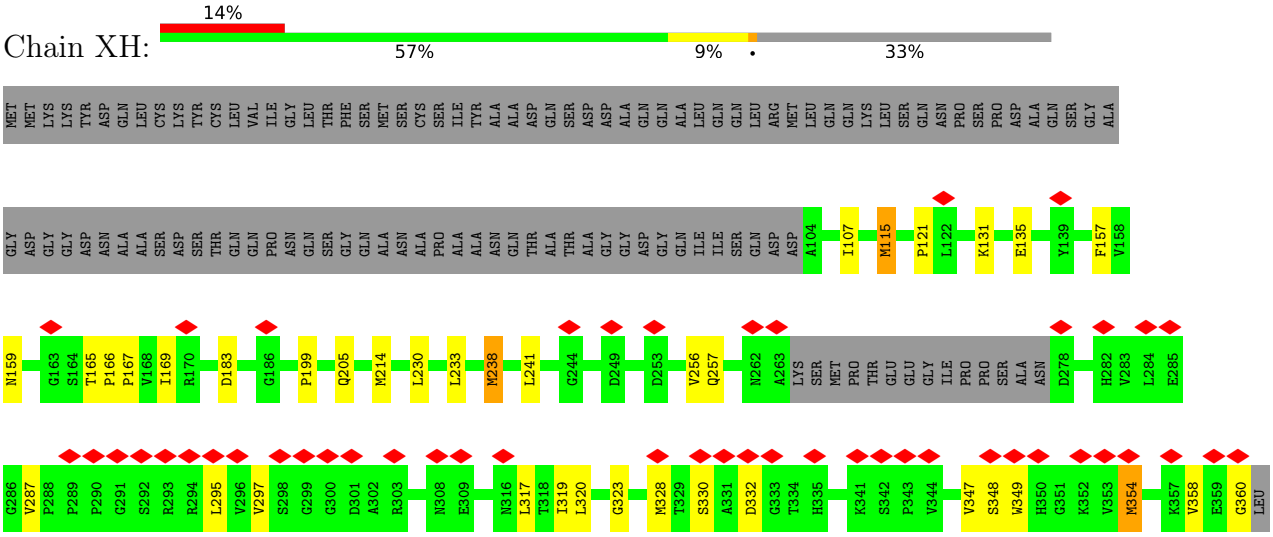
• Molecule 5: Type IV secretion protein IcmK

Chain WH:



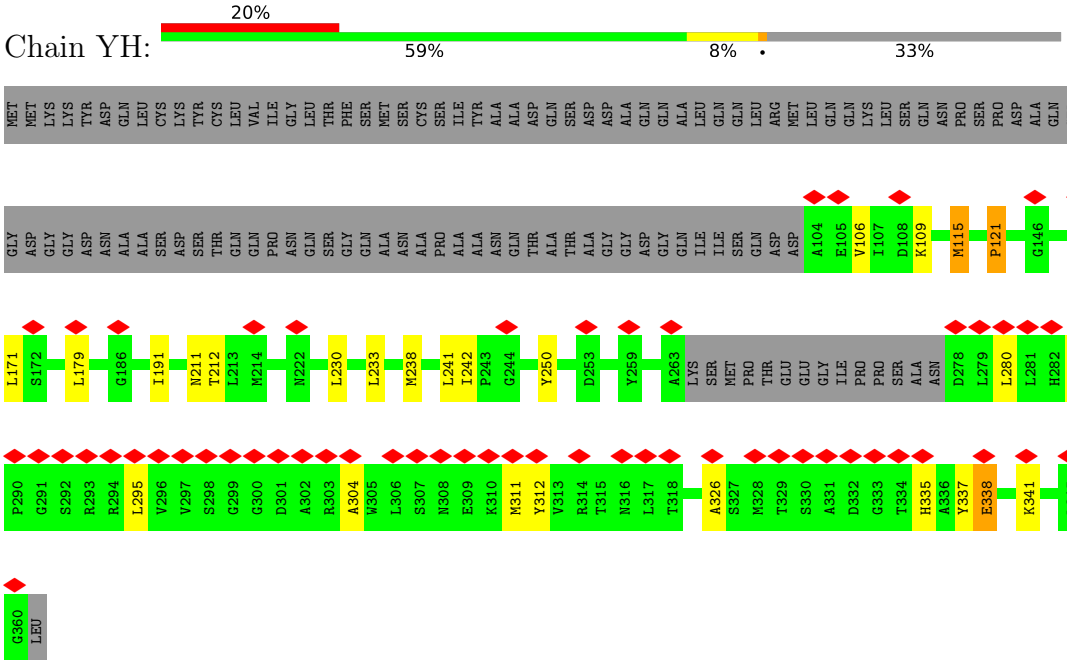
• Molecule 5: Type IV secretion protein IcmK

Chain XH:



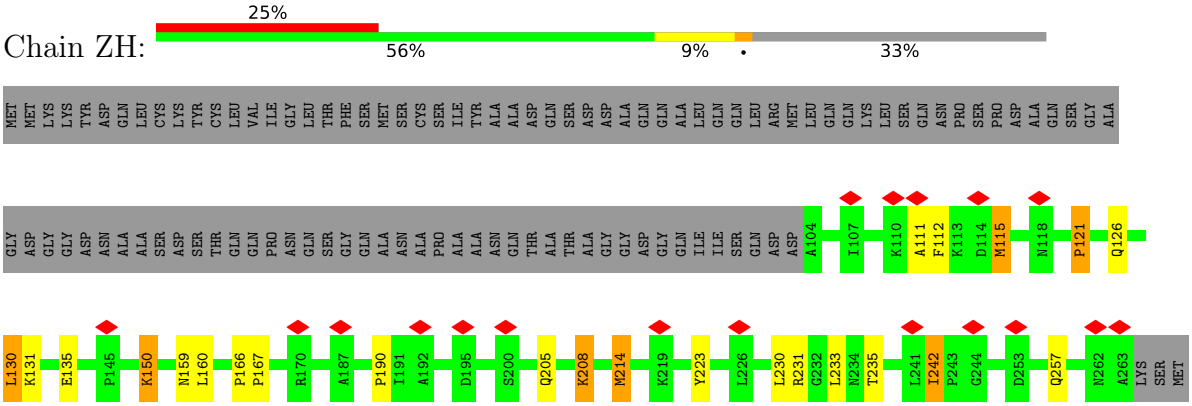
• Molecule 5: Type IV secretion protein IcmK

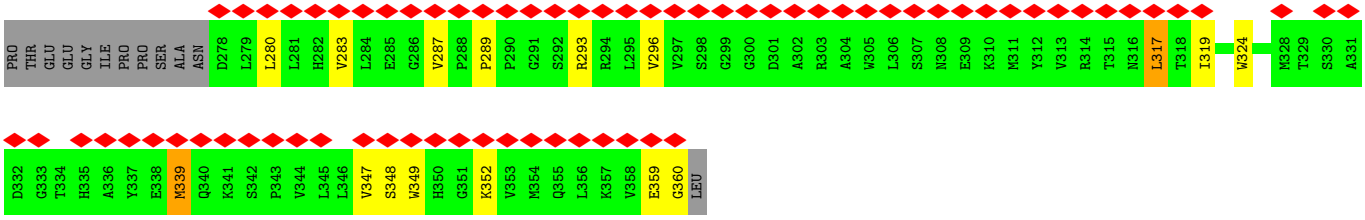
Chain YH:



• Molecule 5: Type IV secretion protein IcmK

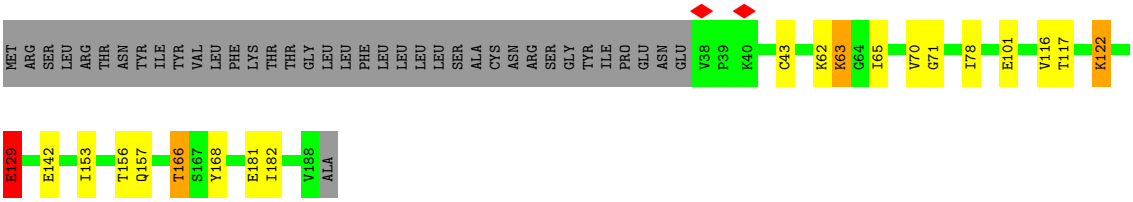
Chain ZH:





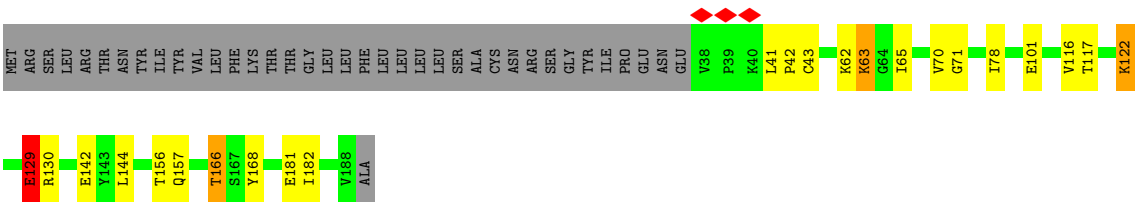
• Molecule 6: Inner membrane lipoprotein YiaD

Chain AK:



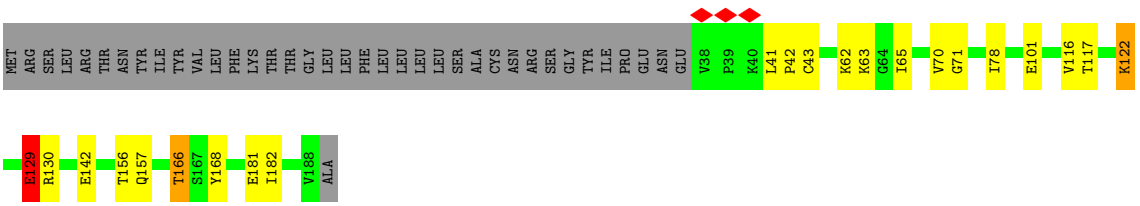
• Molecule 6: Inner membrane lipoprotein YiaD

Chain BK:



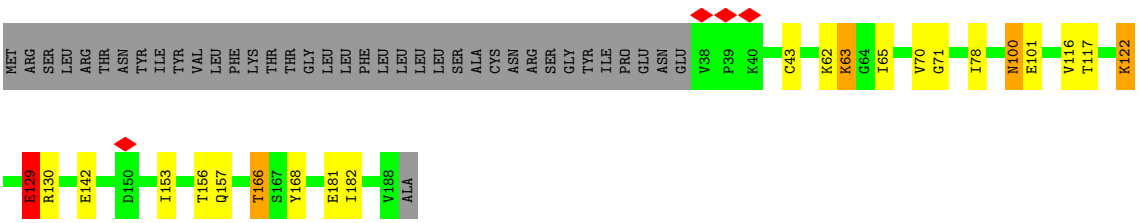
• Molecule 6: Inner membrane lipoprotein YiaD

Chain CK:



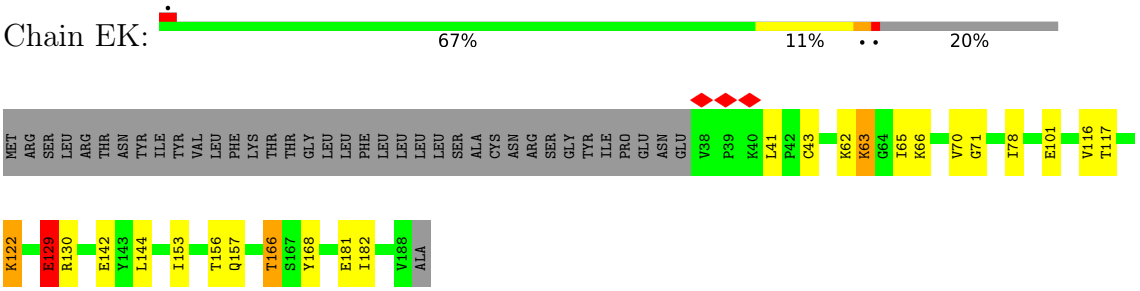
• Molecule 6: Inner membrane lipoprotein YiaD

Chain DK:



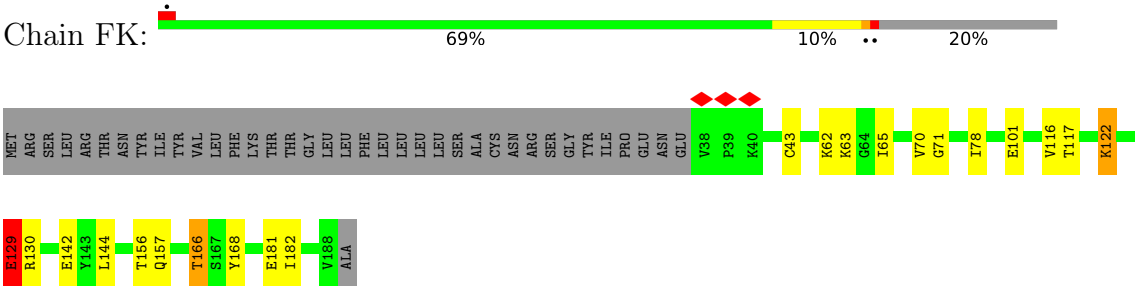
• Molecule 6: Inner membrane lipoprotein YiaD

Chain EK:



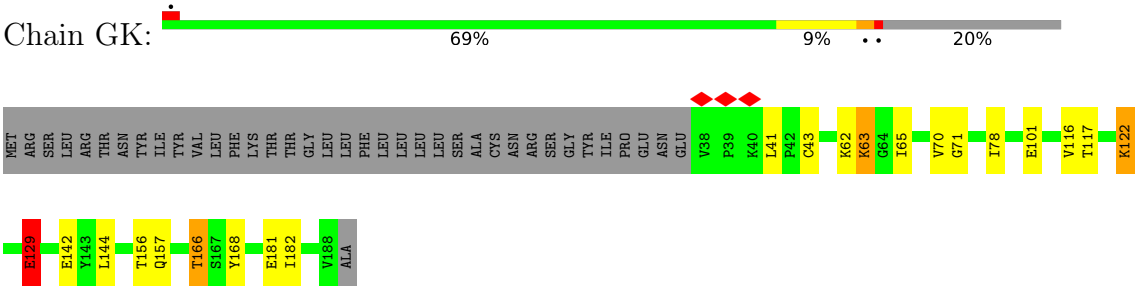
• Molecule 6: Inner membrane lipoprotein YiaD

Chain FK:



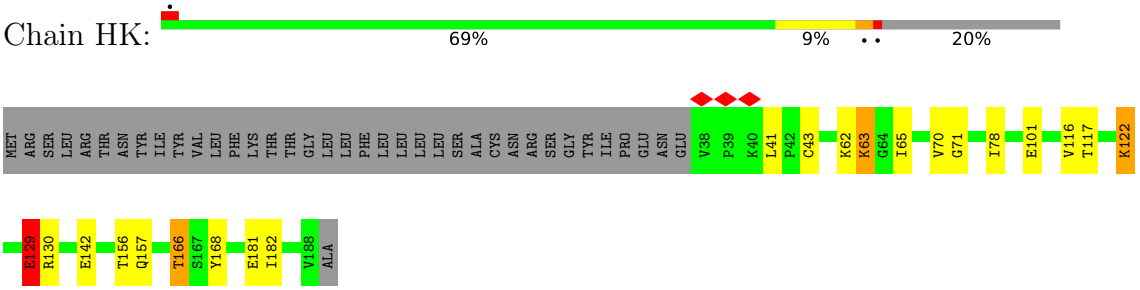
• Molecule 6: Inner membrane lipoprotein YiaD

Chain GK:



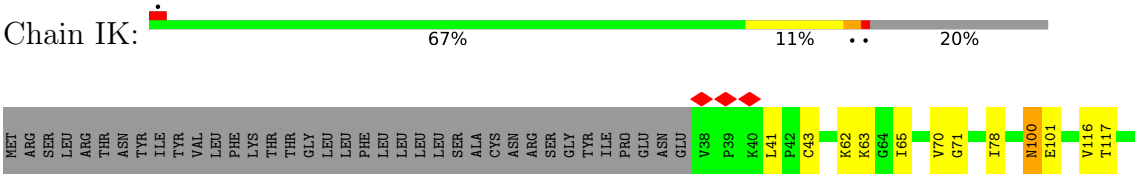
• Molecule 6: Inner membrane lipoprotein YiaD

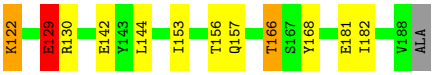
Chain HK:



• Molecule 6: Inner membrane lipoprotein YiaD

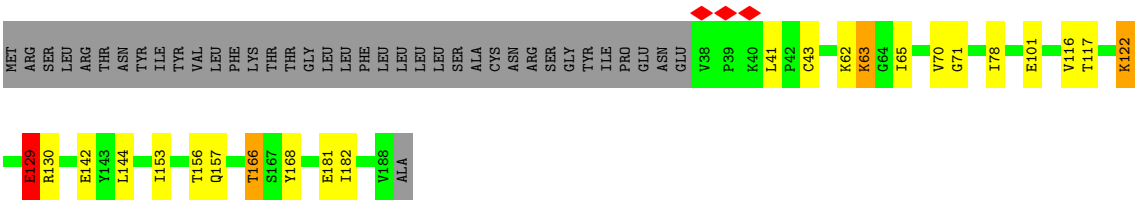
Chain IK:





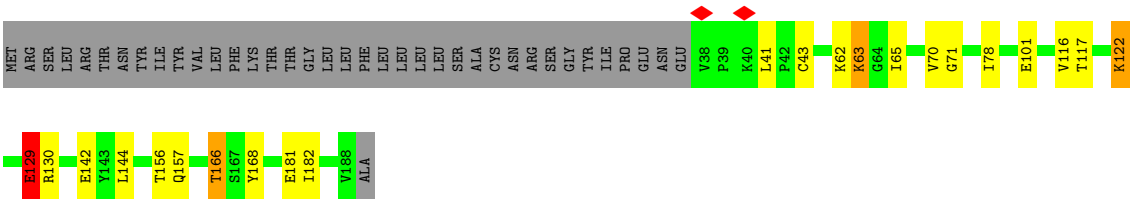
• Molecule 6: Inner membrane lipoprotein YiaD

Chain JK:



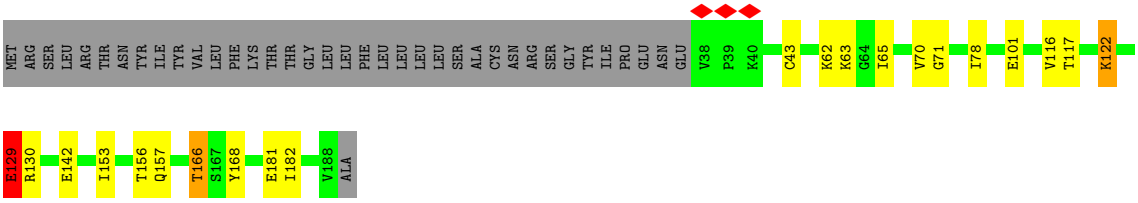
• Molecule 6: Inner membrane lipoprotein YiaD

Chain KK:



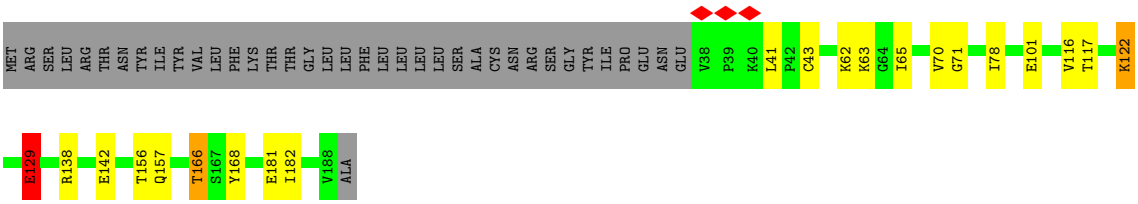
• Molecule 6: Inner membrane lipoprotein YiaD

Chain LK:



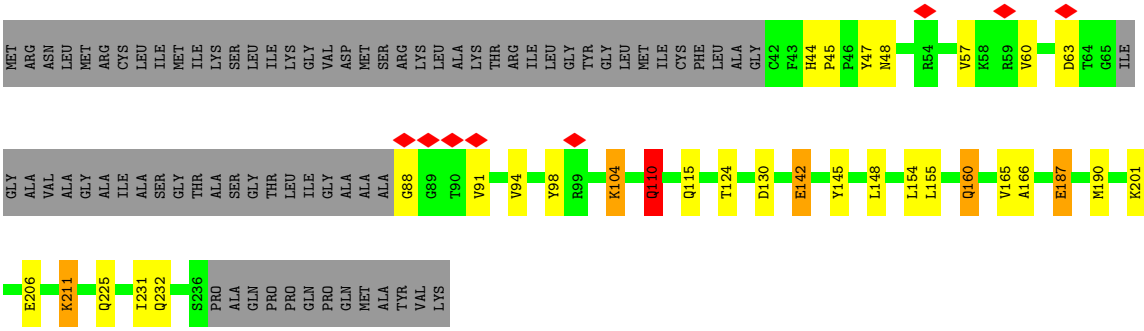
• Molecule 6: Inner membrane lipoprotein YiaD

Chain MK:



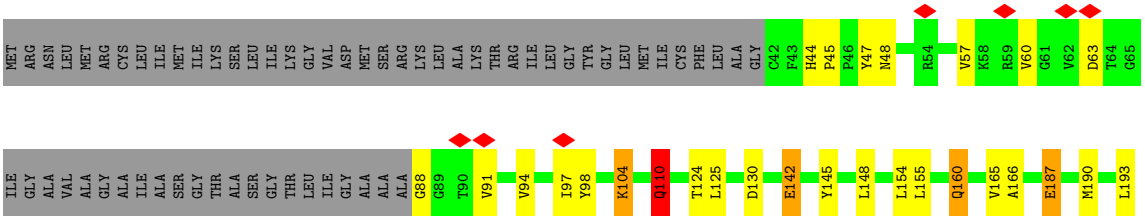
• Molecule 7: Outer membrane protein, OmpA family protein

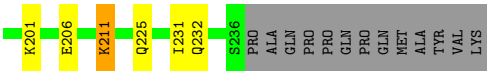
Chain AL:



• Molecule 7: Outer membrane protein, OmpA family protein

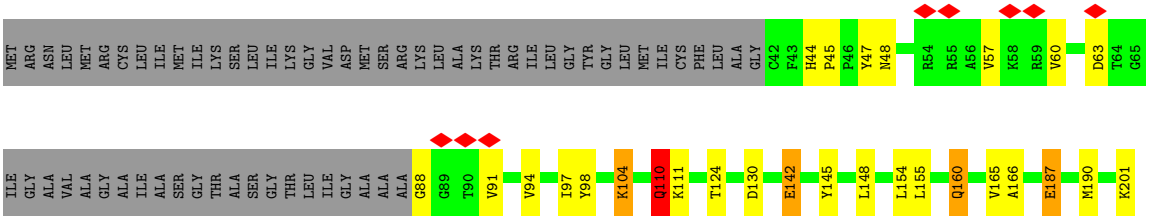
Chain BL:





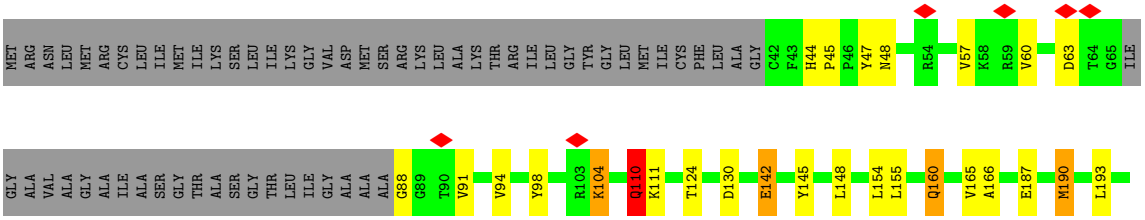
• Molecule 7: Outer membrane protein, OmpA family protein

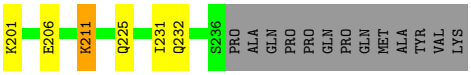
Chain CL:



• Molecule 7: Outer membrane protein, OmpA family protein

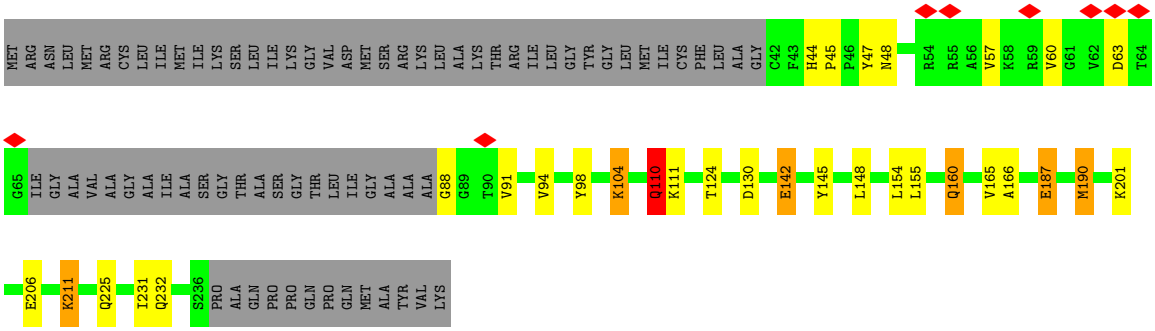
Chain DL:





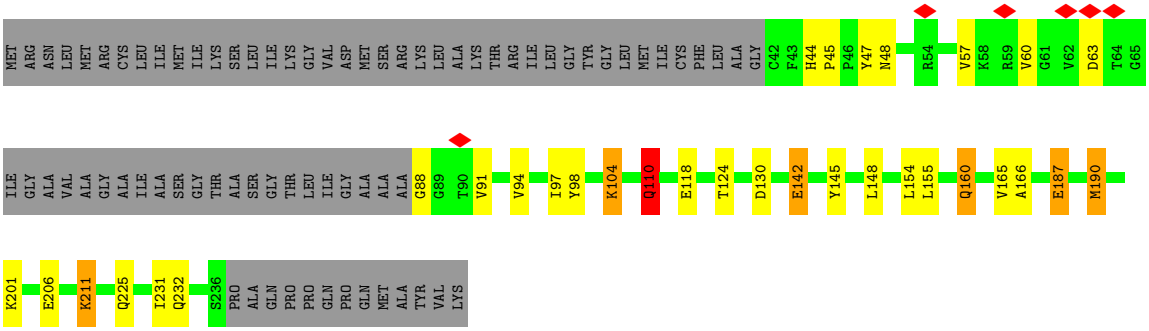
• Molecule 7: Outer membrane protein, OmpA family protein

Chain EL:



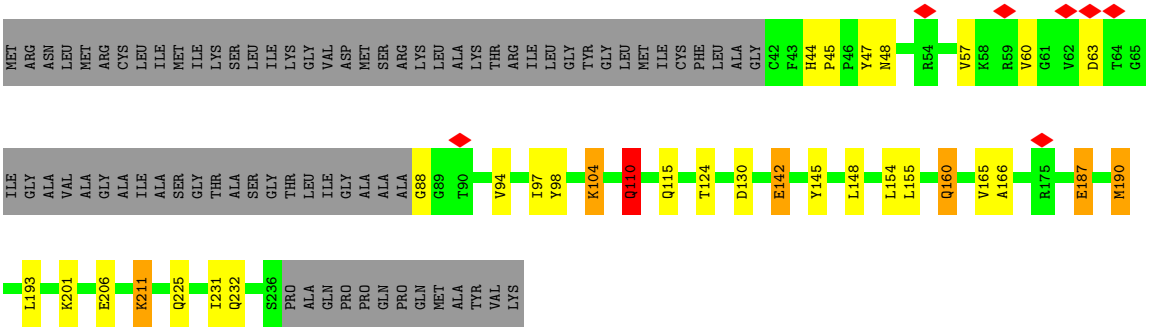
• Molecule 7: Outer membrane protein, OmpA family protein

Chain FL:



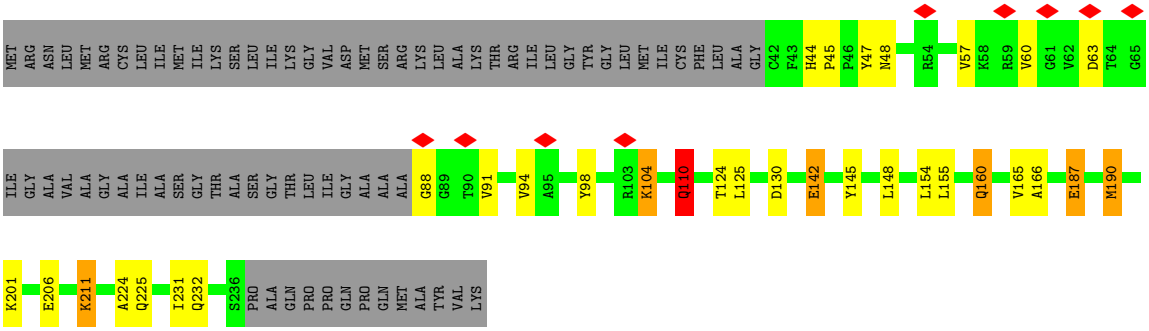
• Molecule 7: Outer membrane protein, OmpA family protein

Chain GL:



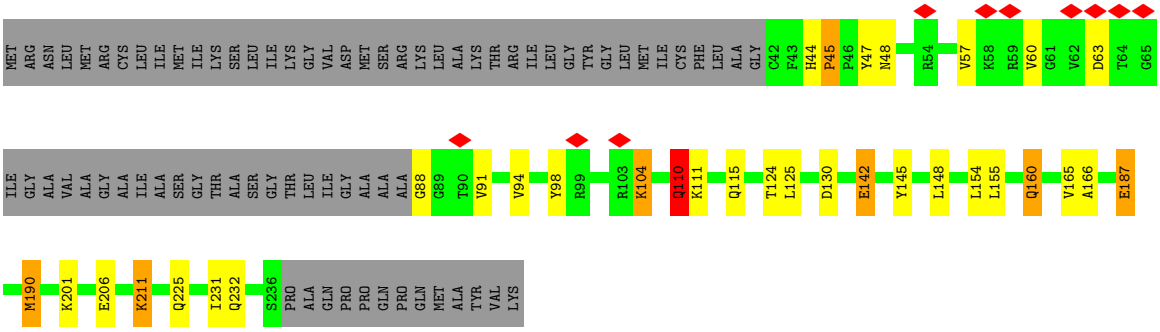
• Molecule 7: Outer membrane protein, OmpA family protein

Chain HL:



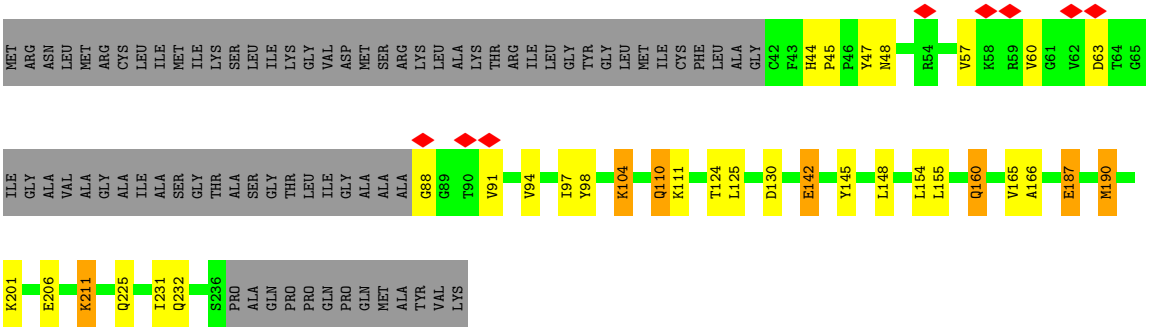
• Molecule 7: Outer membrane protein, OmpA family protein

Chain IL:



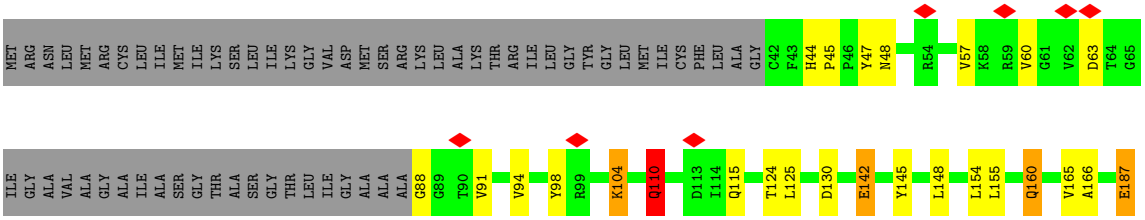
• Molecule 7: Outer membrane protein, OmpA family protein

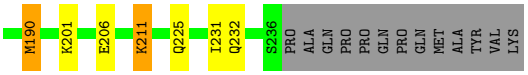
Chain JL:



• Molecule 7: Outer membrane protein, OmpA family protein

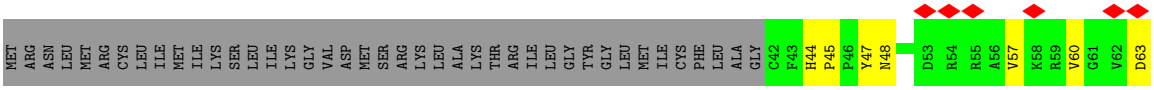
Chain KL:

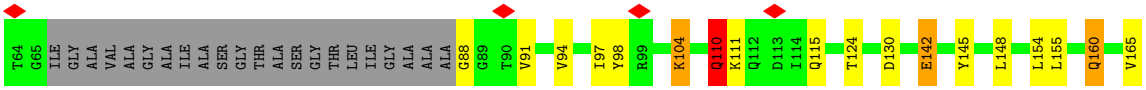


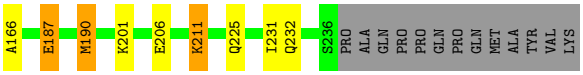


• Molecule 7: Outer membrane protein, OmpA family protein

Chain LL:

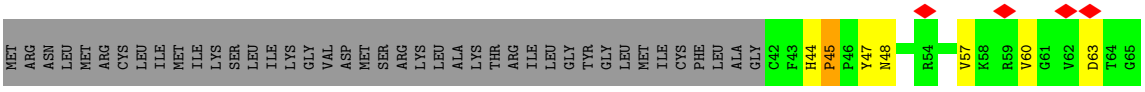


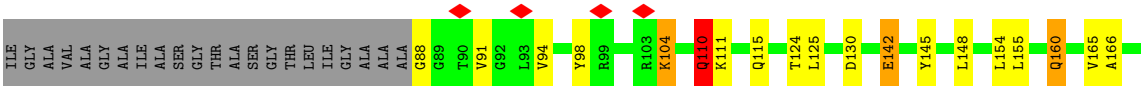


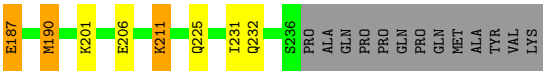


• Molecule 7: Outer membrane protein, OmpA family protein

Chain ML:

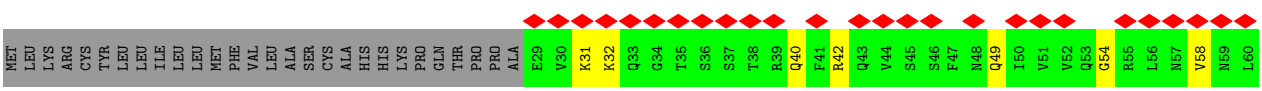


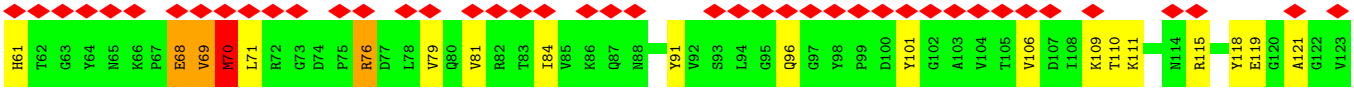


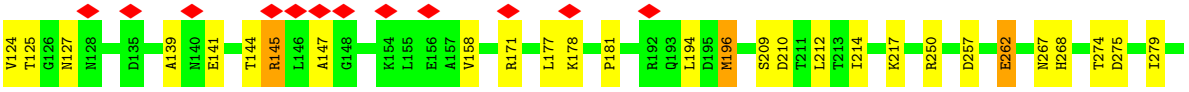


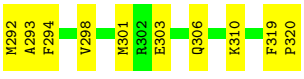
• Molecule 8: DUF2807 domain-containing protein

Chain AM:



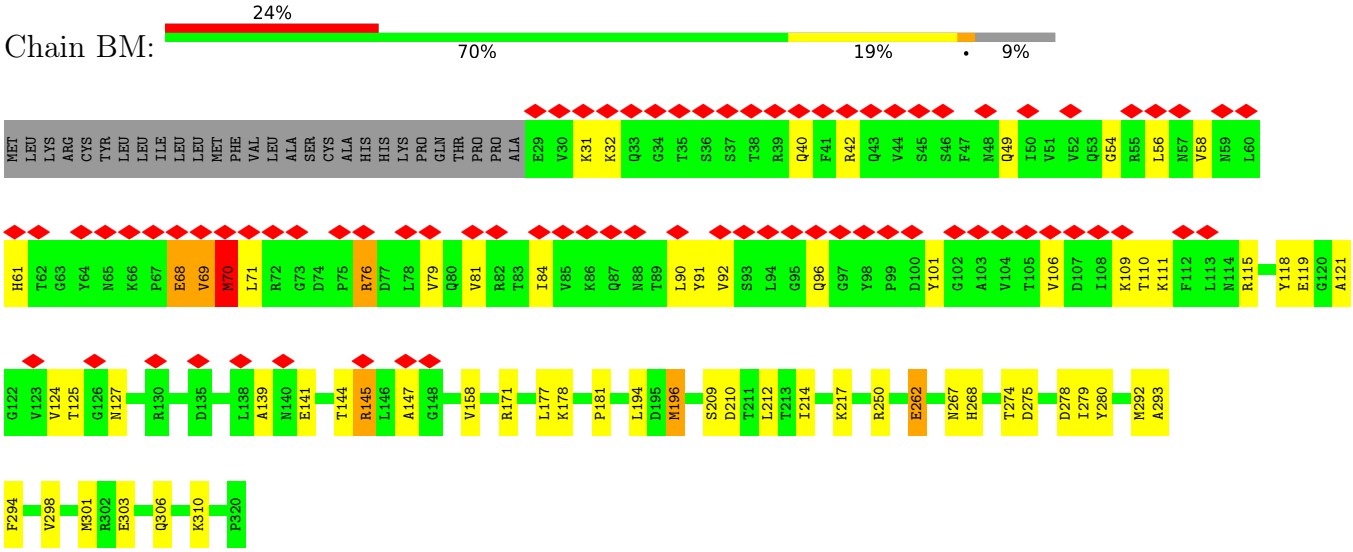






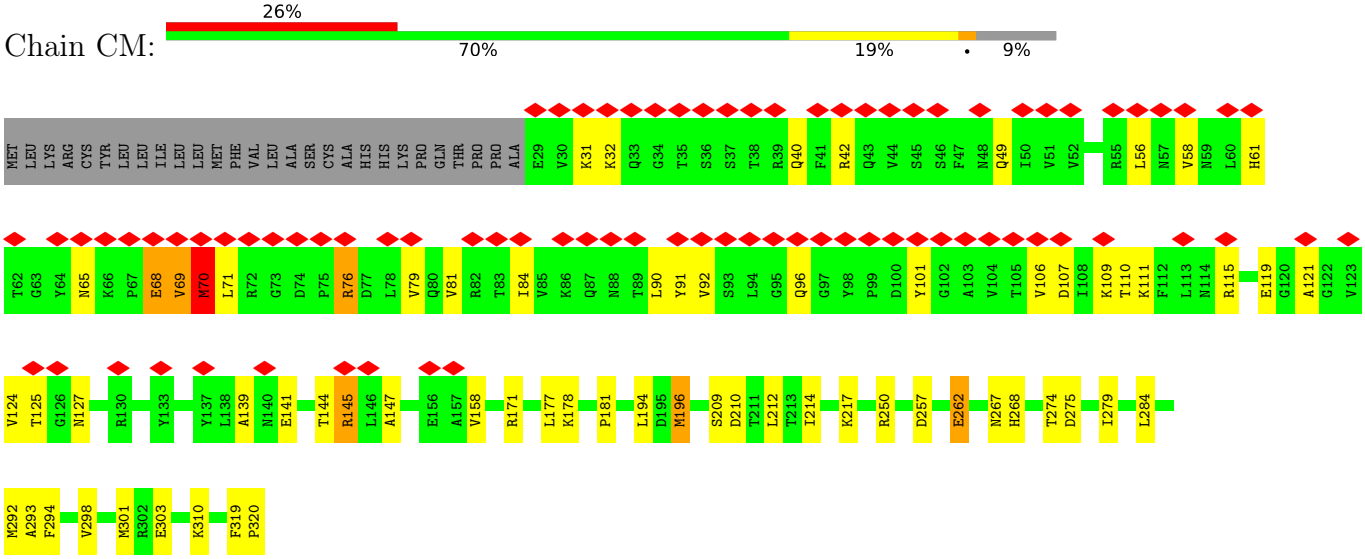
• Molecule 8: DUF2807 domain-containing protein

Chain BM:



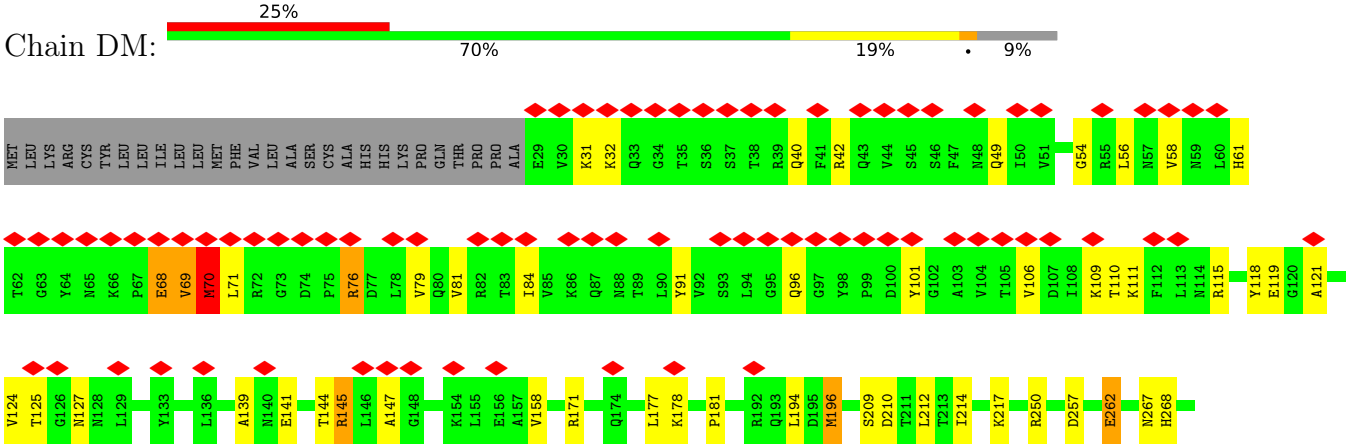
• Molecule 8: DUF2807 domain-containing protein

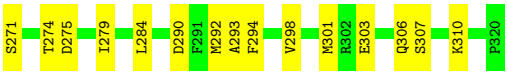
Chain CM:



• Molecule 8: DUF2807 domain-containing protein

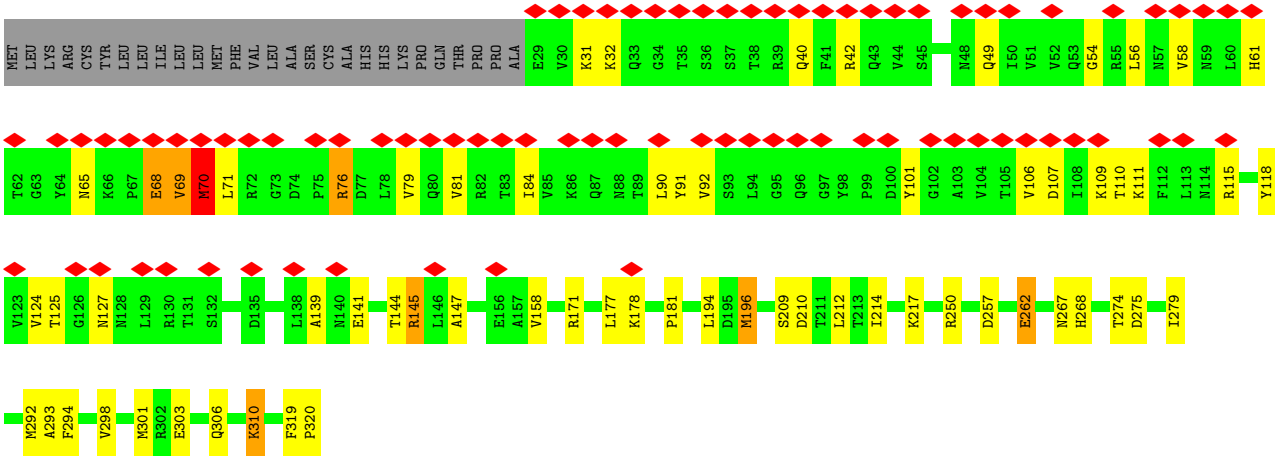
Chain DM:





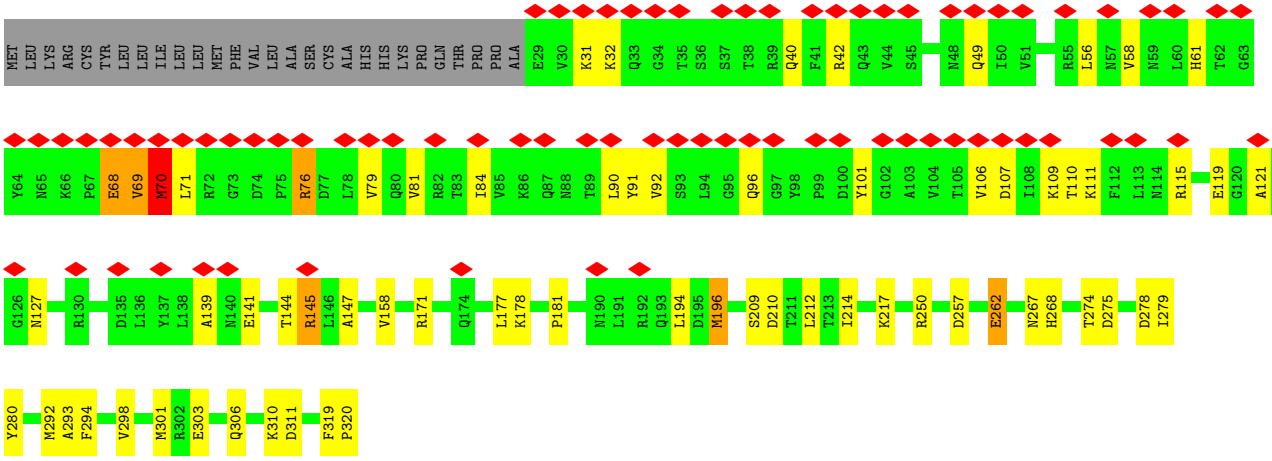
• Molecule 8: DUF2807 domain-containing protein

Chain EM:



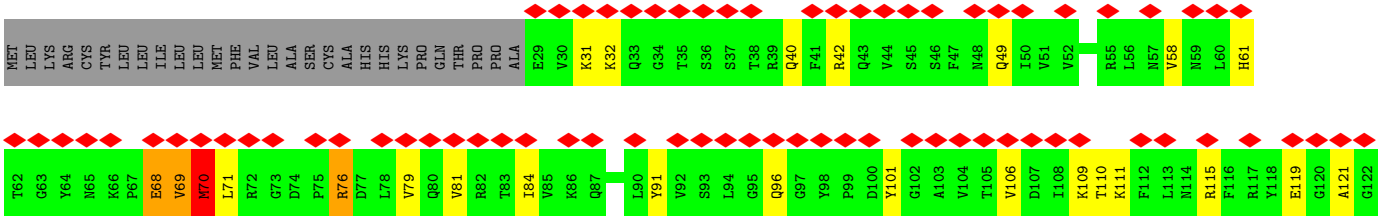
• Molecule 8: DUF2807 domain-containing protein

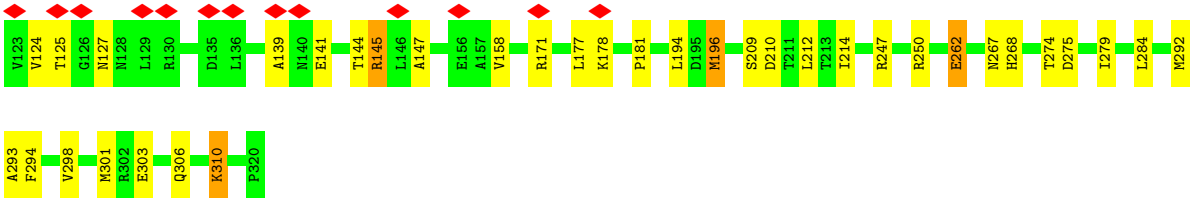
Chain FM:



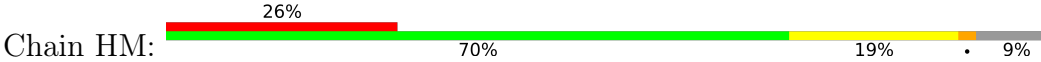
• Molecule 8: DUF2807 domain-containing protein

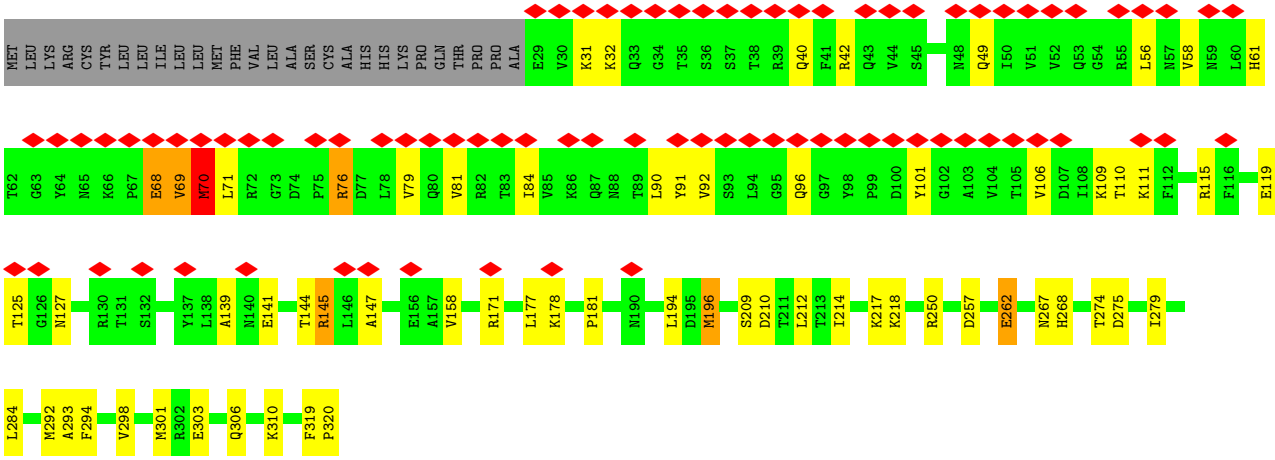
Chain GM:



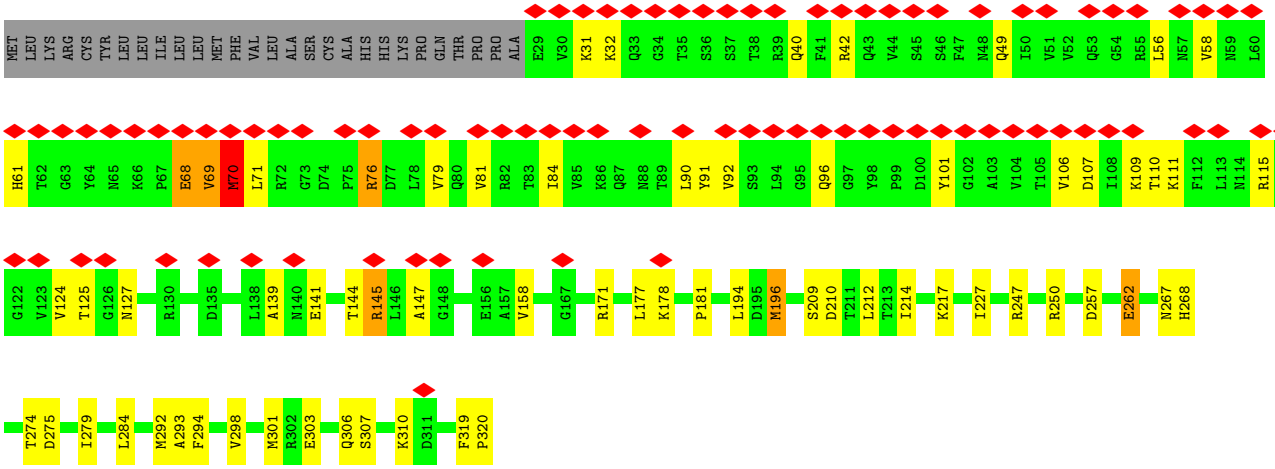


• Molecule 8: DUF2807 domain-containing protein

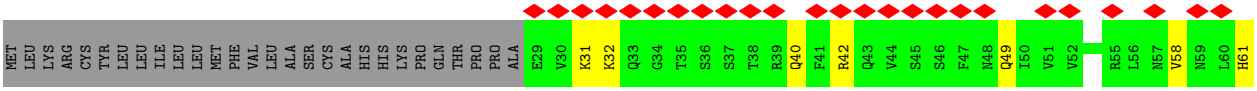


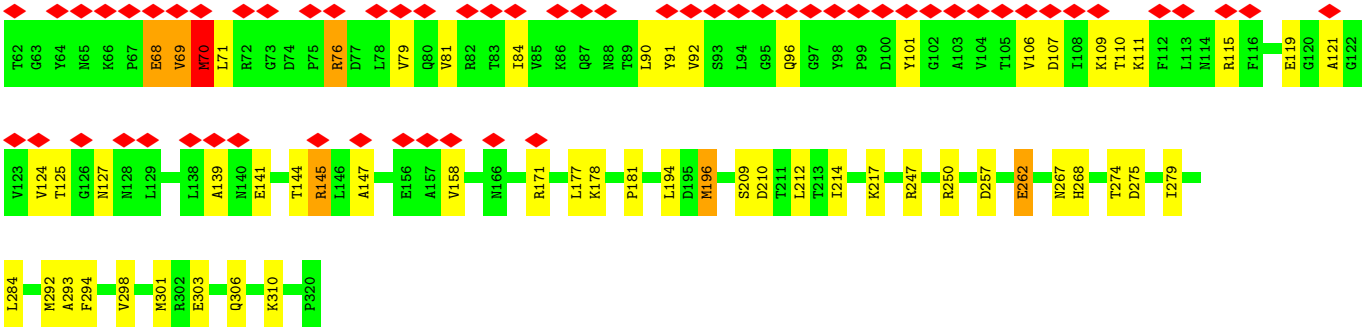


• Molecule 8: DUF2807 domain-containing protein



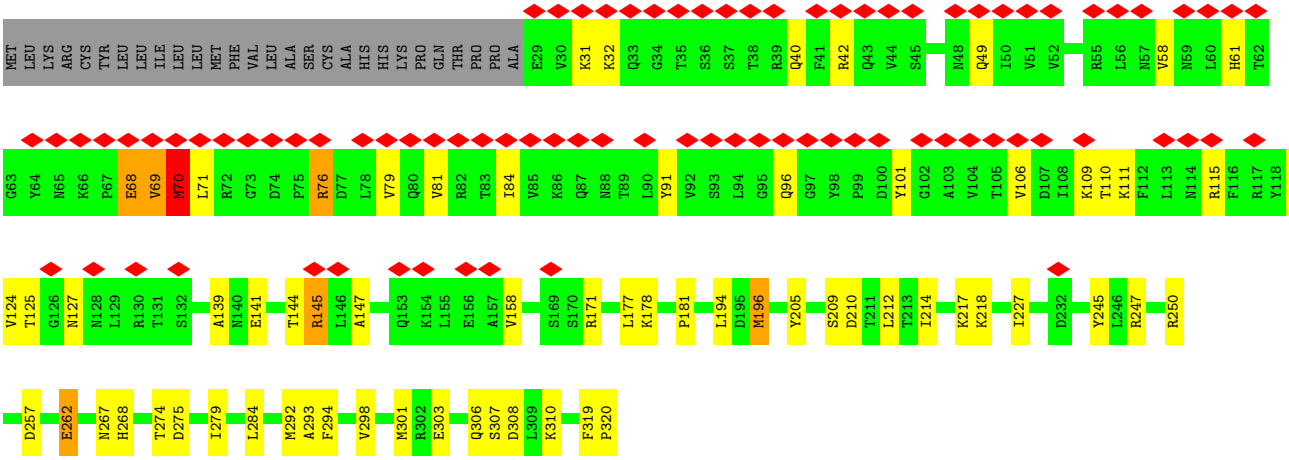
• Molecule 8: DUF2807 domain-containing protein





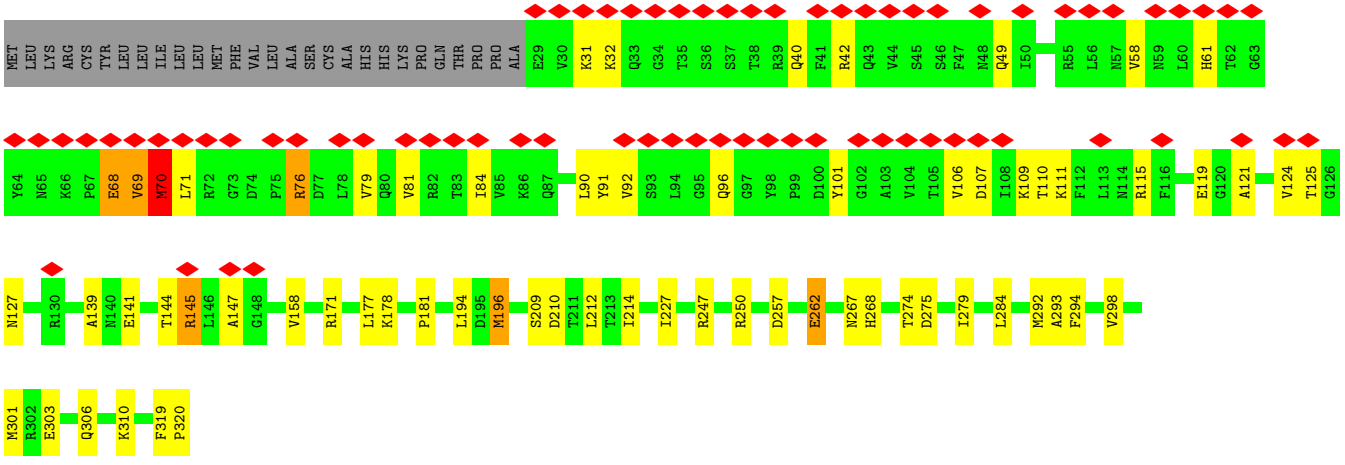
• Molecule 8: DUF2807 domain-containing protein

Chain KM:



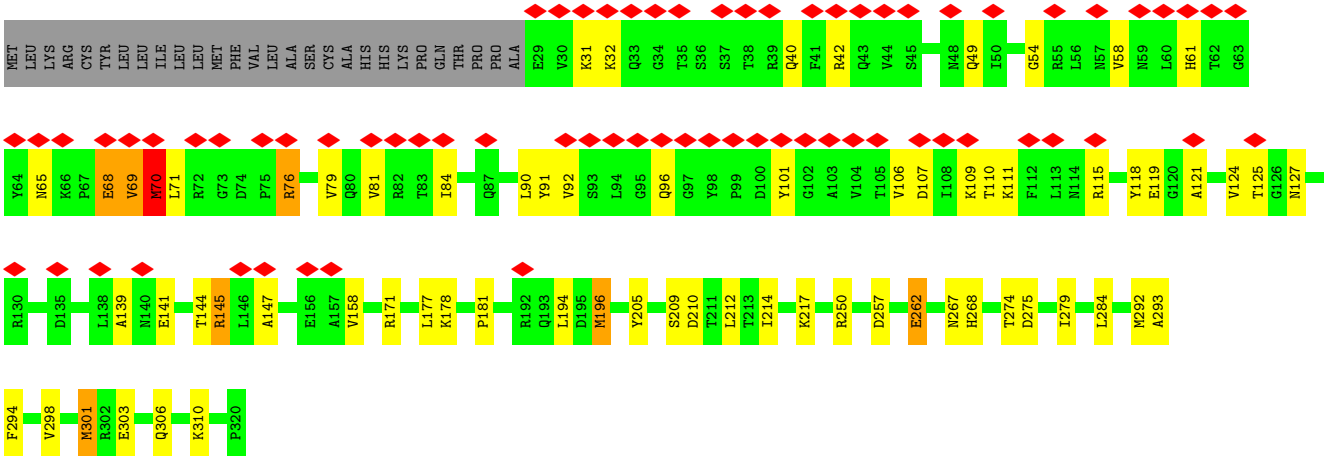
• Molecule 8: DUF2807 domain-containing protein

Chain LM:

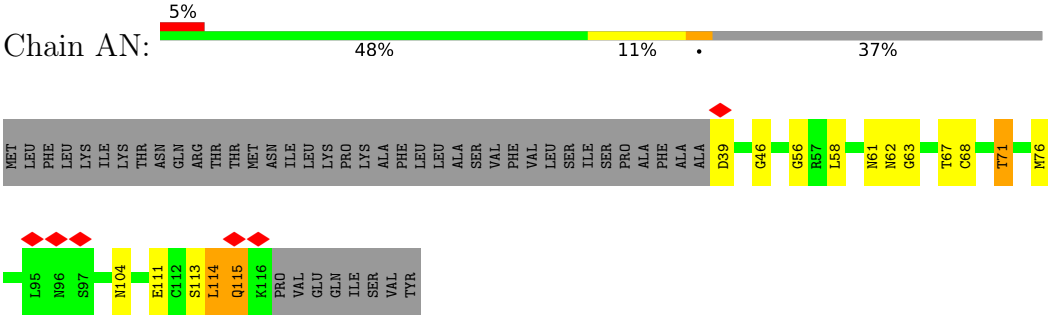


• Molecule 8: DUF2807 domain-containing protein

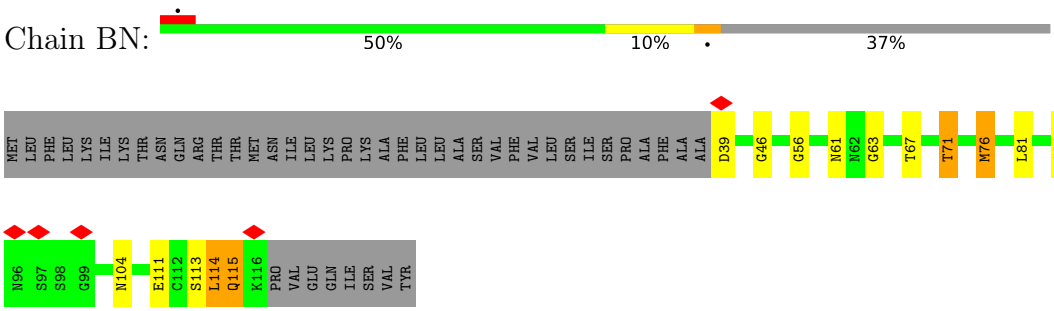
Chain MM:



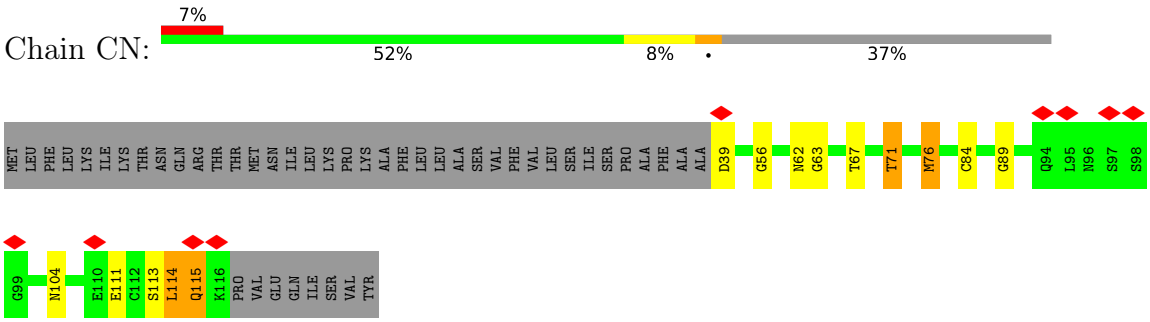
• Molecule 9: Neurogenic locus notch



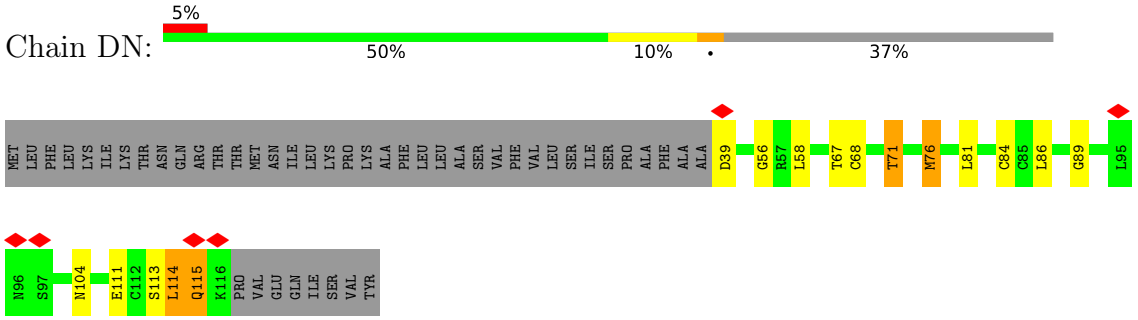
• Molecule 9: Neurogenic locus notch



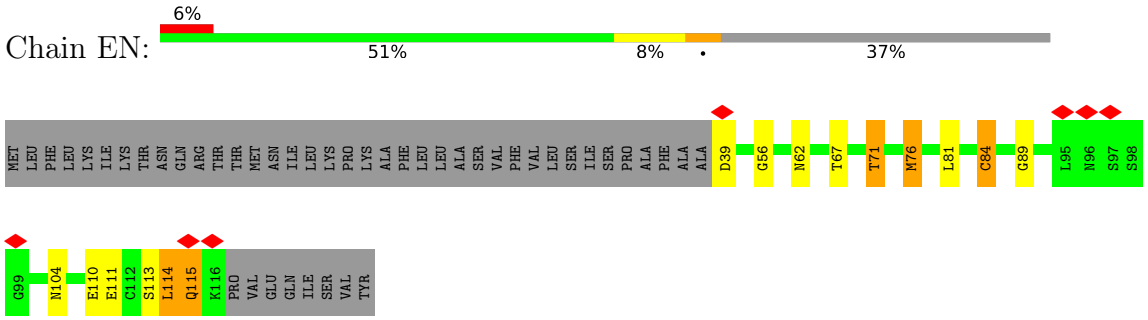
• Molecule 9: Neurogenic locus notch



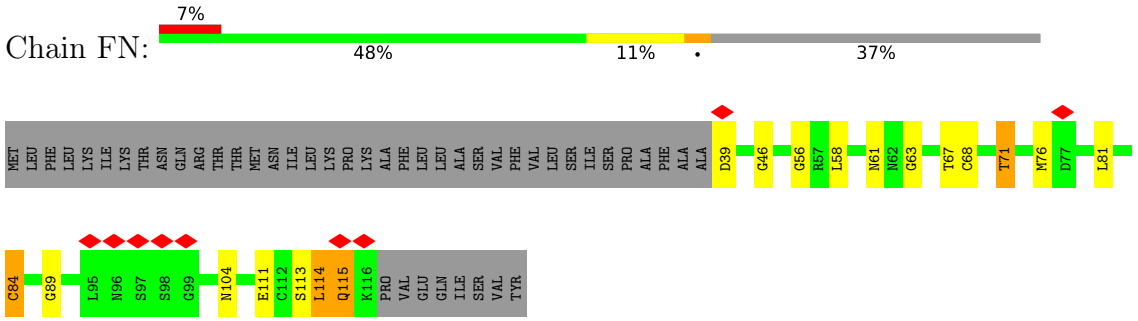
• Molecule 9: Neurogenic locus notch



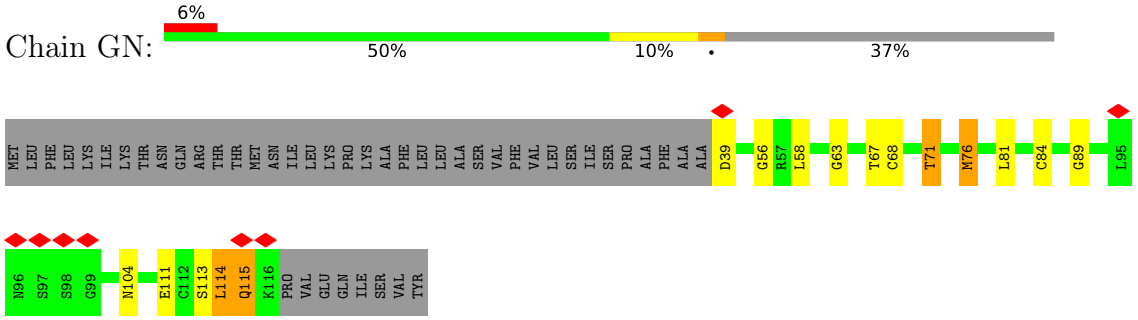
• Molecule 9: Neurogenic locus notch



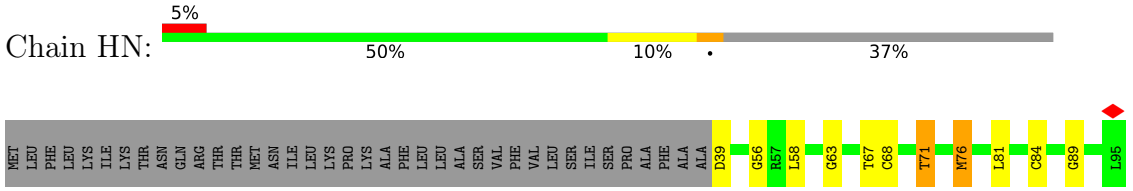
• Molecule 9: Neurogenic locus notch



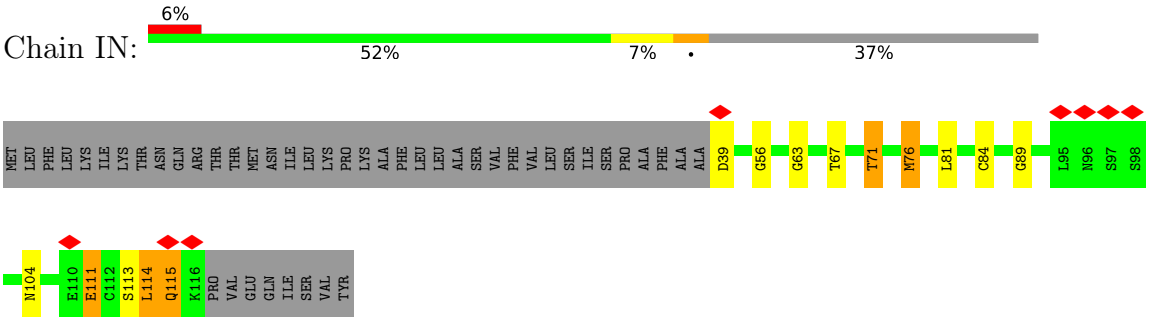
• Molecule 9: Neurogenic locus notch



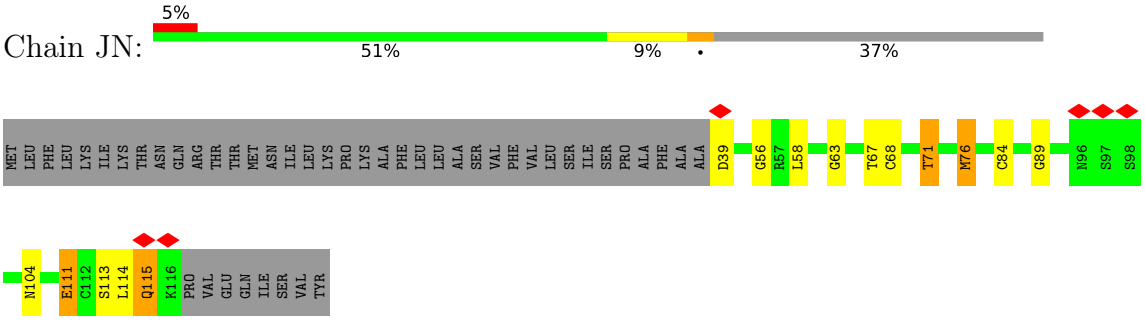
• Molecule 9: Neurogenic locus notch



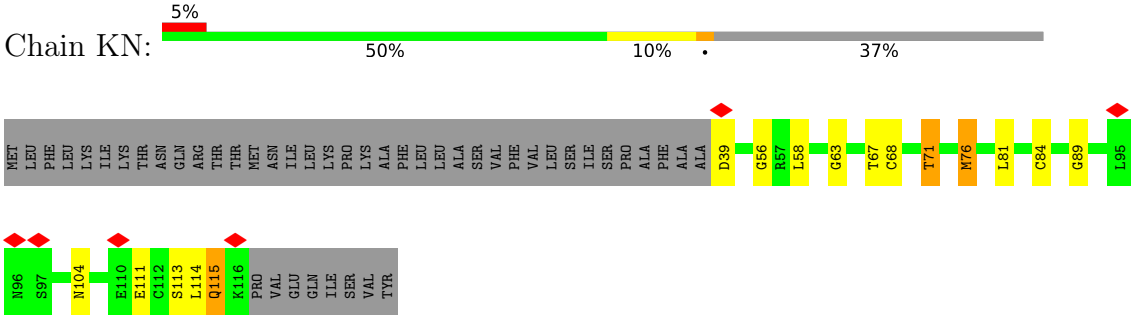
• Molecule 9: Neurogenic locus notch



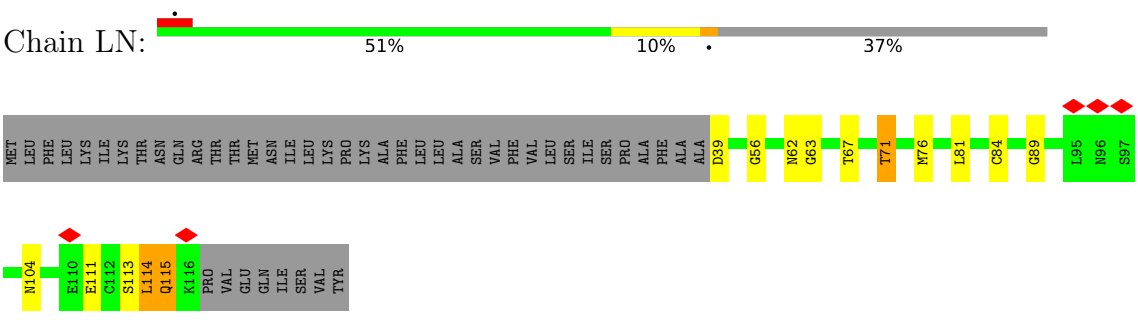
• Molecule 9: Neurogenic locus notch



• Molecule 9: Neurogenic locus notch

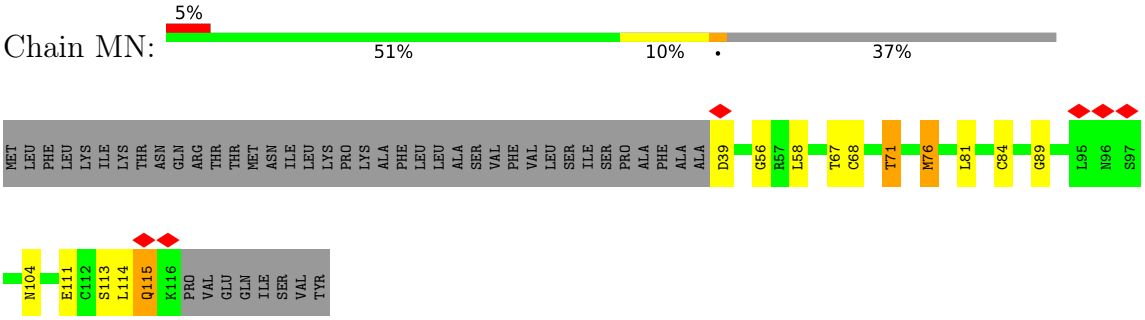


• Molecule 9: Neurogenic locus notch



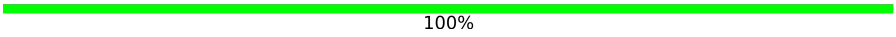
• Molecule 9: Neurogenic locus notch

Chain MN:



• Molecule 10: Unknown protein fragment

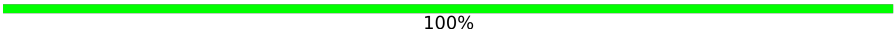
Chain AU:



There are no outlier residues recorded for this chain.

• Molecule 10: Unknown protein fragment

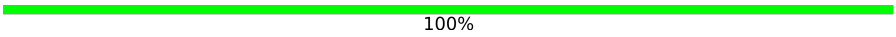
Chain BU:



There are no outlier residues recorded for this chain.

• Molecule 10: Unknown protein fragment

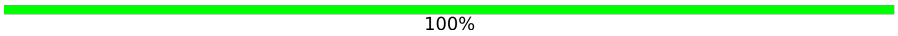
Chain CU:



There are no outlier residues recorded for this chain.

• Molecule 10: Unknown protein fragment

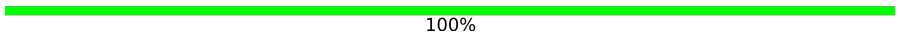
Chain DU:



There are no outlier residues recorded for this chain.

• Molecule 10: Unknown protein fragment

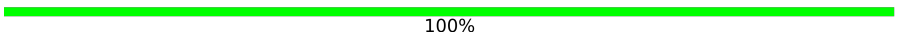
Chain EU:



There are no outlier residues recorded for this chain.

• Molecule 10: Unknown protein fragment

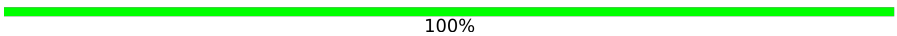
Chain FU:



There are no outlier residues recorded for this chain.

• Molecule 10: Unknown protein fragment

Chain GU:



There are no outlier residues recorded for this chain.

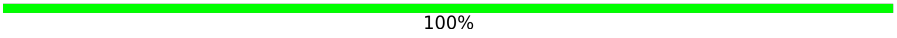
• Molecule 10: Unknown protein fragment

Chain HU:

There are no outlier residues recorded for this chain.

• Molecule 10: Unknown protein fragment

Chain IU:



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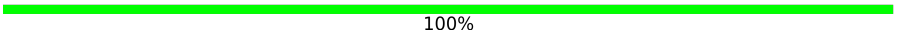
• Molecule 10: Unknown protein fragment

Chain JU:



There are no outlier residues recorded for this chain.

• Molecule 10: Unknown protein fragment

Chain KU:

There are no outlier residues recorded for this chain.

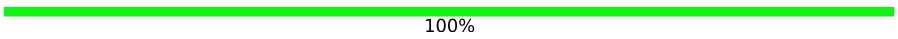
• Molecule 10: Unknown protein fragment

Chain LU:

There are no outlier residues recorded for this chain.

• Molecule 10: Unknown protein fragment

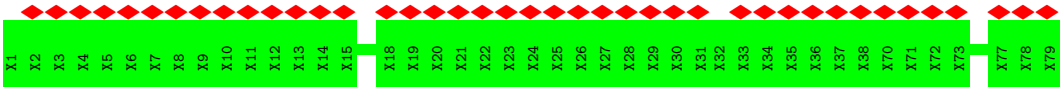
Chain MU:



There are no outlier residues recorded for this chain.

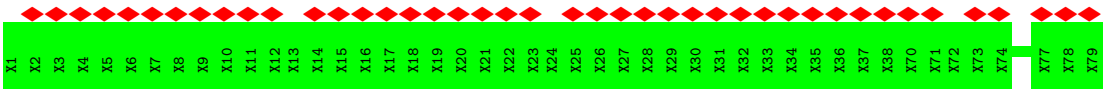
• Molecule 11: Unknown protein fragment

Chain AX:



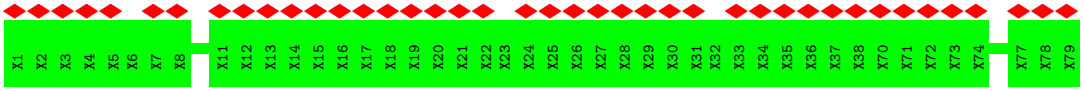
• Molecule 11: Unknown protein fragment

Chain BX:



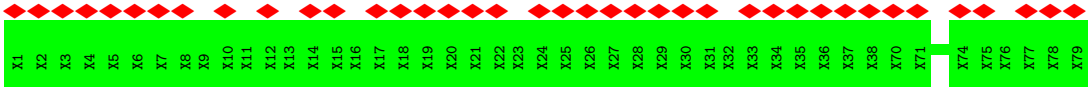
• Molecule 11: Unknown protein fragment

Chain CX:



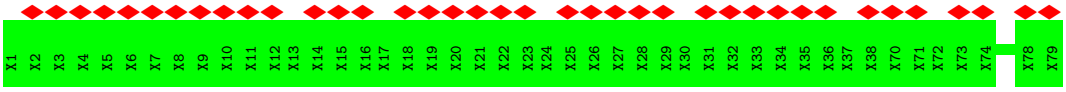
• Molecule 11: Unknown protein fragment

Chain DX:



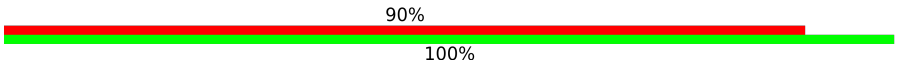
• Molecule 11: Unknown protein fragment

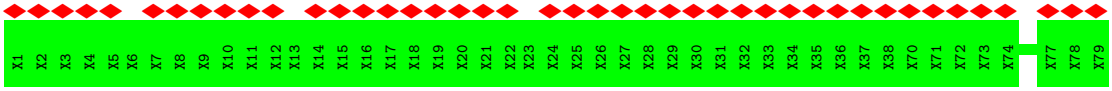
Chain EX:



• Molecule 11: Unknown protein fragment

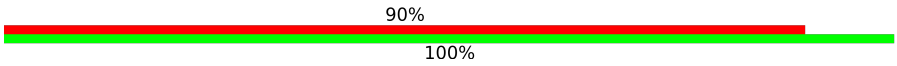
Chain FX:

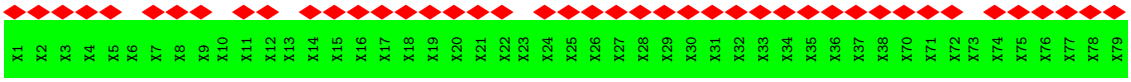




• Molecule 11: Unknown protein fragment

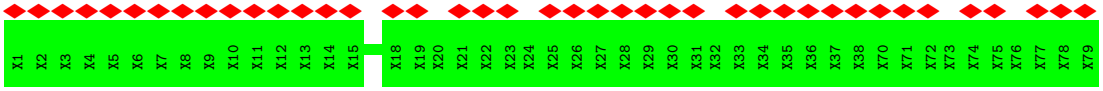
Chain GX:





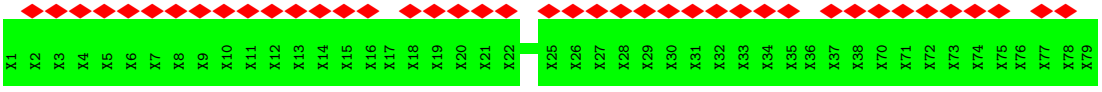
• Molecule 11: Unknown protein fragment

Chain HX:



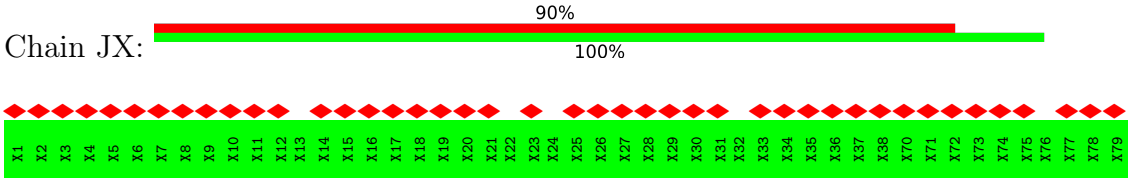
• Molecule 11: Unknown protein fragment

Chain IX:



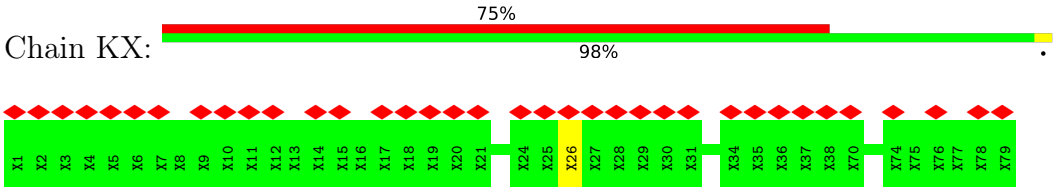
• Molecule 11: Unknown protein fragment

Chain JX:



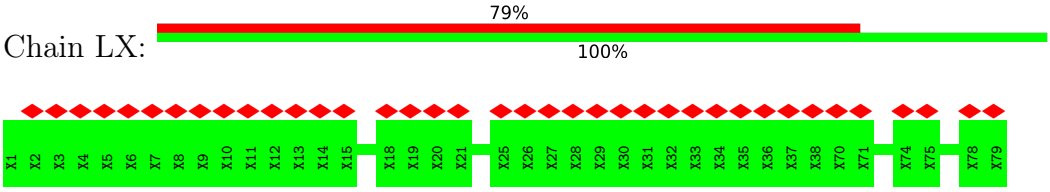
• Molecule 11: Unknown protein fragment

Chain KX:



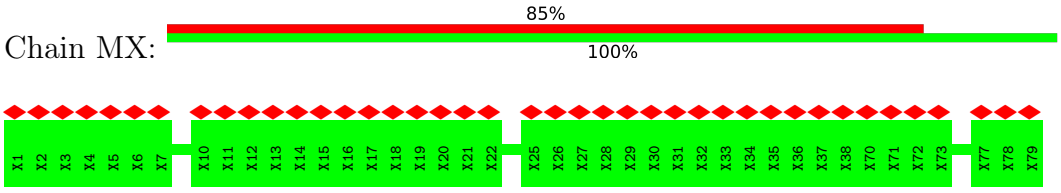
• Molecule 11: Unknown protein fragment

Chain LX:



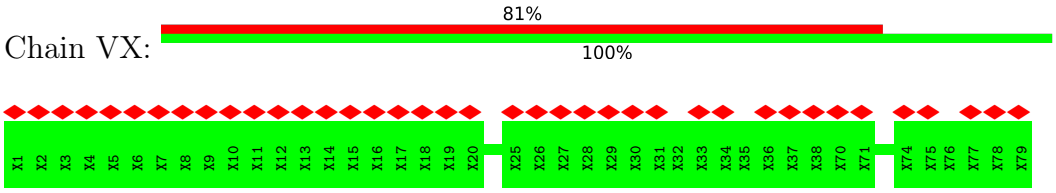
• Molecule 11: Unknown protein fragment

Chain MX:



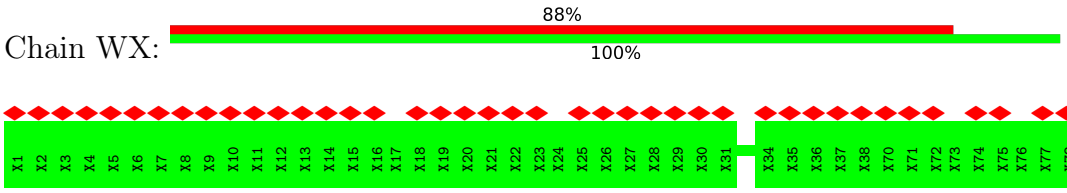
• Molecule 11: Unknown protein fragment

Chain VX:



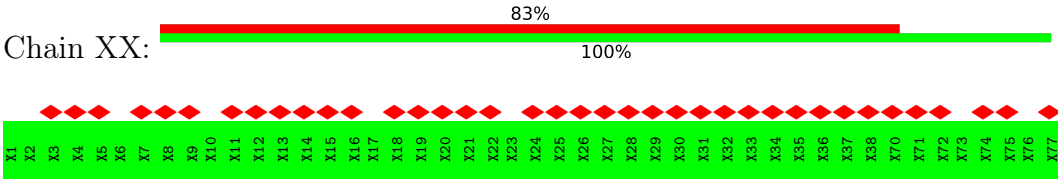
• Molecule 11: Unknown protein fragment

Chain WX:



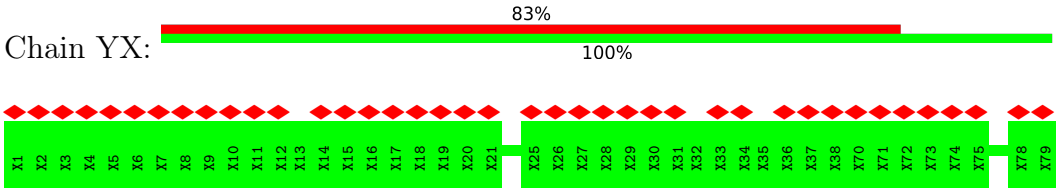
• Molecule 11: Unknown protein fragment

Chain XX:



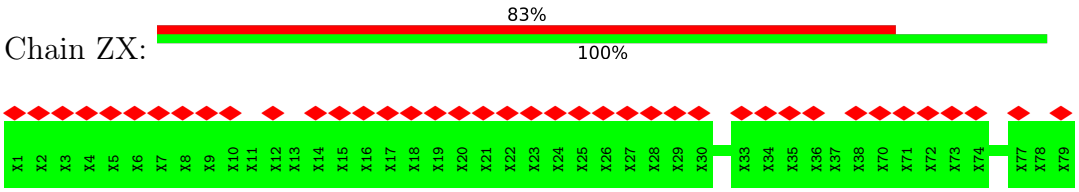
• Molecule 11: Unknown protein fragment

Chain YX:



• Molecule 11: Unknown protein fragment

Chain ZX:



4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	136818	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 ($e^-/Å^2$)	50	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	2.337	Depositor
Minimum map value	-1.017	Depositor
Average map value	0.015	Depositor
Map value standard deviation	0.106	Depositor
Recommended contour level	0.5	Depositor
Map size (Å)	561.0, 561.0, 561.0	wwPDB
Map dimensions	510, 510, 510	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.1, 1.1, 1.1	Depositor

5 Model quality

5.1 Standard geometry

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	#$\|Z\|$ >5	RMSZ	#$\|Z\|$ >5
1	AC	0.46	0/1717	0.90	8/2331 (0.3%)
1	BC	0.47	0/1717	0.90	8/2331 (0.3%)
1	CC	0.47	0/1717	0.90	8/2331 (0.3%)
1	DC	0.47	0/1717	0.91	8/2331 (0.3%)
1	EC	0.47	0/1717	0.90	8/2331 (0.3%)
1	FC	0.47	0/1717	0.90	8/2331 (0.3%)
1	GC	0.46	0/1717	0.90	8/2331 (0.3%)
1	HC	0.46	0/1717	0.90	8/2331 (0.3%)
1	IC	0.46	0/1717	0.90	8/2331 (0.3%)
1	JC	0.46	0/1717	0.90	8/2331 (0.3%)
1	KC	0.46	0/1717	0.90	8/2331 (0.3%)
1	LC	0.46	0/1717	0.90	8/2331 (0.3%)
1	MC	0.47	0/1717	0.90	8/2331 (0.3%)
2	AD	0.59	1/1107 (0.1%)	0.91	1/1502 (0.1%)
2	Ad	0.53	0/1078	0.93	2/1463 (0.1%)
2	BD	0.57	1/1107 (0.1%)	0.89	1/1502 (0.1%)
2	Bd	0.58	0/1078	0.90	3/1463 (0.2%)
2	CD	0.56	0/1107	0.86	3/1502 (0.2%)
2	Cd	0.61	1/1078 (0.1%)	1.04	5/1463 (0.3%)
2	DD	0.59	2/1107 (0.2%)	0.89	3/1502 (0.2%)
2	Dd	0.54	1/1078 (0.1%)	0.92	2/1463 (0.1%)
2	ED	0.60	0/1107	1.00	5/1502 (0.3%)
2	Ed	0.51	0/1078	0.87	0/1463
2	FD	0.50	0/1107	0.80	0/1502
2	Fd	1.03	8/1078 (0.7%)	1.13	13/1463 (0.9%)
2	GD	0.58	1/1107 (0.1%)	0.92	1/1502 (0.1%)
2	Gd	0.58	1/1078 (0.1%)	0.98	3/1463 (0.2%)
2	HD	0.61	1/1107 (0.1%)	0.93	2/1502 (0.1%)
2	Hd	0.56	1/1078 (0.1%)	0.84	0/1463
2	ID	0.54	0/1107	0.88	3/1502 (0.2%)
2	Id	0.53	0/1078	0.91	2/1463 (0.1%)
2	JD	0.52	1/1107 (0.1%)	0.88	1/1502 (0.1%)
2	Jd	0.54	0/1078	0.90	5/1463 (0.3%)
2	KD	0.57	2/1107 (0.2%)	0.88	3/1502 (0.2%)

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	#|Z| >5	RMSZ	#|Z| >5
2	Kd	0.54	0/1078	0.91	3/1463 (0.2%)
2	LD	0.60	1/1107 (0.1%)	0.94	5/1502 (0.3%)
2	Ld	0.55	0/1078	0.95	4/1463 (0.3%)
2	MD	0.57	1/1107 (0.1%)	0.84	0/1502
2	Md	0.59	0/1078	0.85	0/1463
3	AF	0.32	0/490	1.07	2/660 (0.3%)
3	Af	0.18	0/456	0.53	0/615
3	BF	0.32	0/490	1.07	2/660 (0.3%)
3	Bf	0.22	0/456	0.57	0/615
3	CF	0.32	0/490	1.07	2/660 (0.3%)
3	Cf	0.17	0/456	0.49	0/615
3	DF	0.31	0/490	1.07	2/660 (0.3%)
3	Df	0.18	0/456	0.53	0/615
3	EF	0.32	0/490	1.07	2/660 (0.3%)
3	Ef	0.19	0/456	0.54	0/615
3	FF	0.32	0/490	1.07	2/660 (0.3%)
3	Ff	0.20	0/456	0.52	0/615
3	GF	0.32	0/490	1.07	2/660 (0.3%)
3	Gf	0.16	0/456	0.38	0/615
3	HF	0.32	0/490	1.07	2/660 (0.3%)
3	Hf	0.19	0/456	0.58	0/615
3	IF	0.32	0/490	1.07	2/660 (0.3%)
3	If	0.20	0/456	0.57	0/615
3	JF	0.32	0/490	1.07	2/660 (0.3%)
3	Jf	0.17	0/456	0.51	0/615
3	KF	0.32	0/490	1.07	2/660 (0.3%)
3	Kf	0.20	0/456	0.49	0/615
3	LF	0.32	0/490	1.07	2/660 (0.3%)
3	Lf	0.17	0/456	0.43	0/615
3	MF	0.31	0/490	1.07	2/660 (0.3%)
3	Mf	0.18	0/456	0.55	0/615
3	VF	0.32	0/490	1.07	2/660 (0.3%)
3	WF	0.31	0/490	1.07	2/660 (0.3%)
3	XF	0.32	0/490	1.07	2/660 (0.3%)
3	YF	0.32	0/490	1.07	2/660 (0.3%)
3	ZF	0.31	0/490	1.07	2/660 (0.3%)
4	AG	0.82	1/278 (0.4%)	1.82	9/377 (2.4%)
4	BG	0.82	1/278 (0.4%)	1.82	9/377 (2.4%)
4	CG	0.82	1/278 (0.4%)	1.82	9/377 (2.4%)
4	DG	0.82	1/278 (0.4%)	1.82	9/377 (2.4%)
4	EG	0.81	1/278 (0.4%)	1.82	9/377 (2.4%)
4	FG	0.82	1/278 (0.4%)	1.82	9/377 (2.4%)
4	GG	0.82	1/278 (0.4%)	1.82	9/377 (2.4%)

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	#|Z| >5	RMSZ	#|Z| >5
4	HG	0.82	1/278 (0.4%)	1.82	9/377 (2.4%)
4	IG	0.82	1/278 (0.4%)	1.82	9/377 (2.4%)
4	JG	0.82	1/278 (0.4%)	1.82	9/377 (2.4%)
4	KG	0.81	1/278 (0.4%)	1.82	9/377 (2.4%)
4	LG	0.82	1/278 (0.4%)	1.82	9/377 (2.4%)
4	MG	0.82	1/278 (0.4%)	1.82	9/377 (2.4%)
4	VG	0.82	1/278 (0.4%)	1.82	9/377 (2.4%)
4	WG	0.81	1/278 (0.4%)	1.82	9/377 (2.4%)
4	XG	0.81	1/278 (0.4%)	1.82	9/377 (2.4%)
4	YG	0.81	1/278 (0.4%)	1.82	9/377 (2.4%)
4	ZG	0.82	1/278 (0.4%)	1.82	9/377 (2.4%)
5	AH	0.75	5/2033 (0.2%)	0.95	13/2775 (0.5%)
5	BH	0.56	3/2033 (0.1%)	0.92	10/2775 (0.4%)
5	CH	0.65	4/2033 (0.2%)	0.99	5/2775 (0.2%)
5	DH	0.66	4/2033 (0.2%)	0.94	5/2775 (0.2%)
5	EH	0.69	5/2033 (0.2%)	1.04	16/2775 (0.6%)
5	FH	0.66	3/2033 (0.1%)	0.96	10/2775 (0.4%)
5	GH	0.65	3/2033 (0.1%)	0.98	10/2775 (0.4%)
5	HH	0.67	5/2033 (0.2%)	1.00	12/2775 (0.4%)
5	IH	0.65	3/2033 (0.1%)	1.06	14/2775 (0.5%)
5	JH	0.58	2/2033 (0.1%)	0.99	11/2775 (0.4%)
5	KH	0.51	0/2033	0.92	6/2775 (0.2%)
5	LH	0.66	3/2033 (0.1%)	0.97	8/2775 (0.3%)
5	MH	0.59	0/2033	0.95	7/2775 (0.3%)
5	VH	0.45	1/1921 (0.1%)	0.92	6/2620 (0.2%)
5	WH	0.51	3/1921 (0.2%)	0.90	7/2620 (0.3%)
5	XH	0.53	4/1921 (0.2%)	0.97	6/2620 (0.2%)
5	YH	0.60	2/1921 (0.1%)	0.98	4/2620 (0.2%)
5	ZH	0.62	5/1921 (0.3%)	1.03	11/2620 (0.4%)
6	AK	0.48	0/1195	0.89	4/1616 (0.2%)
6	BK	0.48	0/1195	0.89	4/1616 (0.2%)
6	CK	0.48	0/1195	0.89	4/1616 (0.2%)
6	DK	0.48	0/1195	0.89	4/1616 (0.2%)
6	EK	0.48	0/1195	0.89	4/1616 (0.2%)
6	FK	0.48	0/1195	0.89	4/1616 (0.2%)
6	GK	0.48	0/1195	0.89	4/1616 (0.2%)
6	HK	0.48	0/1195	0.89	4/1616 (0.2%)
6	IK	0.48	0/1195	0.89	4/1616 (0.2%)
6	JK	0.48	0/1195	0.89	4/1616 (0.2%)
6	KK	0.48	0/1195	0.89	4/1616 (0.2%)
6	LK	0.48	0/1195	0.89	4/1616 (0.2%)
6	MK	0.48	0/1195	0.89	4/1616 (0.2%)
7	AL	0.48	0/1417	1.19	12/1912 (0.6%)

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	#|Z| >5	RMSZ	#|Z| >5
7	BL	0.47	0/1417	1.19	12/1912 (0.6%)
7	CL	0.47	0/1417	1.19	11/1912 (0.6%)
7	DL	0.48	0/1417	1.19	11/1912 (0.6%)
7	EL	0.47	0/1417	1.19	11/1912 (0.6%)
7	FL	0.47	0/1417	1.19	11/1912 (0.6%)
7	GL	0.47	0/1417	1.19	11/1912 (0.6%)
7	HL	0.48	0/1417	1.19	12/1912 (0.6%)
7	IL	0.48	0/1417	1.19	11/1912 (0.6%)
7	JL	0.47	0/1417	1.19	12/1912 (0.6%)
7	KL	0.47	0/1417	1.19	12/1912 (0.6%)
7	LL	0.48	0/1417	1.19	12/1912 (0.6%)
7	ML	0.47	0/1417	1.19	12/1912 (0.6%)
8	AM	0.42	0/2359	0.93	11/3183 (0.3%)
8	BM	0.42	0/2359	0.93	11/3183 (0.3%)
8	CM	0.42	0/2359	0.93	10/3183 (0.3%)
8	DM	0.42	0/2359	0.93	11/3183 (0.3%)
8	EM	0.42	0/2359	0.93	8/3183 (0.3%)
8	FM	0.42	0/2359	0.93	11/3183 (0.3%)
8	GM	0.42	0/2359	0.93	10/3183 (0.3%)
8	HM	0.42	0/2359	0.93	11/3183 (0.3%)
8	IM	0.42	0/2359	0.93	11/3183 (0.3%)
8	JM	0.42	0/2359	0.93	11/3183 (0.3%)
8	KM	0.42	0/2359	0.93	10/3183 (0.3%)
8	LM	0.42	0/2359	0.93	11/3183 (0.3%)
8	MM	0.42	0/2359	0.93	11/3183 (0.3%)
9	AN	0.45	0/593	1.00	3/799 (0.4%)
9	BN	0.45	0/593	1.01	3/799 (0.4%)
9	CN	0.45	0/593	1.01	3/799 (0.4%)
9	DN	0.45	0/593	1.00	3/799 (0.4%)
9	EN	0.45	0/593	1.01	3/799 (0.4%)
9	FN	0.45	0/593	1.00	3/799 (0.4%)
9	GN	0.46	0/593	1.00	3/799 (0.4%)
9	HN	0.45	0/593	1.00	3/799 (0.4%)
9	IN	0.45	0/593	1.00	3/799 (0.4%)
9	JN	0.45	0/593	1.00	3/799 (0.4%)
9	KN	0.45	0/593	1.01	3/799 (0.4%)
9	LN	0.45	0/593	1.00	3/799 (0.4%)
9	MN	0.45	0/593	1.00	3/799 (0.4%)
All	All	0.52	96/178844 (0.1%)	0.99	911/242314 (0.4%)

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	AD	0	1
2	BD	0	1
2	Bd	0	1
2	CD	0	1
2	Cd	0	1
2	DD	0	1
2	ED	0	1
2	FD	0	1
2	Fd	0	1
2	GD	0	1
2	ID	0	1
2	KD	0	1
2	MD	0	1
5	AH	0	2
5	BH	0	1
5	DH	0	1
5	EH	0	1
5	FH	0	2
5	GH	0	2
5	HH	0	1
5	IH	0	1
5	JH	0	3
5	KH	0	2
5	LH	0	1
5	MH	0	1
5	VH	0	1
5	WH	0	1
5	XH	0	1
5	YH	0	2
5	ZH	0	1
8	AM	0	1
8	BM	0	1
8	CM	0	1
8	DM	0	1
8	EM	0	1
8	FM	0	1
8	GM	0	1
8	HM	0	1
8	IM	0	1
8	JM	0	1
8	KM	0	1
8	LM	0	1

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Mol	Chain	#Chirality outliers	#Planarity outliers
8	MM	0	1
9	AN	0	1
9	BN	0	1
9	CN	0	1
9	DN	0	1
9	EN	0	1
9	FN	0	1
9	GN	0	1
9	HN	0	1
9	IN	0	1
9	JN	0	1
9	KN	0	1
9	LN	0	1
9	MN	0	1
All	All	0	63

All (96) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	AH	205	GLN	CG-CD	-16.68	1.10	1.52
2	Fd	82	ARG	CB-CG	13.11	1.91	1.52
5	IH	115	MET	CB-CG	-12.40	1.15	1.52
5	HH	115	MET	CB-CG	-11.80	1.17	1.52
5	DH	115	MET	CB-CG	-11.62	1.17	1.52
5	HH	115	MET	CG-SD	-11.41	1.52	1.80
5	YH	115	MET	CB-CG	-11.34	1.18	1.52
5	EH	285	GLU	CG-CD	-10.87	1.24	1.52
2	Fd	126	GLU	CB-CG	-10.49	1.21	1.52
5	DH	115	MET	CG-SD	-10.46	1.54	1.80
5	FH	285	GLU	CG-CD	-10.06	1.26	1.52
5	ZH	115	MET	CB-CG	-10.01	1.22	1.52
5	GH	115	MET	SD-CE	9.90	2.04	1.79
5	EH	285	GLU	CB-CG	-9.70	1.23	1.52
5	YH	115	MET	CG-SD	-9.59	1.56	1.80
5	BH	115	MET	SD-CE	-9.53	1.55	1.79
5	AH	205	GLN	CB-CG	-9.51	1.24	1.52
5	LH	115	MET	CG-SD	-9.26	1.57	1.80
2	Fd	82	ARG	CD-NE	-9.22	1.33	1.46
2	Fd	82	ARG	CG-CD	-8.92	1.25	1.52
5	XH	115	MET	CB-CG	-8.88	1.25	1.52
5	XH	115	MET	CG-SD	-8.87	1.58	1.80
5	WH	115	MET	CG-SD	-8.70	1.58	1.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	FH	285	GLU	CB-CG	-8.51	1.26	1.52
2	Fd	126	GLU	CA-C	8.49	1.62	1.52
5	JH	238	MET	CB-CG	-8.42	1.27	1.52
2	Fd	82	ARG	NE-CZ	-8.30	1.24	1.33
5	WH	115	MET	CB-CG	-8.27	1.27	1.52
5	DH	285	GLU	CG-CD	-8.26	1.31	1.52
5	LH	115	MET	CB-CG	-8.11	1.28	1.52
5	CH	115	MET	SD-CE	-7.92	1.59	1.79
5	EH	115	MET	CB-CG	-7.85	1.28	1.52
5	FH	115	MET	CB-CG	-7.75	1.29	1.52
5	CH	115	MET	CB-CG	-7.17	1.30	1.52
2	Fd	127	SER	N-CA	6.88	1.54	1.45
5	IH	115	MET	CA-CB	-6.74	1.42	1.53
5	EH	117	ARG	CB-CG	-6.73	1.32	1.52
5	ZH	115	MET	CG-SD	-6.68	1.64	1.80
5	AH	115	MET	CB-CG	-6.55	1.32	1.52
5	IH	115	MET	CA-C	-6.46	1.44	1.52
5	HH	115	MET	CA-CB	-6.41	1.43	1.53
5	ZH	242	ILE	CG1-CD1	-6.37	1.26	1.51
2	MD	26	LYS	C-N	-6.34	1.23	1.33
2	LD	52	MET	CB-CG	-6.23	1.33	1.52
5	BH	115	MET	CB-CG	-6.16	1.33	1.52
5	EH	117	ARG	CZ-NH1	6.09	1.41	1.32
5	GH	248	VAL	CB-CG1	-6.08	1.32	1.52
2	HD	26	LYS	C-N	-6.07	1.24	1.33
5	AH	205	GLN	CA-CB	-6.01	1.43	1.53
2	JD	26	LYS	C-N	-5.92	1.24	1.33
5	CH	151	PRO	CB-CG	-5.86	1.20	1.49
5	BH	107	ILE	CG1-CD1	-5.85	1.28	1.51
5	VH	115	MET	CG-SD	-5.80	1.66	1.80
2	KD	119	ILE	CG1-CD1	-5.74	1.29	1.51
5	HH	115	MET	CA-C	-5.73	1.45	1.52
5	AH	206	TRP	CB-CG	-5.69	1.32	1.50
2	Fd	125	ASP	CB-CG	5.68	1.66	1.52
4	LG	801	MET	SD-CE	-5.62	1.65	1.79
5	WH	238	MET	SD-CE	-5.62	1.65	1.79
2	AD	26	LYS	C-N	-5.60	1.24	1.33
4	IG	801	MET	SD-CE	-5.60	1.65	1.79
4	GG	801	MET	SD-CE	-5.60	1.65	1.79
4	CG	801	MET	SD-CE	-5.59	1.65	1.79
4	BG	801	MET	SD-CE	-5.58	1.65	1.79
4	DG	801	MET	SD-CE	-5.58	1.65	1.79

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	FG	801	MET	SD-CE	-5.56	1.65	1.79
4	KG	801	MET	SD-CE	-5.56	1.65	1.79
2	Cd	43	GLU	CB-CG	-5.55	1.35	1.52
4	AG	801	MET	SD-CE	-5.55	1.65	1.79
4	ZG	801	MET	SD-CE	-5.55	1.65	1.79
4	MG	801	MET	SD-CE	-5.55	1.65	1.79
4	VG	801	MET	SD-CE	-5.55	1.65	1.79
4	HG	801	MET	SD-CE	-5.54	1.65	1.79
4	JG	801	MET	SD-CE	-5.54	1.65	1.79
4	WG	801	MET	SD-CE	-5.54	1.65	1.79
4	YG	801	MET	SD-CE	-5.54	1.65	1.79
4	EG	801	MET	SD-CE	-5.53	1.65	1.79
4	XG	801	MET	SD-CE	-5.53	1.65	1.79
5	LH	132	GLN	CB-CG	-5.52	1.35	1.52
2	DD	55	MET	SD-CE	-5.50	1.65	1.79
5	DH	285	GLU	CB-CG	-5.49	1.35	1.52
5	JH	115	MET	CB-CG	-5.46	1.36	1.52
5	ZH	130	LEU	CG-CD2	-5.45	1.34	1.52
5	CH	250	TYR	CD2-CE2	-5.39	1.22	1.38
2	KD	26	LYS	C-N	-5.38	1.25	1.33
2	BD	55	MET	SD-CE	-5.32	1.66	1.79
2	GD	26	LYS	C-N	-5.31	1.25	1.33
2	Dd	26	LYS	C-N	-5.30	1.25	1.33
5	XH	115	MET	SD-CE	-5.27	1.66	1.79
5	XH	115	MET	CA-CB	-5.24	1.45	1.53
2	Hd	36	ASP	CB-CG	-5.17	1.39	1.52
5	ZH	208	LYS	CB-CG	-5.14	1.37	1.52
5	GH	115	MET	CG-SD	-5.12	1.68	1.80
5	HH	207	ASP	CB-CG	-5.11	1.39	1.52
2	DD	52	MET	CB-CG	-5.09	1.37	1.52
2	Gd	82	ARG	CB-CG	5.08	1.67	1.52

All (911) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
7	EL	187	GLU	CG-CD-OE2	20.68	165.97	118.40
7	BL	187	GLU	CG-CD-OE2	20.67	165.95	118.40
7	HL	187	GLU	CG-CD-OE2	20.67	165.94	118.40
7	AL	187	GLU	CG-CD-OE2	20.66	165.92	118.40
7	GL	187	GLU	CG-CD-OE2	20.65	165.89	118.40
7	ML	187	GLU	CG-CD-OE2	20.64	165.87	118.40
7	CL	187	GLU	CG-CD-OE2	20.64	165.87	118.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
7	LL	187	GLU	CG-CD-OE2	20.64	165.87	118.40
7	JL	187	GLU	CG-CD-OE2	20.64	165.86	118.40
7	IL	187	GLU	CG-CD-OE2	20.63	165.85	118.40
7	DL	187	GLU	CG-CD-OE2	20.63	165.84	118.40
7	FL	187	GLU	CG-CD-OE2	20.62	165.84	118.40
7	KL	187	GLU	CG-CD-OE2	20.62	165.82	118.40
5	FH	115	MET	CB-CG-SD	-18.56	57.03	112.70
5	YH	115	MET	CG-SD-CE	-16.36	64.91	100.90
7	KL	187	GLU	OE1-CD-OE2	-15.84	84.90	122.90
7	DL	187	GLU	OE1-CD-OE2	-15.82	84.94	122.90
7	FL	187	GLU	OE1-CD-OE2	-15.81	84.95	122.90
7	AL	187	GLU	OE1-CD-OE2	-15.81	84.96	122.90
7	IL	187	GLU	OE1-CD-OE2	-15.81	84.95	122.90
7	CL	187	GLU	OE1-CD-OE2	-15.81	84.96	122.90
7	JL	187	GLU	OE1-CD-OE2	-15.80	84.97	122.90
7	LL	187	GLU	OE1-CD-OE2	-15.80	84.99	122.90
7	GL	187	GLU	OE1-CD-OE2	-15.79	85.00	122.90
7	EL	187	GLU	OE1-CD-OE2	-15.78	85.02	122.90
7	HL	187	GLU	OE1-CD-OE2	-15.78	85.02	122.90
7	ML	187	GLU	OE1-CD-OE2	-15.78	85.02	122.90
7	BL	187	GLU	OE1-CD-OE2	-15.77	85.06	122.90
5	HH	115	MET	CG-SD-CE	-15.61	66.56	100.90
5	IH	115	MET	CG-SD-CE	-15.56	66.67	100.90
5	JH	238	MET	CG-SD-CE	-15.55	66.69	100.90
5	ZH	115	MET	CG-SD-CE	-15.20	67.47	100.90
5	XH	115	MET	CG-SD-CE	-15.00	67.90	100.90
5	DH	115	MET	CG-SD-CE	-14.93	68.05	100.90
2	ED	160	LYS	CD-CE-NZ	-14.72	64.80	111.90
7	IL	187	GLU	CG-CD-OE1	-13.91	86.42	118.40
7	ML	187	GLU	CG-CD-OE1	-13.90	86.42	118.40
7	BL	187	GLU	CG-CD-OE1	-13.90	86.43	118.40
7	FL	187	GLU	CG-CD-OE1	-13.89	86.45	118.40
7	LL	187	GLU	CG-CD-OE1	-13.89	86.46	118.40
7	CL	187	GLU	CG-CD-OE1	-13.89	86.46	118.40
7	GL	187	GLU	CG-CD-OE1	-13.88	86.47	118.40
7	JL	187	GLU	CG-CD-OE1	-13.88	86.47	118.40
7	AL	187	GLU	CG-CD-OE1	-13.88	86.48	118.40
7	DL	187	GLU	CG-CD-OE1	-13.88	86.48	118.40
7	HL	187	GLU	CG-CD-OE1	-13.87	86.49	118.40
5	IH	115	MET	CB-CG-SD	-13.87	71.11	112.70
7	KL	187	GLU	CG-CD-OE1	-13.87	86.51	118.40
7	EL	187	GLU	CG-CD-OE1	-13.86	86.52	118.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
5	CH	115	MET	CA-CB-CG	-13.65	86.80	114.10
5	KH	238	MET	CG-SD-CE	-13.57	71.05	100.90
5	CH	115	MET	CG-SD-CE	-13.52	71.17	100.90
5	LH	115	MET	CG-SD-CE	-13.31	71.62	100.90
5	HH	115	MET	CA-CB-CG	-12.96	88.19	114.10
5	EH	115	MET	CG-SD-CE	-12.43	73.55	100.90
5	VH	115	MET	CB-CG-SD	-11.93	76.90	112.70
4	BG	801	MET	CA-CB-CG	11.83	137.76	114.10
4	AG	801	MET	CA-CB-CG	11.80	137.70	114.10
4	GG	801	MET	CA-CB-CG	11.80	137.70	114.10
4	HG	801	MET	CA-CB-CG	11.80	137.69	114.10
4	CG	801	MET	CA-CB-CG	11.79	137.69	114.10
4	IG	801	MET	CA-CB-CG	11.79	137.69	114.10
4	DG	801	MET	CA-CB-CG	11.79	137.68	114.10
4	VG	801	MET	CA-CB-CG	11.79	137.68	114.10
4	LG	801	MET	CA-CB-CG	11.79	137.67	114.10
4	EG	801	MET	CA-CB-CG	11.78	137.66	114.10
4	JG	801	MET	CA-CB-CG	11.78	137.66	114.10
4	XG	801	MET	CA-CB-CG	11.78	137.66	114.10
4	YG	801	MET	CA-CB-CG	11.78	137.65	114.10
4	FG	801	MET	CA-CB-CG	11.77	137.65	114.10
4	WG	801	MET	CA-CB-CG	11.77	137.65	114.10
4	ZG	801	MET	CA-CB-CG	11.77	137.63	114.10
4	MG	801	MET	CA-CB-CG	11.77	137.63	114.10
4	KG	801	MET	CA-CB-CG	11.76	137.62	114.10
5	DH	115	MET	CA-CB-CG	-11.69	90.71	114.10
5	WH	115	MET	CG-SD-CE	-11.61	75.37	100.90
5	XH	238	MET	CG-SD-CE	-11.57	75.44	100.90
5	EH	208	LYS	CA-CB-CG	11.22	136.54	114.10
5	YH	287	VAL	N-CA-C	-11.13	99.05	108.63
5	GH	115	MET	CB-CG-SD	-11.03	79.61	112.70
5	XH	115	MET	CB-CG-SD	-11.01	79.68	112.70
5	GH	238	MET	CG-SD-CE	-10.73	77.29	100.90
2	Ld	55	MET	CG-SD-CE	-10.71	77.33	100.90
4	FG	796	GLN	CA-CB-CG	10.46	135.01	114.10
4	XG	796	GLN	CA-CB-CG	10.44	134.99	114.10
4	YG	796	GLN	CA-CB-CG	10.44	134.99	114.10
4	VG	796	GLN	CA-CB-CG	10.44	134.98	114.10
4	CG	796	GLN	CA-CB-CG	10.44	134.97	114.10
4	ZG	796	GLN	CA-CB-CG	10.44	134.97	114.10
4	HG	796	GLN	CA-CB-CG	10.44	134.97	114.10
4	AG	796	GLN	CA-CB-CG	10.43	134.96	114.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	EG	796	GLN	CA-CB-CG	10.42	134.94	114.10
4	JG	796	GLN	CA-CB-CG	10.41	134.93	114.10
4	LG	796	GLN	CA-CB-CG	10.41	134.93	114.10
4	MG	796	GLN	CA-CB-CG	10.41	134.92	114.10
4	KG	796	GLN	CA-CB-CG	10.41	134.92	114.10
4	DG	796	GLN	CA-CB-CG	10.40	134.91	114.10
4	IG	796	GLN	CA-CB-CG	10.40	134.91	114.10
4	GG	796	GLN	CA-CB-CG	10.40	134.89	114.10
4	WG	796	GLN	CA-CB-CG	10.39	134.88	114.10
4	BG	796	GLN	CA-CB-CG	10.39	134.88	114.10
2	Fd	82	ARG	CB-CG-CD	-10.28	87.66	111.30
2	Fd	82	ARG	CA-CB-CG	9.98	134.06	114.10
2	Ad	55	MET	CG-SD-CE	-9.85	79.22	100.90
5	BH	115	MET	CG-SD-CE	-9.81	79.33	100.90
5	JH	115	MET	CG-SD-CE	-9.71	79.53	100.90
5	EH	285	GLU	CG-CD-OE2	-9.64	96.23	118.40
5	EH	117	ARG	N-CA-CB	-9.58	95.43	110.28
2	Fd	26	LYS	CD-CE-NZ	-9.55	81.33	111.90
5	FH	285	GLU	CG-CD-OE2	-9.28	97.06	118.40
5	HH	238	MET	CG-SD-CE	-9.28	80.49	100.90
1	MC	123	GLN	CA-CB-CG	9.14	132.39	114.10
1	HC	123	GLN	CA-CB-CG	9.14	132.38	114.10
1	LC	123	GLN	CA-CB-CG	9.14	132.38	114.10
1	FC	123	GLN	CA-CB-CG	9.13	132.37	114.10
1	DC	123	GLN	CA-CB-CG	9.13	132.36	114.10
1	IC	123	GLN	CA-CB-CG	9.12	132.34	114.10
1	CC	123	GLN	CA-CB-CG	9.12	132.34	114.10
1	GC	123	GLN	CA-CB-CG	9.12	132.34	114.10
1	EC	123	GLN	CA-CB-CG	9.12	132.34	114.10
1	AC	123	GLN	CA-CB-CG	9.11	132.32	114.10
1	KC	123	GLN	CA-CB-CG	9.11	132.32	114.10
1	JC	123	GLN	CA-CB-CG	9.11	132.31	114.10
1	BC	123	GLN	CA-CB-CG	9.10	132.31	114.10
1	MC	60	LYS	CA-CB-CG	8.98	132.06	114.10
1	LC	60	LYS	CA-CB-CG	8.98	132.06	114.10
5	VH	287	VAL	N-CA-C	-8.98	100.91	108.63
1	AC	60	LYS	CA-CB-CG	8.97	132.05	114.10
1	GC	60	LYS	CA-CB-CG	8.97	132.04	114.10
7	AL	211	LYS	CB-CG-CD	8.97	131.92	111.30
7	FL	211	LYS	CB-CG-CD	8.97	131.93	111.30
7	EL	211	LYS	CB-CG-CD	8.96	131.92	111.30
1	HC	60	LYS	CA-CB-CG	8.96	132.03	114.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
7	CL	211	LYS	CB-CG-CD	8.96	131.91	111.30
1	CC	60	LYS	CA-CB-CG	8.96	132.02	114.10
7	HL	211	LYS	CB-CG-CD	8.96	131.90	111.30
1	IC	60	LYS	CA-CB-CG	8.96	132.02	114.10
1	KC	60	LYS	CA-CB-CG	8.96	132.01	114.10
7	KL	211	LYS	CB-CG-CD	8.95	131.89	111.30
1	JC	60	LYS	CA-CB-CG	8.95	132.00	114.10
7	BL	211	LYS	CB-CG-CD	8.95	131.88	111.30
7	DL	211	LYS	CB-CG-CD	8.95	131.88	111.30
7	IL	211	LYS	CB-CG-CD	8.94	131.87	111.30
1	BC	60	LYS	CA-CB-CG	8.94	131.99	114.10
1	EC	60	LYS	CA-CB-CG	8.94	131.98	114.10
7	ML	211	LYS	CB-CG-CD	8.94	131.86	111.30
7	GL	211	LYS	CB-CG-CD	8.94	131.85	111.30
7	LL	211	LYS	CB-CG-CD	8.94	131.85	111.30
1	FC	60	LYS	CA-CB-CG	8.92	131.95	114.10
7	JL	211	LYS	CB-CG-CD	8.92	131.82	111.30
1	DC	60	LYS	CA-CB-CG	8.92	131.94	114.10
5	EH	117	ARG	CD-NE-CZ	-8.82	112.05	124.40
2	Fd	82	ARG	CG-CD-NE	-8.73	92.78	112.00
5	DH	285	GLU	CA-CB-CG	-8.57	96.95	114.10
7	CL	110	GLN	CA-CB-CG	8.44	130.98	114.10
7	FL	110	GLN	CA-CB-CG	8.42	130.94	114.10
7	IL	110	GLN	CA-CB-CG	8.41	130.93	114.10
7	GL	110	GLN	CA-CB-CG	8.41	130.93	114.10
7	EL	110	GLN	CA-CB-CG	8.41	130.92	114.10
7	BL	110	GLN	CA-CB-CG	8.39	130.89	114.10
7	LL	110	GLN	CA-CB-CG	8.39	130.88	114.10
7	HL	110	GLN	CA-CB-CG	8.39	130.87	114.10
7	JL	110	GLN	CA-CB-CG	8.39	130.87	114.10
7	ML	110	GLN	CA-CB-CG	8.39	130.87	114.10
7	AL	110	GLN	CA-CB-CG	8.38	130.86	114.10
7	KL	110	GLN	CA-CB-CG	8.38	130.86	114.10
7	DL	110	GLN	CA-CB-CG	8.38	130.86	114.10
5	HH	115	MET	CB-CG-SD	-8.36	87.64	112.70
5	IH	150	LYS	CB-CG-CD	8.28	130.35	111.30
5	AH	205	GLN	CA-CB-CG	-8.28	97.54	114.10
5	IH	208	LYS	N-CA-CB	-8.27	97.72	110.46
4	BG	813	LYS	CA-CB-CG	8.24	130.59	114.10
4	XG	813	LYS	CA-CB-CG	8.24	130.59	114.10
4	MG	813	LYS	CA-CB-CG	8.24	130.58	114.10
4	GG	813	LYS	CA-CB-CG	8.23	130.57	114.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	EG	813	LYS	CA-CB-CG	8.23	130.56	114.10
4	FG	813	LYS	CA-CB-CG	8.23	130.56	114.10
4	LG	813	LYS	CA-CB-CG	8.23	130.56	114.10
4	AG	813	LYS	CA-CB-CG	8.22	130.55	114.10
4	HG	813	LYS	CA-CB-CG	8.22	130.55	114.10
4	WG	813	LYS	CA-CB-CG	8.22	130.55	114.10
4	YG	813	LYS	CA-CB-CG	8.22	130.55	114.10
4	VG	813	LYS	CA-CB-CG	8.22	130.54	114.10
4	JG	813	LYS	CA-CB-CG	8.22	130.53	114.10
4	ZG	813	LYS	CA-CB-CG	8.21	130.51	114.10
4	CG	813	LYS	CA-CB-CG	8.21	130.51	114.10
4	KG	813	LYS	CA-CB-CG	8.21	130.51	114.10
4	DG	813	LYS	CA-CB-CG	8.20	130.50	114.10
4	IG	813	LYS	CA-CB-CG	8.20	130.49	114.10
5	EH	115	MET	CA-CB-CG	-8.18	97.75	114.10
5	IH	115	MET	N-CA-C	8.12	120.21	111.36
5	EH	285	GLU	CA-CB-CG	-8.03	98.04	114.10
5	IH	150	LYS	CA-CB-CG	7.97	130.04	114.10
5	YH	115	MET	CA-CB-CG	-7.93	98.23	114.10
5	AH	115	MET	CG-SD-CE	-7.90	83.52	100.90
8	DM	196	MET	CA-CB-CG	7.76	129.62	114.10
8	FM	196	MET	CA-CB-CG	7.75	129.60	114.10
8	BM	196	MET	CA-CB-CG	7.75	129.60	114.10
2	Cd	52	MET	CA-CB-CG	7.75	129.59	114.10
8	GM	196	MET	CA-CB-CG	7.75	129.59	114.10
8	CM	196	MET	CA-CB-CG	7.74	129.59	114.10
8	IM	196	MET	CA-CB-CG	7.74	129.59	114.10
8	LM	196	MET	CA-CB-CG	7.74	129.57	114.10
8	MM	196	MET	CA-CB-CG	7.73	129.56	114.10
8	AM	196	MET	CA-CB-CG	7.73	129.56	114.10
8	HM	196	MET	CA-CB-CG	7.72	129.54	114.10
1	EC	60	LYS	CB-CG-CD	7.72	129.05	111.30
3	HF	267	GLU	CA-CB-CG	7.71	129.53	114.10
8	EM	196	MET	CA-CB-CG	7.71	129.52	114.10
1	FC	60	LYS	CB-CG-CD	7.71	129.03	111.30
8	JM	196	MET	CA-CB-CG	7.71	129.52	114.10
1	CC	60	LYS	CB-CG-CD	7.70	129.01	111.30
3	XF	267	GLU	CA-CB-CG	7.70	129.50	114.10
3	JF	267	GLU	CA-CB-CG	7.70	129.49	114.10
1	KC	60	LYS	CB-CG-CD	7.70	129.00	111.30
8	KM	196	MET	CA-CB-CG	7.70	129.49	114.10
3	EF	267	GLU	CA-CB-CG	7.69	129.48	114.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	JC	60	LYS	CB-CG-CD	7.69	128.99	111.30
1	DC	60	LYS	CB-CG-CD	7.69	128.98	111.30
3	GF	267	GLU	CA-CB-CG	7.69	129.47	114.10
3	ZF	267	GLU	CA-CB-CG	7.69	129.47	114.10
1	AC	60	LYS	CB-CG-CD	7.68	128.97	111.30
1	BC	60	LYS	CB-CG-CD	7.68	128.98	111.30
3	VF	267	GLU	CA-CB-CG	7.68	129.47	114.10
1	HC	60	LYS	CB-CG-CD	7.68	128.97	111.30
1	IC	60	LYS	CB-CG-CD	7.68	128.96	111.30
1	LC	60	LYS	CB-CG-CD	7.68	128.96	111.30
3	YF	267	GLU	CA-CB-CG	7.68	129.46	114.10
3	AF	267	GLU	CA-CB-CG	7.68	129.45	114.10
3	DF	267	GLU	CA-CB-CG	7.67	129.45	114.10
3	LF	267	GLU	CA-CB-CG	7.67	129.45	114.10
3	BF	267	GLU	CA-CB-CG	7.67	129.44	114.10
3	WF	267	GLU	CA-CB-CG	7.67	129.44	114.10
3	CF	267	GLU	CA-CB-CG	7.67	129.43	114.10
3	FF	267	GLU	CA-CB-CG	7.67	129.43	114.10
1	GC	60	LYS	CB-CG-CD	7.66	128.93	111.30
3	IF	267	GLU	CA-CB-CG	7.66	129.42	114.10
3	KF	267	GLU	CA-CB-CG	7.66	129.41	114.10
1	MC	60	LYS	CB-CG-CD	7.66	128.91	111.30
3	MF	267	GLU	CA-CB-CG	7.66	129.41	114.10
4	WG	796	GLN	N-CA-CB	7.46	121.20	110.16
4	KG	796	GLN	N-CA-CB	7.43	121.16	110.16
4	BG	796	GLN	N-CA-CB	7.43	121.15	110.16
4	AG	796	GLN	N-CA-CB	7.42	121.14	110.16
4	IG	796	GLN	N-CA-CB	7.42	121.14	110.16
4	XG	796	GLN	N-CA-CB	7.41	121.13	110.16
4	GG	796	GLN	N-CA-CB	7.41	121.13	110.16
4	JG	796	GLN	N-CA-CB	7.41	121.12	110.16
4	DG	796	GLN	N-CA-CB	7.41	121.12	110.16
4	CG	796	GLN	N-CA-CB	7.40	121.12	110.16
4	EG	796	GLN	N-CA-CB	7.40	121.11	110.16
4	MG	796	GLN	N-CA-CB	7.40	121.11	110.16
4	ZG	796	GLN	N-CA-CB	7.40	121.11	110.16
4	HG	796	GLN	N-CA-CB	7.40	121.11	110.16
4	LG	796	GLN	N-CA-CB	7.40	121.11	110.16
4	FG	796	GLN	N-CA-CB	7.39	121.09	110.16
4	YG	796	GLN	N-CA-CB	7.38	121.09	110.16
4	VG	796	GLN	N-CA-CB	7.38	121.08	110.16
5	ZH	208	LYS	N-CA-CB	-7.37	99.11	110.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
5	IH	208	LYS	CA-CB-CG	7.37	128.83	114.10
2	Gd	82	ARG	CG-CD-NE	-7.31	95.91	112.00
8	AM	76	ARG	CB-CG-CD	-7.31	94.49	111.30
8	BM	76	ARG	CB-CG-CD	-7.31	94.49	111.30
8	CM	76	ARG	CB-CG-CD	-7.31	94.49	111.30
8	EM	76	ARG	CB-CG-CD	-7.31	94.49	111.30
8	MM	76	ARG	CB-CG-CD	-7.30	94.51	111.30
8	KM	76	ARG	CB-CG-CD	-7.30	94.51	111.30
8	EM	262	GLU	CA-CB-CG	7.29	128.69	114.10
8	IM	262	GLU	CA-CB-CG	7.29	128.69	114.10
2	Cd	125	ASP	CA-C-N	-7.29	108.94	121.85
2	Cd	125	ASP	C-N-CA	-7.29	108.94	121.85
8	DM	76	ARG	CB-CG-CD	-7.29	94.54	111.30
8	JM	76	ARG	CB-CG-CD	-7.29	94.54	111.30
8	GM	262	GLU	CA-CB-CG	7.29	128.67	114.10
8	CM	262	GLU	CA-CB-CG	7.28	128.67	114.10
8	GM	76	ARG	CB-CG-CD	-7.28	94.55	111.30
8	JM	262	GLU	CA-CB-CG	7.28	128.66	114.10
8	KM	262	GLU	CA-CB-CG	7.28	128.67	114.10
8	LM	76	ARG	CB-CG-CD	-7.28	94.55	111.30
8	IM	76	ARG	CB-CG-CD	-7.28	94.56	111.30
8	FM	76	ARG	CB-CG-CD	-7.28	94.56	111.30
8	DM	262	GLU	CA-CB-CG	7.28	128.65	114.10
8	FM	262	GLU	CA-CB-CG	7.28	128.65	114.10
8	HM	262	GLU	CA-CB-CG	7.28	128.65	114.10
8	LM	262	GLU	CA-CB-CG	7.27	128.64	114.10
8	AM	262	GLU	CA-CB-CG	7.27	128.63	114.10
8	MM	262	GLU	CA-CB-CG	7.26	128.63	114.10
8	HM	76	ARG	CB-CG-CD	-7.26	94.59	111.30
8	BM	262	GLU	CA-CB-CG	7.25	128.59	114.10
2	BD	55	MET	CG-SD-CE	-7.25	84.96	100.90
5	HH	138	GLU	N-CA-CB	7.16	120.75	110.16
2	LD	52	MET	CB-CG-SD	-7.11	91.36	112.70
2	Fd	126	GLU	CG-CD-OE1	7.09	134.71	118.40
5	CH	238	MET	CB-CG-SD	-7.07	91.48	112.70
5	EH	138	GLU	N-CA-CB	7.05	120.60	110.16
7	IL	211	LYS	CA-CB-CG	7.05	128.20	114.10
7	KL	211	LYS	CA-CB-CG	7.05	128.20	114.10
7	LL	211	LYS	CA-CB-CG	7.04	128.18	114.10
7	JL	211	LYS	CA-CB-CG	7.04	128.17	114.10
7	GL	211	LYS	CA-CB-CG	7.04	128.17	114.10
7	EL	211	LYS	CA-CB-CG	7.03	128.17	114.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
7	ML	211	LYS	CA-CB-CG	7.03	128.17	114.10
7	FL	211	LYS	CA-CB-CG	7.02	128.15	114.10
7	HL	211	LYS	CA-CB-CG	7.02	128.15	114.10
7	BL	211	LYS	CA-CB-CG	7.02	128.15	114.10
7	AL	211	LYS	CA-CB-CG	7.01	128.13	114.10
7	DL	211	LYS	CA-CB-CG	7.01	128.13	114.10
5	MH	311	MET	CB-CG-SD	-7.01	91.67	112.70
5	GH	311	MET	CB-CG-SD	-7.00	91.69	112.70
7	CL	211	LYS	CA-CB-CG	6.99	128.09	114.10
5	JH	208	LYS	CD-CE-NZ	-6.97	89.59	111.90
4	IG	795	GLN	CA-CB-CG	6.96	128.02	114.10
4	DG	795	GLN	CA-CB-CG	6.95	127.99	114.10
2	Gd	82	ARG	CA-CB-CG	6.95	127.99	114.10
4	AG	795	GLN	CA-CB-CG	6.95	127.99	114.10
4	WG	795	GLN	CA-CB-CG	6.94	127.98	114.10
4	ZG	795	GLN	CA-CB-CG	6.94	127.98	114.10
4	CG	795	GLN	CA-CB-CG	6.94	127.98	114.10
4	FG	795	GLN	CA-CB-CG	6.94	127.97	114.10
4	MG	795	GLN	CA-CB-CG	6.94	127.97	114.10
5	KH	130	LEU	CA-CB-CG	6.94	140.58	116.30
4	KG	795	GLN	CA-CB-CG	6.93	127.97	114.10
5	IH	115	MET	CB-CA-C	-6.93	99.06	110.85
4	JG	795	GLN	CA-CB-CG	6.93	127.96	114.10
4	VG	795	GLN	CA-CB-CG	6.93	127.96	114.10
4	YG	795	GLN	CA-CB-CG	6.93	127.96	114.10
4	BG	795	GLN	CA-CB-CG	6.93	127.96	114.10
4	HG	795	GLN	CA-CB-CG	6.92	127.95	114.10
4	XG	795	GLN	CA-CB-CG	6.92	127.95	114.10
2	HD	52	MET	CB-CG-SD	-6.92	91.93	112.70
4	GG	795	GLN	CA-CB-CG	6.92	127.94	114.10
4	EG	795	GLN	CA-CB-CG	6.91	127.92	114.10
4	LG	795	GLN	CA-CB-CG	6.91	127.91	114.10
5	WH	238	MET	CG-SD-CE	-6.90	85.72	100.90
2	Ld	52	MET	CA-CB-CG	6.84	127.79	114.10
6	HK	129	GLU	CA-CB-CG	6.82	127.74	114.10
6	IK	129	GLU	CA-CB-CG	6.79	127.69	114.10
6	MK	129	GLU	CA-CB-CG	6.79	127.69	114.10
5	HH	135	GLU	N-CA-CB	6.79	120.81	110.28
6	CK	129	GLU	CA-CB-CG	6.79	127.68	114.10
6	FK	129	GLU	CA-CB-CG	6.79	127.68	114.10
6	JK	129	GLU	CA-CB-CG	6.78	127.67	114.10
6	LK	129	GLU	CA-CB-CG	6.78	127.67	114.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
6	EK	129	GLU	CA-CB-CG	6.78	127.65	114.10
6	AK	129	GLU	CA-CB-CG	6.77	127.64	114.10
6	GK	129	GLU	CA-CB-CG	6.77	127.64	114.10
6	DK	129	GLU	CA-CB-CG	6.77	127.64	114.10
5	IH	115	MET	N-CA-CB	6.77	120.17	110.16
2	AD	59	GLU	CB-CA-C	-6.76	99.36	110.85
5	LH	216	GLN	N-CA-CB	6.76	122.59	110.83
6	KK	129	GLU	CA-CB-CG	6.76	127.61	114.10
5	GH	115	MET	CG-SD-CE	-6.75	86.05	100.90
6	BK	129	GLU	CA-CB-CG	6.74	127.58	114.10
5	VH	328	MET	CA-CB-CG	6.73	127.55	114.10
5	FH	115	MET	CB-CA-C	-6.71	99.44	110.85
5	ZH	115	MET	N-CA-CB	-6.69	100.28	110.12
5	ZH	150	LYS	CG-CD-CE	-6.69	95.92	111.30
2	DD	52	MET	CG-SD-CE	-6.68	86.20	100.90
5	JH	138	GLU	N-CA-CB	6.67	120.04	110.16
5	EH	138	GLU	CA-CB-CG	6.65	127.39	114.10
2	GD	107	ARG	CG-CD-NE	6.62	126.57	112.00
5	DH	135	GLU	N-CA-CB	6.61	120.53	110.28
5	WH	287	VAL	N-CA-C	-6.61	102.94	108.63
5	FH	285	GLU	CA-CB-CG	-6.60	100.91	114.10
2	Fd	126	GLU	CB-CG-CD	6.51	123.67	112.60
2	ID	52	MET	CG-SD-CE	-6.51	86.58	100.90
5	KH	115	MET	CG-SD-CE	-6.51	86.58	100.90
8	DM	145	ARG	CA-CB-CG	6.48	127.06	114.10
8	KM	145	ARG	CA-CB-CG	6.47	127.04	114.10
8	BM	145	ARG	CA-CB-CG	6.47	127.04	114.10
8	LM	145	ARG	CA-CB-CG	6.47	127.04	114.10
8	EM	145	ARG	CA-CB-CG	6.47	127.03	114.10
8	AM	145	ARG	CA-CB-CG	6.46	127.02	114.10
8	CM	145	ARG	CA-CB-CG	6.45	127.01	114.10
8	GM	145	ARG	CA-CB-CG	6.45	127.00	114.10
8	FM	145	ARG	CA-CB-CG	6.45	127.00	114.10
5	IH	214	MET	CA-CB-CG	6.45	126.99	114.10
8	HM	145	ARG	CA-CB-CG	6.44	126.99	114.10
8	MM	145	ARG	CA-CB-CG	6.44	126.99	114.10
8	IM	145	ARG	CA-CB-CG	6.44	126.98	114.10
8	JM	145	ARG	CA-CB-CG	6.43	126.97	114.10
5	AH	205	GLN	N-CA-CB	-6.40	100.39	110.69
5	LH	189	TRP	CA-CB-CG	6.39	125.74	113.60
5	ZH	339	MET	CG-SD-CE	-6.38	86.85	100.90
2	KD	119	ILE	CA-CB-CG2	6.36	121.31	110.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
8	CM	70	MET	CA-CB-CG	6.34	126.78	114.10
5	AH	141	LYS	CG-CD-CE	-6.34	96.72	111.30
8	HM	70	MET	CA-CB-CG	6.34	126.77	114.10
8	HM	196	MET	CB-CG-SD	6.33	131.69	112.70
8	IM	196	MET	CB-CG-SD	6.33	131.68	112.70
8	JM	196	MET	CB-CG-SD	6.33	131.68	112.70
8	KM	196	MET	CB-CG-SD	6.33	131.69	112.70
8	DM	70	MET	CA-CB-CG	6.33	126.75	114.10
8	FM	70	MET	CA-CB-CG	6.32	126.74	114.10
8	KM	70	MET	CA-CB-CG	6.32	126.74	114.10
8	MM	70	MET	CA-CB-CG	6.32	126.74	114.10
8	EM	196	MET	CB-CG-SD	6.31	131.64	112.70
8	GM	196	MET	CB-CG-SD	6.31	131.64	112.70
8	CM	196	MET	CB-CG-SD	6.31	131.64	112.70
8	AM	196	MET	CB-CG-SD	6.31	131.63	112.70
8	MM	196	MET	CB-CG-SD	6.31	131.63	112.70
8	LM	196	MET	CB-CG-SD	6.31	131.62	112.70
8	DM	196	MET	CB-CG-SD	6.30	131.61	112.70
5	GH	115	MET	N-CA-C	6.30	118.23	111.36
5	BH	354	MET	CG-SD-CE	-6.30	87.05	100.90
2	ED	52	MET	CG-SD-CE	-6.30	87.05	100.90
8	FM	196	MET	CB-CG-SD	6.30	131.59	112.70
8	GM	70	MET	CA-CB-CG	6.30	126.69	114.10
8	BM	196	MET	CB-CG-SD	6.29	131.58	112.70
8	BM	70	MET	CA-CB-CG	6.29	126.67	114.10
7	CL	160	GLN	CA-CB-CG	6.29	126.67	114.10
7	LL	160	GLN	CA-CB-CG	6.29	126.67	114.10
8	JM	70	MET	CA-CB-CG	6.28	126.67	114.10
8	EM	70	MET	CA-CB-CG	6.28	126.66	114.10
8	IM	70	MET	CA-CB-CG	6.28	126.65	114.10
8	LM	70	MET	CA-CB-CG	6.28	126.65	114.10
8	AM	70	MET	CA-CB-CG	6.27	126.65	114.10
7	DL	160	GLN	CA-CB-CG	6.27	126.63	114.10
7	GL	160	GLN	CA-CB-CG	6.26	126.61	114.10
7	HL	160	GLN	CA-CB-CG	6.26	126.61	114.10
7	IL	160	GLN	CA-CB-CG	6.25	126.61	114.10
7	AL	160	GLN	CA-CB-CG	6.25	126.61	114.10
7	KL	160	GLN	CA-CB-CG	6.25	126.61	114.10
5	LH	266	MET	CB-CA-C	-6.25	100.11	112.05
5	BH	115	MET	CA-CB-CG	-6.25	101.60	114.10
7	BL	160	GLN	CA-CB-CG	6.25	126.60	114.10
5	GH	248	VAL	CB-CA-C	6.25	119.46	110.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
7	ML	160	GLN	CA-CB-CG	6.25	126.60	114.10
7	EL	160	GLN	CA-CB-CG	6.25	126.59	114.10
7	FL	160	GLN	CA-CB-CG	6.24	126.59	114.10
2	Id	55	MET	CG-SD-CE	-6.24	87.17	100.90
7	JL	160	GLN	CA-CB-CG	6.23	126.56	114.10
9	JN	84	CYS	CA-CB-SG	6.19	128.63	114.40
9	EN	84	CYS	CA-CB-SG	6.19	128.63	114.40
9	KN	84	CYS	CA-CB-SG	6.19	128.63	114.40
9	AN	84	CYS	CA-CB-SG	6.18	128.61	114.40
9	CN	84	CYS	CA-CB-SG	6.18	128.61	114.40
9	FN	84	CYS	CA-CB-SG	6.18	128.61	114.40
9	LN	84	CYS	CA-CB-SG	6.18	128.61	114.40
9	DN	84	CYS	CA-CB-SG	6.17	128.60	114.40
9	HN	84	CYS	CA-CB-SG	6.17	128.60	114.40
5	EH	285	GLU	OE1-CD-OE2	6.17	137.72	122.90
9	GN	84	CYS	CA-CB-SG	6.17	128.59	114.40
5	XH	287	VAL	N-CA-C	-6.17	103.32	108.63
5	AH	208	LYS	N-CA-CB	-6.17	100.96	110.46
9	BN	84	CYS	CA-CB-SG	6.17	128.59	114.40
9	MN	84	CYS	CA-CB-SG	6.17	128.58	114.40
9	IN	84	CYS	CA-CB-SG	6.16	128.57	114.40
2	Cd	43	GLU	N-CA-CB	6.16	119.28	110.16
9	IN	115	GLN	CA-CB-CG	6.13	126.36	114.10
9	AN	115	GLN	CA-CB-CG	6.12	126.34	114.10
2	Bd	26	LYS	N-CA-C	-6.12	98.93	108.90
9	KN	115	GLN	CA-CB-CG	6.11	126.33	114.10
9	MN	115	GLN	CA-CB-CG	6.11	126.33	114.10
5	EH	135	GLU	N-CA-CB	6.11	119.75	110.28
9	BN	115	GLN	CA-CB-CG	6.11	126.32	114.10
9	DN	115	GLN	CA-CB-CG	6.11	126.32	114.10
9	FN	115	GLN	CA-CB-CG	6.11	126.32	114.10
5	EH	208	LYS	N-CA-CB	-6.11	101.05	110.46
9	JN	115	GLN	CA-CB-CG	6.11	126.31	114.10
5	BH	214	MET	CA-CB-CG	6.10	126.31	114.10
9	CN	115	GLN	CA-CB-CG	6.10	126.29	114.10
9	EN	115	GLN	CA-CB-CG	6.10	126.30	114.10
9	LN	115	GLN	CA-CB-CG	6.10	126.29	114.10
5	FH	115	MET	N-CA-CB	6.09	119.17	110.16
3	ZF	243	MET	CA-CB-CG	6.09	126.27	114.10
9	HN	115	GLN	CA-CB-CG	6.08	126.27	114.10
3	FF	243	MET	CA-CB-CG	6.08	126.27	114.10
3	VF	243	MET	CA-CB-CG	6.08	126.25	114.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	YF	243	MET	CA-CB-CG	6.08	126.25	114.10
4	ZG	813	LYS	CB-CA-C	-6.08	99.00	110.67
3	JF	243	MET	CA-CB-CG	6.07	126.25	114.10
3	CF	243	MET	CA-CB-CG	6.07	126.24	114.10
3	DF	243	MET	CA-CB-CG	6.07	126.24	114.10
3	EF	243	MET	CA-CB-CG	6.07	126.24	114.10
4	CG	813	LYS	CB-CA-C	-6.07	99.02	110.67
3	AF	243	MET	CA-CB-CG	6.07	126.23	114.10
4	AG	813	LYS	CB-CA-C	-6.07	99.03	110.67
4	VG	813	LYS	CB-CA-C	-6.07	99.03	110.67
4	DG	813	LYS	CB-CA-C	-6.06	99.03	110.67
3	GF	243	MET	CA-CB-CG	6.06	126.22	114.10
3	IF	243	MET	CA-CB-CG	6.06	126.22	114.10
3	HF	243	MET	CA-CB-CG	6.06	126.22	114.10
4	LG	813	LYS	CB-CA-C	-6.06	99.03	110.67
3	MF	243	MET	CA-CB-CG	6.06	126.22	114.10
4	EG	813	LYS	CB-CA-C	-6.06	99.04	110.67
4	JG	813	LYS	CB-CA-C	-6.06	99.04	110.67
4	WG	813	LYS	CB-CA-C	-6.06	99.04	110.67
4	HG	813	LYS	CB-CA-C	-6.06	99.04	110.67
3	KF	243	MET	CA-CB-CG	6.05	126.21	114.10
3	XF	243	MET	CA-CB-CG	6.05	126.20	114.10
3	BF	243	MET	CA-CB-CG	6.05	126.20	114.10
4	KG	813	LYS	CB-CA-C	-6.05	99.05	110.67
4	YG	813	LYS	CB-CA-C	-6.05	99.05	110.67
9	GN	115	GLN	CA-CB-CG	6.05	126.20	114.10
2	ED	52	MET	CB-CG-SD	-6.05	94.56	112.70
4	MG	813	LYS	CB-CA-C	-6.05	99.06	110.67
2	Kd	52	MET	CA-CB-CG	6.04	126.19	114.10
4	GG	813	LYS	CB-CA-C	-6.04	99.07	110.67
4	IG	813	LYS	CB-CA-C	-6.04	99.07	110.67
3	LF	243	MET	CA-CB-CG	6.04	126.18	114.10
4	FG	813	LYS	CB-CA-C	-6.03	99.09	110.67
3	WF	243	MET	CA-CB-CG	6.03	126.16	114.10
4	XG	813	LYS	CB-CA-C	-6.03	99.10	110.67
4	BG	813	LYS	CB-CA-C	-6.02	99.10	110.67
5	ZH	287	VAL	N-CA-C	-6.01	103.46	108.63
5	HH	132	GLN	N-CA-CB	5.98	119.00	110.16
2	ED	52	MET	CA-CB-CG	-5.97	102.16	114.10
2	Fd	82	ARG	CA-C-N	5.96	132.13	122.53
2	Fd	82	ARG	C-N-CA	5.96	132.13	122.53
2	CD	52	MET	CA-CB-CG	-5.96	102.17	114.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
5	MH	179	LEU	CB-CA-C	5.95	120.02	109.72
2	CD	52	MET	CB-CG-SD	-5.93	94.90	112.70
5	HH	150	LYS	CA-CB-CG	5.91	125.93	114.10
5	EH	238	MET	CG-SD-CE	-5.91	87.89	100.90
5	YH	338	GLU	N-CA-CB	5.91	119.85	110.57
5	MH	207	ASP	OD1-CG-OD2	-5.90	108.74	122.90
2	Bd	52	MET	CA-CB-CG	5.90	125.89	114.10
2	Cd	126	GLU	CG-CD-OE2	5.88	131.93	118.40
2	Fd	127	SER	N-CA-CB	5.87	118.63	110.17
2	HD	55	MET	CG-SD-CE	-5.87	87.98	100.90
5	FH	205	GLN	N-CA-CB	5.85	120.11	110.69
5	XH	328	MET	CG-SD-CE	-5.84	88.05	100.90
2	Gd	52	MET	CA-CB-CG	5.80	125.70	114.10
5	FH	251	ARG	N-CA-CB	5.80	120.33	110.71
2	DD	55	MET	CG-SD-CE	-5.79	88.16	100.90
2	DD	52	MET	CA-CB-CG	-5.78	102.54	114.10
2	LD	52	MET	CA-CB-CG	-5.73	102.64	114.10
2	Fd	74	PRO	CA-C-N	5.71	132.45	121.54
2	Fd	74	PRO	C-N-CA	5.71	132.45	121.54
5	AH	110	LYS	CB-CA-C	5.71	120.55	110.85
2	Kd	126	GLU	N-CA-CB	-5.69	101.79	111.69
2	Id	52	MET	CA-CB-CG	5.68	125.45	114.10
5	JH	354	MET	CG-SD-CE	-5.68	88.41	100.90
2	Dd	55	MET	CG-SD-CE	-5.67	88.42	100.90
5	XH	354	MET	CG-SD-CE	-5.67	88.42	100.90
5	JH	285	GLU	CA-CB-CG	-5.67	102.77	114.10
2	Dd	52	MET	CG-SD-CE	5.66	113.36	100.90
5	KH	263	ALA	CA-C-N	5.65	131.35	121.52
5	KH	263	ALA	C-N-CA	5.65	131.35	121.52
5	BH	130	LEU	CA-CB-CG	5.65	136.06	116.30
2	Fd	52	MET	CG-SD-CE	-5.64	88.48	100.90
5	ZH	214	MET	CA-CB-CG	5.64	125.38	114.10
5	IH	196	LEU	N-CA-CB	5.63	120.15	110.68
5	CH	122	LEU	CD1-CG-CD2	-5.63	98.41	110.80
5	LH	266	MET	N-CA-CB	5.62	119.15	110.43
1	IC	69	GLN	CB-CG-CD	5.60	122.13	112.60
1	MC	69	GLN	CB-CG-CD	5.60	122.12	112.60
1	DC	69	GLN	CB-CG-CD	5.60	122.11	112.60
5	GH	135	GLU	N-CA-CB	5.59	118.95	110.28
1	FC	69	GLN	CB-CG-CD	5.59	122.11	112.60
1	AC	69	GLN	CB-CG-CD	5.59	122.10	112.60
2	Jd	52	MET	CG-SD-CE	-5.59	88.61	100.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	HC	69	GLN	CB-CG-CD	5.59	122.10	112.60
1	LC	69	GLN	CB-CG-CD	5.59	122.10	112.60
1	CC	210	ILE	CB-CG1-CD1	-5.58	102.08	113.80
1	IC	210	ILE	CB-CG1-CD1	-5.58	102.07	113.80
1	CC	69	GLN	CB-CG-CD	5.58	122.09	112.60
1	JC	69	GLN	CB-CG-CD	5.58	122.09	112.60
2	Jd	125	ASP	CA-C-N	-5.58	110.92	121.63
2	Jd	125	ASP	C-N-CA	-5.58	110.92	121.63
1	KC	69	GLN	CB-CG-CD	5.58	122.08	112.60
1	KC	210	ILE	CB-CG1-CD1	-5.57	102.10	113.80
5	FH	285	GLU	N-CA-CB	-5.57	101.14	110.39
1	BC	69	GLN	CB-CG-CD	5.57	122.06	112.60
1	HC	210	ILE	CB-CG1-CD1	-5.57	102.11	113.80
1	BC	210	ILE	CB-CG1-CD1	-5.57	102.11	113.80
1	FC	210	ILE	CB-CG1-CD1	-5.57	102.11	113.80
1	GC	69	GLN	CB-CG-CD	5.57	122.06	112.60
5	JH	311	MET	CB-CG-SD	-5.57	96.00	112.70
5	AH	135	GLU	N-CA-CB	5.56	118.90	110.28
1	EC	69	GLN	CB-CG-CD	5.56	122.05	112.60
5	WH	115	MET	CA-CB-CG	-5.55	102.99	114.10
1	DC	210	ILE	CB-CG1-CD1	-5.55	102.14	113.80
1	AC	210	ILE	CB-CG1-CD1	-5.55	102.15	113.80
2	LD	52	MET	CG-SD-CE	-5.55	88.69	100.90
1	LC	210	ILE	CB-CG1-CD1	-5.55	102.15	113.80
5	AH	115	MET	N-CA-CB	-5.54	101.96	110.16
1	EC	210	ILE	CB-CG1-CD1	-5.54	102.17	113.80
1	JC	210	ILE	CB-CG1-CD1	-5.54	102.17	113.80
1	MC	210	ILE	CB-CG1-CD1	-5.54	102.17	113.80
5	JH	107	ILE	CA-CB-CG1	5.54	119.81	110.40
1	GC	210	ILE	CB-CG1-CD1	-5.53	102.18	113.80
2	LD	55	MET	CA-CB-CG	5.53	125.16	114.10
5	ZH	208	LYS	CA-CB-CG	5.53	125.16	114.10
5	BH	115	MET	CB-CG-SD	-5.52	96.14	112.70
6	IK	142	GLU	CA-CB-CG	-5.52	103.06	114.10
6	CK	142	GLU	CA-CB-CG	-5.52	103.07	114.10
5	HH	208	LYS	N-CA-CB	-5.51	101.97	110.46
5	KH	250	TYR	N-CA-CB	-5.51	101.99	110.20
5	AH	205	GLN	N-CA-C	5.51	118.46	109.59
6	AK	142	GLU	CA-CB-CG	-5.51	103.08	114.10
5	CH	122	LEU	CB-CG-CD2	-5.51	94.18	110.70
6	MK	142	GLU	CA-CB-CG	-5.51	103.09	114.10
5	VH	287	VAL	CB-CA-C	5.51	116.06	110.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
6	KK	122	LYS	CG-CD-CE	5.50	123.95	111.30
6	KK	142	GLU	CA-CB-CG	-5.50	103.10	114.10
1	DC	61	GLU	CA-CB-CG	5.50	125.10	114.10
6	JK	122	LYS	CG-CD-CE	5.50	123.95	111.30
5	EH	130	LEU	CA-CB-CG	5.50	135.55	116.30
6	HK	142	GLU	CA-CB-CG	-5.50	103.11	114.10
4	GG	822	GLU	CB-CG-CD	5.50	121.94	112.60
1	GC	61	GLU	CA-CB-CG	5.49	125.09	114.10
6	LK	129	GLU	CB-CG-CD	5.49	121.94	112.60
6	GK	142	GLU	CA-CB-CG	-5.49	103.12	114.10
6	LK	142	GLU	CA-CB-CG	-5.49	103.12	114.10
1	BC	61	GLU	CA-CB-CG	5.49	125.08	114.10
1	CC	61	GLU	CA-CB-CG	5.49	125.08	114.10
1	EC	61	GLU	CA-CB-CG	5.49	125.08	114.10
4	KG	797	ARG	CB-CG-CD	5.49	123.92	111.30
4	CG	822	GLU	CB-CG-CD	5.49	121.93	112.60
6	DK	122	LYS	CG-CD-CE	5.49	123.92	111.30
6	HK	122	LYS	CG-CD-CE	5.49	123.92	111.30
6	CK	122	LYS	CG-CD-CE	5.48	123.91	111.30
4	CG	797	ARG	CB-CG-CD	5.48	123.91	111.30
5	IH	214	MET	CB-CG-SD	5.48	129.15	112.70
6	EK	142	GLU	CA-CB-CG	-5.48	103.14	114.10
1	FC	61	GLU	CA-CB-CG	5.48	125.06	114.10
6	GK	122	LYS	CG-CD-CE	5.48	123.91	111.30
1	JC	61	GLU	CA-CB-CG	5.48	125.06	114.10
6	BK	142	GLU	CA-CB-CG	-5.48	103.15	114.10
6	FK	122	LYS	CG-CD-CE	5.48	123.90	111.30
6	JK	142	GLU	CA-CB-CG	-5.48	103.14	114.10
4	BG	797	ARG	CB-CG-CD	5.48	123.89	111.30
6	FK	142	GLU	CA-CB-CG	-5.47	103.15	114.10
4	AG	797	ARG	CB-CG-CD	5.47	123.89	111.30
6	AK	122	LYS	CG-CD-CE	5.47	123.89	111.30
4	DG	822	GLU	CB-CG-CD	5.47	121.90	112.60
4	EG	797	ARG	CB-CG-CD	5.47	123.89	111.30
4	EG	822	GLU	CB-CG-CD	5.47	121.91	112.60
5	GH	151	PRO	N-CA-C	5.47	119.19	110.21
4	HG	797	ARG	CB-CG-CD	5.47	123.88	111.30
1	IC	61	GLU	CA-CB-CG	5.47	125.04	114.10
6	DK	142	GLU	CA-CB-CG	-5.47	103.16	114.10
6	EK	122	LYS	CG-CD-CE	5.47	123.88	111.30
4	FG	797	ARG	CB-CG-CD	5.47	123.88	111.30
4	JG	797	ARG	CB-CG-CD	5.47	123.88	111.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	GG	797	ARG	CB-CG-CD	5.47	123.87	111.30
4	XG	797	ARG	CB-CG-CD	5.47	123.87	111.30
6	BK	129	GLU	CB-CG-CD	5.46	121.89	112.60
1	MC	61	GLU	CA-CB-CG	5.46	125.03	114.10
4	YG	797	ARG	CB-CG-CD	5.46	123.87	111.30
6	LK	122	LYS	CG-CD-CE	5.46	123.86	111.30
4	AG	822	GLU	CB-CG-CD	5.46	121.88	112.60
1	KC	61	GLU	CA-CB-CG	5.46	125.03	114.10
4	XG	822	GLU	CB-CG-CD	5.46	121.89	112.60
4	WG	822	GLU	CB-CG-CD	5.46	121.88	112.60
6	BK	122	LYS	CG-CD-CE	5.46	123.85	111.30
1	LC	61	GLU	CA-CB-CG	5.46	125.01	114.10
4	YG	822	GLU	CB-CG-CD	5.46	121.88	112.60
6	AK	129	GLU	CB-CG-CD	5.46	121.87	112.60
4	FG	822	GLU	CB-CG-CD	5.46	121.87	112.60
4	IG	797	ARG	CB-CG-CD	5.46	123.85	111.30
4	VG	797	ARG	CB-CG-CD	5.46	123.85	111.30
4	JG	822	GLU	CB-CG-CD	5.45	121.87	112.60
2	ED	160	LYS	CA-CB-CG	5.45	125.00	114.10
4	MG	822	GLU	CB-CG-CD	5.45	121.87	112.60
4	MG	797	ARG	CB-CG-CD	5.45	123.84	111.30
1	HC	157	LYS	CA-CB-CG	5.45	125.00	114.10
6	MK	122	LYS	CG-CD-CE	5.45	123.83	111.30
4	WG	797	ARG	CB-CG-CD	5.45	123.83	111.30
4	ZG	797	ARG	CB-CG-CD	5.45	123.83	111.30
6	IK	122	LYS	CG-CD-CE	5.45	123.83	111.30
1	AC	157	LYS	CA-CB-CG	5.45	124.99	114.10
1	HC	61	GLU	CA-CB-CG	5.45	124.99	114.10
4	HG	822	GLU	CB-CG-CD	5.45	121.86	112.60
4	DG	797	ARG	CB-CG-CD	5.44	123.82	111.30
6	DK	129	GLU	CB-CG-CD	5.44	121.85	112.60
4	ZG	822	GLU	CB-CG-CD	5.44	121.85	112.60
6	KK	129	GLU	CB-CG-CD	5.44	121.85	112.60
4	LG	822	GLU	CB-CG-CD	5.44	121.85	112.60
1	AC	61	GLU	CA-CB-CG	5.44	124.98	114.10
6	MK	129	GLU	CB-CG-CD	5.44	121.85	112.60
4	VG	822	GLU	CB-CG-CD	5.44	121.85	112.60
6	GK	129	GLU	CB-CG-CD	5.44	121.84	112.60
6	JK	129	GLU	CB-CG-CD	5.44	121.84	112.60
4	KG	822	GLU	CB-CG-CD	5.44	121.84	112.60
4	BG	822	GLU	CB-CG-CD	5.43	121.83	112.60
4	LG	797	ARG	CB-CG-CD	5.43	123.79	111.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	Ld	52	MET	CB-CG-SD	-5.43	96.41	112.70
1	MC	157	LYS	CA-CB-CG	5.42	124.95	114.10
1	KC	157	LYS	CA-CB-CG	5.42	124.94	114.10
1	IC	157	LYS	CA-CB-CG	5.42	124.94	114.10
6	IK	129	GLU	CB-CG-CD	5.42	121.81	112.60
1	LC	157	LYS	CA-CB-CG	5.42	124.94	114.10
7	JL	211	LYS	CG-CD-CE	5.42	123.77	111.30
1	BC	157	LYS	CA-CB-CG	5.42	124.94	114.10
6	CK	129	GLU	CB-CG-CD	5.42	121.81	112.60
4	IG	822	GLU	CB-CG-CD	5.42	121.81	112.60
1	DC	157	LYS	CA-CB-CG	5.42	124.93	114.10
6	EK	129	GLU	CB-CG-CD	5.41	121.80	112.60
1	GC	157	LYS	CA-CB-CG	5.41	124.91	114.10
6	HK	129	GLU	CB-CG-CD	5.41	121.79	112.60
5	LH	238	MET	N-CA-C	5.41	117.77	109.07
1	JC	157	LYS	CA-CB-CG	5.40	124.91	114.10
1	CC	157	LYS	CA-CB-CG	5.40	124.90	114.10
1	EC	157	LYS	CA-CB-CG	5.40	124.90	114.10
6	FK	129	GLU	CB-CG-CD	5.40	121.78	112.60
5	BH	214	MET	CB-CG-SD	5.40	128.90	112.70
1	FC	157	LYS	CA-CB-CG	5.40	124.89	114.10
2	Ad	52	MET	CB-CG-SD	-5.40	96.51	112.70
7	GL	211	LYS	CG-CD-CE	5.40	123.71	111.30
7	CL	211	LYS	CG-CD-CE	5.39	123.70	111.30
5	MH	285	GLU	CA-CB-CG	-5.39	103.31	114.10
7	ML	211	LYS	CG-CD-CE	5.39	123.70	111.30
7	BL	211	LYS	CG-CD-CE	5.39	123.69	111.30
7	EL	211	LYS	CG-CD-CE	5.38	123.67	111.30
7	LL	211	LYS	CG-CD-CE	5.38	123.67	111.30
7	HL	211	LYS	CG-CD-CE	5.37	123.65	111.30
5	IH	321	SER	N-CA-C	5.37	121.68	109.81
7	IL	211	LYS	CG-CD-CE	5.37	123.65	111.30
7	KL	211	LYS	CG-CD-CE	5.37	123.65	111.30
7	AL	211	LYS	CG-CD-CE	5.37	123.64	111.30
7	DL	211	LYS	CG-CD-CE	5.37	123.64	111.30
7	FL	211	LYS	CG-CD-CE	5.37	123.64	111.30
2	ID	55	MET	CG-SD-CE	-5.36	89.10	100.90
5	AH	110	LYS	CA-CB-CG	5.36	124.82	114.10
2	Fd	26	LYS	N-CA-C	-5.36	100.67	109.40
8	LM	96	GLN	CA-C-N	5.36	128.49	121.85
8	LM	96	GLN	C-N-CA	5.36	128.49	121.85
2	Jd	52	MET	CB-CA-C	-5.35	101.91	110.79

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	ID	52	MET	CA-CB-CG	-5.35	103.40	114.10
2	JD	52	MET	CB-CG-SD	-5.35	96.66	112.70
5	HH	115	MET	N-CA-C	5.34	117.18	111.36
8	MM	96	GLN	CA-C-N	5.33	128.47	121.85
8	MM	96	GLN	C-N-CA	5.33	128.47	121.85
8	BM	96	GLN	CA-C-N	5.33	128.46	121.85
8	BM	96	GLN	C-N-CA	5.33	128.46	121.85
5	VH	328	MET	N-CA-CB	-5.32	102.43	111.69
8	AM	96	GLN	CA-C-N	5.31	128.44	121.85
8	AM	96	GLN	C-N-CA	5.31	128.44	121.85
5	DH	309	GLU	CA-CB-CG	5.31	124.72	114.10
8	IM	96	GLN	CA-C-N	5.31	128.44	121.85
8	IM	96	GLN	C-N-CA	5.31	128.44	121.85
2	LD	123	THR	OG1-CB-CG2	-5.31	98.68	109.30
8	GM	96	GLN	CA-C-N	5.31	128.43	121.85
8	GM	96	GLN	C-N-CA	5.31	128.43	121.85
8	KM	96	GLN	CA-C-N	5.30	128.42	121.85
8	KM	96	GLN	C-N-CA	5.30	128.42	121.85
8	CM	96	GLN	CA-C-N	5.30	128.42	121.85
8	CM	96	GLN	C-N-CA	5.30	128.42	121.85
8	JM	96	GLN	CA-C-N	5.29	128.41	121.85
8	JM	96	GLN	C-N-CA	5.29	128.41	121.85
8	FM	96	GLN	CA-C-N	5.29	128.40	121.85
8	FM	96	GLN	C-N-CA	5.29	128.40	121.85
8	DM	96	GLN	CA-C-N	5.28	128.40	121.85
8	DM	96	GLN	C-N-CA	5.28	128.40	121.85
5	FH	208	LYS	CD-CE-NZ	-5.28	95.00	111.90
5	EH	309	GLU	CA-CB-CG	5.28	124.66	114.10
2	Bd	55	MET	CG-SD-CE	-5.27	89.30	100.90
5	LH	321	SER	N-CA-C	5.27	121.47	109.81
8	HM	96	GLN	CA-C-N	5.27	128.39	121.85
8	HM	96	GLN	C-N-CA	5.27	128.39	121.85
5	MH	280	LEU	CA-CB-CG	5.27	134.75	116.30
5	HH	115	MET	CB-CA-C	-5.27	101.89	110.85
5	GH	160	LEU	CB-CG-CD2	-5.27	94.89	110.70
5	ZH	242	ILE	CB-CG1-CD1	-5.26	102.74	113.80
7	DL	45	PRO	CA-C-N	-5.26	113.27	119.84
7	DL	45	PRO	C-N-CA	-5.26	113.27	119.84
8	JM	68	GLU	CG-CD-OE1	5.25	130.48	118.40
8	BM	68	GLU	CG-CD-OE1	5.25	130.47	118.40
7	KL	45	PRO	CA-C-N	-5.25	113.28	119.84
7	KL	45	PRO	C-N-CA	-5.25	113.28	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	Ld	130	GLU	CB-CA-C	5.25	120.74	110.67
8	HM	68	GLU	CG-CD-OE1	5.24	130.46	118.40
8	CM	68	GLU	CG-CD-OE1	5.24	130.45	118.40
8	LM	68	GLU	CG-CD-OE1	5.24	130.45	118.40
8	GM	68	GLU	CG-CD-OE1	5.24	130.45	118.40
7	HL	45	PRO	CA-C-N	-5.24	113.29	119.84
7	HL	45	PRO	C-N-CA	-5.24	113.29	119.84
8	FM	68	GLU	CG-CD-OE1	5.24	130.44	118.40
5	JH	321	SER	N-CA-C	5.24	121.38	109.81
8	DM	68	GLU	CG-CD-OE1	5.23	130.43	118.40
8	KM	68	GLU	CG-CD-OE1	5.23	130.43	118.40
5	GH	216	GLN	N-CA-CB	5.23	119.93	110.83
5	WH	120	TYR	CA-CB-CG	-5.23	104.49	113.90
8	IM	68	GLU	CG-CD-OE1	5.22	130.41	118.40
5	AH	321	SER	N-CA-C	5.22	121.34	109.81
8	AM	68	GLU	CG-CD-OE1	5.22	130.40	118.40
8	EM	68	GLU	CG-CD-OE1	5.22	130.40	118.40
8	MM	68	GLU	CG-CD-OE1	5.22	130.40	118.40
5	LH	158	VAL	CG1-CB-CG2	-5.21	99.33	110.80
7	JL	45	PRO	CA-C-N	-5.21	113.32	119.84
7	JL	45	PRO	C-N-CA	-5.21	113.32	119.84
7	CL	45	PRO	CA-C-N	-5.21	113.32	119.84
7	CL	45	PRO	C-N-CA	-5.21	113.32	119.84
7	LL	45	PRO	CA-C-N	-5.21	113.33	119.84
7	LL	45	PRO	C-N-CA	-5.21	113.33	119.84
4	HG	795	GLN	CB-CG-CD	5.20	121.44	112.60
7	FL	45	PRO	CA-C-N	-5.19	113.35	119.84
7	FL	45	PRO	C-N-CA	-5.19	113.35	119.84
7	EL	45	PRO	CA-C-N	-5.19	113.36	119.84
7	EL	45	PRO	C-N-CA	-5.19	113.36	119.84
5	ZH	208	LYS	CB-CG-CD	-5.19	99.37	111.30
9	CN	111	GLU	CB-CA-C	5.18	119.66	110.85
4	GG	795	GLN	CB-CG-CD	5.18	121.41	112.60
7	GL	45	PRO	CA-C-N	-5.18	113.36	119.84
7	GL	45	PRO	C-N-CA	-5.18	113.36	119.84
5	FH	208	LYS	N-CA-CB	-5.18	102.48	110.46
4	YG	795	GLN	CB-CG-CD	5.18	121.41	112.60
7	IL	45	PRO	CA-C-N	-5.18	113.37	119.84
7	IL	45	PRO	C-N-CA	-5.18	113.37	119.84
5	IH	231	ARG	N-CA-CB	5.17	117.82	110.16
7	ML	45	PRO	CA-C-N	-5.17	113.37	119.84
7	ML	45	PRO	C-N-CA	-5.17	113.37	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	LG	795	GLN	CB-CG-CD	5.17	121.39	112.60
4	XG	795	GLN	CB-CG-CD	5.17	121.39	112.60
4	WG	795	GLN	CB-CG-CD	5.17	121.38	112.60
7	AL	45	PRO	CA-C-N	-5.16	113.39	119.84
7	AL	45	PRO	C-N-CA	-5.16	113.39	119.84
7	BL	45	PRO	CA-C-N	-5.16	113.39	119.84
7	BL	45	PRO	C-N-CA	-5.16	113.39	119.84
4	FG	795	GLN	CB-CG-CD	5.16	121.38	112.60
9	JN	111	GLU	CB-CA-C	5.16	119.62	110.85
4	KG	795	GLN	CB-CG-CD	5.16	121.37	112.60
9	KN	111	GLU	CB-CA-C	5.16	119.62	110.85
4	DG	795	GLN	CB-CG-CD	5.16	121.37	112.60
4	BG	795	GLN	CB-CG-CD	5.15	121.36	112.60
4	EG	795	GLN	CB-CG-CD	5.15	121.36	112.60
4	JG	795	GLN	CB-CG-CD	5.15	121.36	112.60
5	VH	309	GLU	CA-CB-CG	5.15	124.41	114.10
9	FN	111	GLU	CB-CA-C	5.15	119.61	110.85
9	HN	111	GLU	CB-CA-C	5.15	119.61	110.85
4	ZG	795	GLN	CB-CG-CD	5.15	121.36	112.60
9	DN	111	GLU	CB-CA-C	5.15	119.60	110.85
9	EN	111	GLU	CB-CA-C	5.14	119.60	110.85
4	IG	795	GLN	CB-CG-CD	5.14	121.34	112.60
4	VG	795	GLN	CB-CG-CD	5.14	121.34	112.60
9	AN	111	GLU	CB-CA-C	5.14	119.59	110.85
8	IM	70	MET	CB-CG-SD	5.14	128.12	112.70
9	IN	111	GLU	CB-CA-C	5.14	119.59	110.85
5	WH	328	MET	CA-CB-CG	5.14	124.37	114.10
9	BN	111	GLU	CB-CA-C	5.13	119.58	110.85
5	HH	321	SER	N-CA-C	5.13	121.16	109.81
4	AG	795	GLN	CB-CG-CD	5.13	121.32	112.60
9	GN	111	GLU	CB-CA-C	5.13	119.57	110.85
4	CG	795	GLN	CB-CG-CD	5.13	121.32	112.60
8	LM	70	MET	CB-CG-SD	5.13	128.09	112.70
9	MN	111	GLU	CB-CA-C	5.13	119.57	110.85
8	GM	70	MET	CB-CG-SD	5.12	128.07	112.70
5	ZH	317	LEU	CB-CG-CD1	-5.12	95.32	110.70
8	EM	70	MET	CB-CG-SD	5.12	128.07	112.70
8	AM	70	MET	CB-CG-SD	5.12	128.06	112.70
8	JM	70	MET	CB-CG-SD	5.12	128.06	112.70
2	KD	55	MET	CA-CB-CG	5.12	124.33	114.10
4	MG	795	GLN	CB-CG-CD	5.12	121.30	112.60
7	JL	142	GLU	CA-CB-CG	5.11	124.33	114.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
8	DM	70	MET	CB-CG-SD	5.11	128.03	112.70
8	BM	70	MET	CB-CG-SD	5.11	128.03	112.70
8	MM	70	MET	CB-CG-SD	5.11	128.03	112.70
9	LN	111	GLU	CB-CA-C	5.11	119.53	110.85
5	AH	132	GLN	N-CA-CB	5.10	117.62	110.12
7	HL	142	GLU	CA-CB-CG	5.10	124.31	114.10
8	CM	70	MET	CB-CG-SD	5.10	128.01	112.70
8	FM	70	MET	CB-CG-SD	5.10	128.00	112.70
8	KM	70	MET	CB-CG-SD	5.10	128.00	112.70
7	AL	142	GLU	CA-CB-CG	5.10	124.29	114.10
5	BH	107	ILE	CA-CB-CG1	5.09	119.06	110.40
7	LL	142	GLU	CA-CB-CG	5.09	124.29	114.10
5	MH	150	LYS	CB-CG-CD	5.09	123.02	111.30
7	IL	142	GLU	CA-CB-CG	5.09	124.29	114.10
8	HM	70	MET	CB-CG-SD	5.09	127.97	112.70
5	MH	179	LEU	CA-CB-CG	5.09	134.11	116.30
7	DL	142	GLU	CA-CB-CG	5.09	124.27	114.10
5	WH	356	LEU	N-CA-CB	5.08	119.78	111.08
2	Jd	52	MET	CA-CB-CG	5.08	124.26	114.10
5	JH	193	ALA	CA-C-N	-5.08	113.99	122.21
5	JH	193	ALA	C-N-CA	-5.08	113.99	122.21
7	BL	142	GLU	CA-CB-CG	5.07	124.25	114.10
1	JC	157	LYS	CB-CG-CD	5.07	122.97	111.30
1	CC	157	LYS	CB-CG-CD	5.07	122.97	111.30
1	EC	157	LYS	CB-CG-CD	5.07	122.96	111.30
7	KL	142	GLU	CA-CB-CG	5.07	124.24	114.10
7	EL	142	GLU	CA-CB-CG	5.07	124.24	114.10
1	KC	157	LYS	CB-CG-CD	5.07	122.95	111.30
1	BC	157	LYS	CB-CG-CD	5.06	122.95	111.30
1	GC	157	LYS	CB-CG-CD	5.06	122.95	111.30
1	DC	157	LYS	CB-CG-CD	5.06	122.94	111.30
1	FC	157	LYS	CB-CG-CD	5.06	122.94	111.30
1	LC	157	LYS	CB-CG-CD	5.06	122.94	111.30
1	MC	157	LYS	CB-CG-CD	5.06	122.94	111.30
7	ML	142	GLU	CA-CB-CG	5.06	124.22	114.10
1	AC	157	LYS	CB-CG-CD	5.06	122.94	111.30
7	CL	142	GLU	CA-CB-CG	5.06	124.22	114.10
7	GL	142	GLU	CA-CB-CG	5.06	124.21	114.10
5	BH	220	LEU	CA-CB-CG	5.05	133.99	116.30
7	FL	142	GLU	CA-CB-CG	5.05	124.20	114.10
1	IC	157	LYS	CB-CG-CD	5.05	122.91	111.30
1	HC	157	LYS	CB-CG-CD	5.04	122.88	111.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
7	AL	110	GLN	CB-CG-CD	5.03	121.16	112.60
2	Kd	130	GLU	CB-CG-CD	-5.03	104.04	112.60
8	MM	262	GLU	CB-CA-C	-5.03	101.81	110.16
5	EH	321	SER	N-CA-C	5.03	120.93	109.81
7	BL	110	GLN	CB-CG-CD	5.03	121.15	112.60
5	AH	208	LYS	CA-CB-CG	5.03	124.15	114.10
8	BM	262	GLU	CB-CA-C	-5.03	101.82	110.16
7	HL	110	GLN	CB-CG-CD	5.03	121.14	112.60
8	IM	262	GLU	CB-CA-C	-5.03	101.82	110.16
8	HM	262	GLU	CB-CA-C	-5.02	101.82	110.16
7	LL	110	GLN	CB-CG-CD	5.02	121.14	112.60
7	KL	110	GLN	CB-CG-CD	5.02	121.13	112.60
8	LM	262	GLU	CB-CA-C	-5.02	101.83	110.16
2	KD	119	ILE	CB-CG1-CD1	-5.02	103.26	113.80
8	AM	262	GLU	CB-CA-C	-5.02	101.83	110.16
7	ML	110	GLN	CB-CG-CD	5.01	121.12	112.60
7	JL	110	GLN	CB-CG-CD	5.01	121.12	112.60
2	CD	93	GLU	CB-CG-CD	5.01	121.11	112.60
8	DM	262	GLU	CB-CA-C	-5.01	101.84	110.16
8	JM	145	ARG	CB-CG-CD	5.00	122.81	111.30
5	BH	223	TYR	N-CA-C	5.00	117.75	109.85
8	FM	262	GLU	CB-CA-C	-5.00	101.86	110.16

There are no chirality outliers.

All (63) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	AD	27	LYS	Peptide
5	AH	121	PRO	Mainchain
5	AH	137	SER	Mainchain
8	AM	68	GLU	Sidechain
9	AN	115	GLN	Peptide
2	BD	27	LYS	Peptide
5	BH	121	PRO	Mainchain
8	BM	68	GLU	Sidechain
9	BN	115	GLN	Peptide
2	Bd	26	LYS	Mainchain
2	CD	27	LYS	Peptide
8	CM	68	GLU	Sidechain
9	CN	115	GLN	Peptide
2	Cd	39	ILE	Mainchain
2	DD	78	ASN	Peptide

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Mol	Chain	Res	Type	Group
5	DH	131	LYS	Mainchain
8	DM	68	GLU	Sidechain
9	DN	115	GLN	Peptide
2	ED	27	LYS	Peptide
5	EH	131	LYS	Mainchain
8	EM	68	GLU	Sidechain
9	EN	115	GLN	Peptide
2	FD	27	LYS	Peptide
5	FH	121	PRO	Mainchain
5	FH	285	GLU	Sidechain
8	FM	68	GLU	Sidechain
9	FN	115	GLN	Peptide
2	Fd	125	ASP	Mainchain
2	GD	27	LYS	Peptide
5	GH	121	PRO	Mainchain
5	GH	263	ALA	Peptide
8	GM	68	GLU	Sidechain
9	GN	115	GLN	Peptide
5	HH	121	PRO	Mainchain
8	HM	68	GLU	Sidechain
9	HN	115	GLN	Peptide
2	ID	27	LYS	Peptide
5	IH	121	PRO	Mainchain
8	IM	68	GLU	Sidechain
9	IN	115	GLN	Peptide
5	JH	121	PRO	Mainchain
5	JH	125	GLU	Sidechain
5	JH	267	PRO	Peptide
8	JM	68	GLU	Sidechain
9	JN	115	GLN	Peptide
2	KD	27	LYS	Peptide
5	KH	121	PRO	Mainchain
5	KH	263	ALA	Peptide
8	KM	68	GLU	Sidechain
9	KN	115	GLN	Peptide
5	LH	121	PRO	Mainchain
8	LM	68	GLU	Sidechain
9	LN	115	GLN	Peptide
2	MD	27	LYS	Peptide
5	MH	121	PRO	Mainchain
8	MM	68	GLU	Sidechain
9	MN	115	GLN	Peptide

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Mol	Chain	Res	Type	Group
5	VH	134	TYR	Mainchain
5	WH	121	PRO	Mainchain
5	XH	121	PRO	Mainchain
5	YH	121	PRO	Mainchain
5	YH	350	HIS	Sidechain
5	ZH	121	PRO	Mainchain

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	AC	1685	0	1684	32	0
1	BC	1685	0	1684	41	0
1	CC	1685	0	1684	46	0
1	DC	1685	0	1684	42	0
1	EC	1685	0	1684	31	0
1	FC	1685	0	1684	24	0
1	GC	1685	0	1684	36	0
1	HC	1685	0	1684	36	0
1	IC	1685	0	1684	30	0
1	JC	1685	0	1684	30	0
1	KC	1685	0	1684	32	0
1	LC	1685	0	1684	25	0
1	MC	1685	0	1684	33	0
2	AD	1086	0	1121	15	0
2	Ad	1058	0	1092	8	0
2	BD	1086	0	1121	14	0
2	Bd	1058	0	1092	16	0
2	CD	1086	0	1121	20	0
2	Cd	1058	0	1092	12	0
2	DD	1086	0	1121	17	0
2	Dd	1058	0	1092	8	0
2	ED	1086	0	1119	18	0
2	Ed	1058	0	1092	12	0
2	FD	1086	0	1121	10	0
2	Fd	1058	0	1092	22	0
2	GD	1086	0	1121	13	0
2	Gd	1058	0	1092	12	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	HD	1086	0	1121	14	0
2	Hd	1058	0	1092	8	0
2	ID	1086	0	1121	18	0
2	Id	1058	0	1092	18	0
2	JD	1086	0	1121	13	0
2	Jd	1058	0	1092	6	0
2	KD	1086	0	1121	17	0
2	Kd	1058	0	1092	10	0
2	LD	1086	0	1121	16	0
2	Ld	1058	0	1092	4	0
2	MD	1086	0	1121	18	0
2	Md	1058	0	1092	10	0
3	AF	483	0	502	8	0
3	Af	449	0	474	6	0
3	BF	483	0	502	9	0
3	Bf	449	0	474	5	0
3	CF	483	0	502	8	0
3	Cf	449	0	474	4	0
3	DF	483	0	502	7	0
3	Df	449	0	474	5	0
3	EF	483	0	502	8	0
3	Ef	449	0	474	3	0
3	FF	483	0	502	9	0
3	Ff	449	0	474	5	0
3	GF	483	0	502	15	0
3	Gf	449	0	474	4	0
3	HF	483	0	502	8	0
3	Hf	449	0	474	15	0
3	IF	483	0	502	8	0
3	If	449	0	474	5	0
3	JF	483	0	502	10	0
3	Jf	449	0	474	6	0
3	KF	483	0	502	7	0
3	Kf	449	0	474	15	0
3	LF	483	0	502	7	0
3	Lf	449	0	474	7	0
3	MF	483	0	502	7	0
3	Mf	449	0	474	4	0
3	VF	483	0	502	8	0
3	WF	483	0	502	8	0
3	XF	483	0	502	15	0
3	YF	483	0	502	7	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	ZF	483	0	502	10	0
4	AG	276	0	263	17	0
4	BG	276	0	263	8	0
4	CG	276	0	263	10	0
4	DG	276	0	263	15	0
4	EG	276	0	263	9	0
4	FG	276	0	263	9	0
4	GG	276	0	263	14	0
4	HG	276	0	263	12	0
4	IG	276	0	263	8	0
4	JG	276	0	263	16	0
4	KG	276	0	263	13	0
4	LG	276	0	263	11	0
4	MG	276	0	263	9	0
4	VG	276	0	263	10	0
4	WG	276	0	263	11	0
4	XG	276	0	263	9	0
4	YG	276	0	263	6	0
4	ZG	276	0	263	14	0
5	AH	1983	0	2009	35	0
5	BH	1983	0	2009	32	0
5	CH	1983	0	2009	44	0
5	DH	1983	0	2009	23	0
5	EH	1983	0	2009	41	0
5	FH	1983	0	2009	28	0
5	GH	1983	0	2009	43	0
5	HH	1983	0	2009	33	0
5	IH	1983	0	2009	43	0
5	JH	1983	0	2009	31	0
5	KH	1983	0	2009	24	0
5	LH	1983	0	2009	36	0
5	MH	1983	0	2009	30	0
5	VH	1875	0	1900	14	0
5	WH	1875	0	1900	26	0
5	XH	1875	0	1900	27	0
5	YH	1875	0	1900	20	0
5	ZH	1875	0	1900	31	0
6	AK	1175	0	1221	9	0
6	BK	1175	0	1221	13	0
6	CK	1175	0	1221	11	0
6	DK	1175	0	1221	11	0
6	EK	1175	0	1221	13	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	FK	1175	0	1221	9	0
6	GK	1175	0	1221	10	0
6	HK	1175	0	1221	10	0
6	IK	1175	0	1221	12	0
6	JK	1175	0	1221	12	0
6	KK	1175	0	1221	11	0
6	LK	1175	0	1221	9	0
6	MK	1175	0	1221	9	0
7	AL	1388	0	1378	15	0
7	BL	1388	0	1378	17	0
7	CL	1388	0	1378	16	0
7	DL	1388	0	1378	16	0
7	EL	1388	0	1378	16	0
7	FL	1388	0	1378	17	0
7	GL	1388	0	1378	17	0
7	HL	1388	0	1378	17	0
7	IL	1388	0	1378	19	0
7	JL	1388	0	1378	17	0
7	KL	1388	0	1378	17	0
7	LL	1388	0	1378	18	0
7	ML	1388	0	1378	19	0
8	AM	2320	0	2343	36	0
8	BM	2320	0	2343	37	0
8	CM	2320	0	2343	39	0
8	DM	2320	0	2343	37	0
8	EM	2320	0	2343	41	0
8	FM	2320	0	2343	43	0
8	GM	2320	0	2343	36	0
8	HM	2320	0	2343	46	0
8	IM	2320	0	2343	41	0
8	JM	2320	0	2343	38	0
8	KM	2320	0	2343	47	0
8	LM	2320	0	2343	38	0
8	MM	2320	0	2343	40	0
9	AN	582	0	530	8	0
9	BN	582	0	530	9	0
9	CN	582	0	530	8	0
9	DN	582	0	530	7	0
9	EN	582	0	530	9	0
9	FN	582	0	530	8	0
9	GN	582	0	530	7	0
9	HN	582	0	530	7	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
9	IN	582	0	530	7	0
9	JN	582	0	530	6	0
9	KN	582	0	530	6	0
9	LN	582	0	530	6	0
9	MN	582	0	530	6	0
10	AU	45	0	12	0	0
10	BU	45	0	12	0	0
10	CU	45	0	12	0	0
10	DU	45	0	12	0	0
10	EU	45	0	12	0	0
10	FU	45	0	12	0	0
10	GU	45	0	12	0	0
10	HU	45	0	12	0	0
10	IU	45	0	12	0	0
10	JU	45	0	12	0	0
10	KU	45	0	12	0	0
10	LU	45	0	12	0	0
10	MU	45	0	12	0	0
11	AX	240	0	54	0	0
11	BX	240	0	54	0	0
11	CX	240	0	54	0	0
11	DX	240	0	54	0	0
11	EX	240	0	54	0	0
11	FX	240	0	54	0	0
11	GX	240	0	54	0	0
11	HX	240	0	54	0	0
11	IX	240	0	54	0	0
11	JX	240	0	54	0	0
11	KX	240	0	54	1	0
11	LX	240	0	54	0	0
11	MX	240	0	54	0	0
11	VX	240	0	54	0	0
11	WX	240	0	54	0	0
11	XX	240	0	54	0	0
11	YX	240	0	54	0	0
11	ZX	240	0	54	0	0
All	All	180380	0	178472	2286	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 6.

All (2286) 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:Fd:82:ARG:CB	2:Fd:82:ARG:CG	1.91	1.47
5:GH:115:MET:SD	5:GH:115:MET:CE	2.04	1.46
1:BC:177:CYS:HB2	1:CC:35:GLN:NE2	1.38	1.39
5:GH:115:MET:CE	5:GH:115:MET:HB2	1.71	1.21
1:BC:177:CYS:CB	1:CC:35:GLN:NE2	2.04	1.18
1:DC:157:LYS:NZ	1:EC:28:THR:OG1	1.77	1.17
2:Fd:82:ARG:CB	2:Fd:82:ARG:CD	2.24	1.14
8:HM:306:GLN:CG	3:Hf:253:ARG:NH2	2.13	1.12
5:GH:115:MET:HB2	5:GH:115:MET:HE3	1.31	1.05
1:BC:177:CYS:CB	1:CC:35:GLN:HE21	1.67	1.03
8:HM:306:GLN:HG2	3:Hf:253:ARG:NH2	1.73	1.02
8:HM:306:GLN:HG2	3:Hf:253:ARG:HH22	1.25	0.99
1:LC:176:TRP:CZ3	1:MC:35:GLN:NE2	2.31	0.99
8:HM:306:GLN:CG	3:Hf:253:ARG:HH21	1.77	0.96
8:HM:306:GLN:CD	3:Hf:253:ARG:HH21	1.72	0.96
1:CC:180:THR:HG21	1:DC:34:LEU:HD22	1.46	0.95
5:GH:115:MET:CE	5:GH:115:MET:CB	2.45	0.94
1:CC:157:LYS:NZ	1:DC:28:THR:OG1	1.99	0.94
1:FC:177:CYS:HB2	1:GC:35:GLN:HE22	1.30	0.94
1:DC:180:THR:HG21	1:EC:34:LEU:HD22	1.48	0.93
1:BC:177:CYS:HB2	1:CC:35:GLN:HE22	0.97	0.93
1:FC:177:CYS:CB	1:GC:35:GLN:HE22	1.80	0.93
1:GC:177:CYS:HB2	1:HC:35:GLN:NE2	1.86	0.90
2:Fd:82:ARG:CB	2:Fd:82:ARG:HD2	2.02	0.89
8:HM:306:GLN:NE2	3:Hf:253:ARG:HH21	1.70	0.89
8:KM:247:ARG:HD3	3:Kf:213:GLN:O	1.73	0.88
5:ZH:111:ALA:O	5:ZH:115:MET:HB2	1.76	0.86
1:BC:157:LYS:NZ	1:CC:28:THR:OG1	2.08	0.86
1:BC:180:THR:HG21	1:CC:34:LEU:HD22	1.58	0.85
1:BC:180:THR:HG21	1:CC:34:LEU:CD2	2.07	0.84
5:GH:115:MET:CE	5:GH:115:MET:CG	2.55	0.84
1:GC:177:CYS:HB2	1:HC:35:GLN:HE22	1.44	0.83
1:GC:157:LYS:HZ3	1:HC:28:THR:HB	1.40	0.83
8:HM:306:GLN:NE2	3:Hf:253:ARG:NH2	2.27	0.83
1:DC:157:LYS:CE	1:EC:28:THR:OG1	2.26	0.82
4:FG:818:GLN:NE2	5:FH:205:GLN:OE1	2.12	0.82
2:Fd:82:ARG:NH2	2:Fd:83:ALA:H	1.78	0.82
1:FC:177:CYS:HB2	1:GC:35:GLN:NE2	1.96	0.79
1:GC:180:THR:HG21	1:HC:34:LEU:HD22	1.65	0.79
1:CC:180:THR:HG21	1:DC:34:LEU:CD2	2.12	0.79
1:DC:180:THR:HG21	1:EC:34:LEU:CD2	2.12	0.78
1:BC:180:THR:CG2	1:CC:34:LEU:HD22	2.12	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:DD:24:LYS:O	7:DL:110:GLN:NE2	2.16	0.78
8:HM:306:GLN:HE21	3:Hf:253:ARG:NH2	1.82	0.78
4:DG:818:GLN:NE2	5:DH:205:GLN:OE1	2.18	0.77
5:GH:190:PRO:HB2	5:GH:231:ARG:HD2	1.67	0.77
5:FH:208:LYS:HD2	4:WG:823:GLY:HA2	1.64	0.77
1:CC:177:CYS:HB3	1:DC:35:GLN:HE22	1.48	0.77
1:GC:180:THR:HG21	1:HC:34:LEU:CD2	2.15	0.76
5:DH:238:MET:HG3	5:EH:250:TYR:HB3	1.68	0.76
2:HD:24:LYS:O	7:HL:110:GLN:NE2	2.18	0.76
1:GC:157:LYS:NZ	1:HC:28:THR:HB	2.01	0.75
5:WH:344:VAL:HG13	5:WH:355:GLN:HE21	1.51	0.75
3:BF:231:THR:HG21	5:CH:151:PRO:HG2	1.70	0.74
5:BH:115:MET:HE3	4:ZG:801:MET:HB3	1.69	0.74
5:EH:111:ALA:O	5:EH:115:MET:HB2	1.87	0.74
2:FD:107:ARG:HH11	2:FD:153:VAL:HG11	1.53	0.74
8:JM:247:ARG:HD3	3:Jf:213:GLN:O	1.88	0.74
2:Fd:82:ARG:HH21	2:Fd:83:ALA:H	1.35	0.74
1:GC:177:CYS:CB	1:HC:35:GLN:NE2	2.51	0.74
5:KH:266:MET:HE3	5:KH:267:PRO:HD2	1.69	0.73
4:AG:801:MET:HB3	5:CH:115:MET:HE3	1.68	0.73
1:AC:34:LEU:HD22	1:MC:180:THR:HG21	1.70	0.73
5:LH:236:PRO:HD2	5:MH:251:ARG:HH22	1.54	0.73
5:LH:263:ALA:HB1	5:LH:266:MET:HG3	1.71	0.73
1:IC:180:THR:HG21	1:JC:34:LEU:HD22	1.70	0.72
2:ED:107:ARG:NH1	2:Ed:130:GLU:OE2	2.22	0.72
2:ID:24:LYS:O	7:IL:110:GLN:NE2	2.22	0.72
11:KX:26:UNK:O	5:LH:132:GLN:NE2	2.14	0.72
5:AH:205:GLN:HB3	5:AH:214:MET:HB2	1.71	0.72
7:CL:187:GLU:OE2	8:CM:268:HIS:ND1	2.22	0.71
2:Fd:82:ARG:HD2	2:Fd:82:ARG:HB2	1.71	0.71
5:ZH:348:SER:HA	5:ZH:352:LYS:O	1.89	0.71
1:HC:180:THR:HG21	1:IC:34:LEU:HD22	1.71	0.71
5:WH:202:PHE:HB3	5:WH:215:ILE:HD11	1.72	0.71
4:AG:808:LEU:HD11	5:AH:141:LYS:HE3	1.72	0.71
1:CC:157:LYS:HZ3	1:DC:28:THR:CB	2.04	0.71
4:DG:797:ARG:CZ	5:EH:117:ARG:HH22	2.03	0.70
5:BH:279:LEU:HD11	5:BH:289:PRO:HB3	1.73	0.70
1:CC:176:TRP:CZ3	1:DC:35:GLN:OE1	2.45	0.70
2:FD:24:LYS:O	7:FL:110:GLN:NE2	2.24	0.70
5:GH:115:MET:HE3	5:GH:115:MET:CB	2.13	0.70
8:BM:177:LEU:HD13	8:BM:181:PRO:HG2	1.74	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:KM:177:LEU:HD13	8:KM:181:PRO:HG2	1.74	0.70
8:AM:177:LEU:HD13	8:AM:181:PRO:HG2	1.74	0.70
8:DM:177:LEU:HD13	8:DM:181:PRO:HG2	1.74	0.70
8:FM:177:LEU:HD13	8:FM:181:PRO:HG2	1.74	0.70
8:HM:177:LEU:HD13	8:HM:181:PRO:HG2	1.74	0.70
8:CM:177:LEU:HD13	8:CM:181:PRO:HG2	1.74	0.69
7:EL:187:GLU:OE2	8:EM:268:HIS:ND1	2.25	0.69
8:EM:177:LEU:HD13	8:EM:181:PRO:HG2	1.74	0.69
8:LM:177:LEU:HD13	8:LM:181:PRO:HG2	1.74	0.69
3:VF:229:THR:OG1	5:VH:165:THR:OG1	2.10	0.69
7:GL:187:GLU:OE2	8:GM:268:HIS:ND1	2.25	0.69
5:KH:189:TRP:HE1	5:KH:232:GLY:H	1.39	0.69
2:BD:107:ARG:HH11	2:BD:153:VAL:HG11	1.58	0.69
2:Hd:124:LYS:NZ	2:ID:54:GLU:OE2	2.25	0.69
3:JF:214:ALA:HA	5:JH:245:GLN:HE21	1.56	0.69
8:IM:177:LEU:HD13	8:IM:181:PRO:HG2	1.74	0.69
8:MM:177:LEU:HD13	8:MM:181:PRO:HG2	1.74	0.69
8:FM:306:GLN:OE1	3:Ff:251:GLN:HG2	1.92	0.69
1:MC:35:GLN:HA	1:MC:35:GLN:HE21	1.58	0.69
2:Fd:82:ARG:CB	2:Fd:82:ARG:NE	2.55	0.69
5:FH:253:ASP:HB3	5:FH:255:ARG:HH22	1.57	0.69
4:DG:791:GLN:N	4:DG:793:GLU:OE1	2.27	0.68
2:Dd:124:LYS:NZ	2:ED:54:GLU:OE2	2.22	0.68
4:AG:791:GLN:N	4:AG:793:GLU:OE1	2.27	0.68
8:JM:177:LEU:HD13	8:JM:181:PRO:HG2	1.74	0.68
4:ZG:791:GLN:N	4:ZG:793:GLU:OE1	2.27	0.68
4:EG:791:GLN:N	4:EG:793:GLU:OE1	2.27	0.68
8:GM:177:LEU:HD13	8:GM:181:PRO:HG2	1.74	0.68
8:HM:306:GLN:HG3	3:Hf:253:ARG:NH2	2.06	0.68
4:XG:791:GLN:N	4:XG:793:GLU:OE1	2.27	0.68
1:BC:177:CYS:CA	1:CC:35:GLN:HE21	2.06	0.68
4:BG:791:GLN:N	4:BG:793:GLU:OE1	2.27	0.68
2:ED:77:TYR:OH	6:EK:130:ARG:NH1	2.27	0.68
4:GG:791:GLN:N	4:GG:793:GLU:OE1	2.27	0.68
4:YG:791:GLN:N	4:YG:793:GLU:OE1	2.27	0.68
8:AM:306:GLN:OE1	3:Af:251:GLN:HG2	1.92	0.68
4:CG:791:GLN:N	4:CG:793:GLU:OE1	2.27	0.68
4:IG:791:GLN:N	4:IG:793:GLU:OE1	2.27	0.68
1:LC:176:TRP:HZ3	1:MC:35:GLN:NE2	1.89	0.68
4:FG:791:GLN:N	4:FG:793:GLU:OE1	2.27	0.68
4:JG:791:GLN:N	4:JG:793:GLU:OE1	2.27	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:MG:791:GLN:N	4:MG:793:GLU:OE1	2.27	0.68
4:WG:791:GLN:N	4:WG:793:GLU:OE1	2.27	0.68
2:Fd:82:ARG:NH2	2:Fd:83:ALA:O	2.25	0.67
4:LG:791:GLN:N	4:LG:793:GLU:OE1	2.27	0.67
2:BD:107:ARG:NE	2:Bd:130:GLU:OE2	2.25	0.67
1:GC:57:MET:N	1:GC:57:MET:SD	2.68	0.67
4:HG:801:MET:HB3	5:IH:115:MET:HE3	1.74	0.67
5:WH:328:MET:HE3	5:WH:329:THR:H	1.59	0.67
1:DC:57:MET:N	1:DC:57:MET:SD	2.68	0.67
1:FC:57:MET:N	1:FC:57:MET:SD	2.68	0.67
5:HH:208:LYS:HD2	4:XG:823:GLY:HA2	1.76	0.67
5:JH:208:LYS:HD2	4:KG:823:GLY:HA2	1.77	0.67
4:KG:791:GLN:N	4:KG:793:GLU:OE1	2.27	0.67
7:LL:187:GLU:OE2	8:LM:268:HIS:ND1	2.26	0.67
4:VG:791:GLN:N	4:VG:793:GLU:OE1	2.27	0.67
1:AC:57:MET:N	1:AC:57:MET:SD	2.68	0.67
1:FC:69:GLN:HE22	9:FN:114:LEU:HA	1.59	0.67
4:HG:791:GLN:N	4:HG:793:GLU:OE1	2.27	0.67
5:IH:230:LEU:HB2	5:IH:233:LEU:HD13	1.75	0.67
5:IH:311:MET:HE3	5:IH:341:LYS:HA	1.75	0.67
2:GD:24:LYS:O	7:GL:110:GLN:NE2	2.27	0.67
8:KM:227:ILE:HD11	3:Kf:213:GLN:HB3	1.76	0.67
1:BC:57:MET:N	1:BC:57:MET:SD	2.68	0.67
1:BC:66:VAL:HG22	1:CC:34:LEU:HD11	1.77	0.67
7:DL:124:THR:HG22	7:DL:232:GLN:HG2	1.77	0.67
8:EM:306:GLN:OE1	3:Ef:251:GLN:HG2	1.95	0.67
1:LC:57:MET:N	1:LC:57:MET:SD	2.68	0.67
5:LH:238:MET:HE1	5:MH:250:TYR:CD2	2.29	0.67
7:BL:124:THR:HG22	7:BL:232:GLN:HG2	1.77	0.67
1:MC:57:MET:SD	1:MC:57:MET:N	2.68	0.67
5:AH:115:MET:HE1	4:MG:802:LEU:HA	1.77	0.67
5:AH:158:VAL:HG23	5:AH:254:LEU:HD12	1.75	0.67
1:EC:57:MET:SD	1:EC:57:MET:N	2.68	0.67
1:IC:57:MET:N	1:IC:57:MET:SD	2.68	0.67
7:ML:124:THR:HG22	7:ML:232:GLN:HG2	1.77	0.67
7:AL:124:THR:HG22	7:AL:232:GLN:HG2	1.77	0.67
1:EC:157:LYS:HZ3	1:FC:28:THR:HB	1.58	0.67
7:EL:124:THR:HG22	7:EL:232:GLN:HG2	1.77	0.67
1:HC:57:MET:SD	1:HC:57:MET:N	2.68	0.67
4:HG:801:MET:HB3	5:IH:115:MET:CE	2.25	0.67
7:IL:124:THR:HG22	7:IL:232:GLN:HG2	1.77	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:KL:187:GLU:OE2	8:KM:268:HIS:ND1	2.28	0.67
5:ZH:349:TRP:HB3	5:ZH:352:LYS:HE3	1.77	0.67
5:BH:169:ILE:HD11	5:BH:239:LEU:HD23	1.77	0.66
1:CC:57:MET:N	1:CC:57:MET:SD	2.68	0.66
5:CH:245:GLN:HE21	5:CH:246:LYS:H	1.43	0.66
7:FL:124:THR:HG22	7:FL:232:GLN:HG2	1.77	0.66
7:JL:124:THR:HG22	7:JL:232:GLN:HG2	1.77	0.66
5:AH:229:ARG:HG2	5:AH:236:PRO:HB3	1.76	0.66
2:CD:130:GLU:OE2	2:CD:133:ARG:NH1	2.28	0.66
1:CC:66:VAL:HG22	1:DC:34:LEU:HD11	1.77	0.66
1:JC:57:MET:N	1:JC:57:MET:SD	2.68	0.66
2:Ed:98:ILE:HG23	2:Ed:128:LEU:HD12	1.78	0.66
1:KC:57:MET:N	1:KC:57:MET:SD	2.68	0.66
7:CL:124:THR:HG22	7:CL:232:GLN:HG2	1.77	0.66
5:JH:238:MET:HE1	5:KH:250:TYR:CD2	2.31	0.66
2:Jd:93:GLU:OE2	1:KC:264:ALA:N	2.28	0.66
2:MD:24:LYS:O	7:ML:110:GLN:NE2	2.28	0.66
7:HL:124:THR:HG22	7:HL:232:GLN:HG2	1.77	0.66
1:GC:97:ASN:HD22	1:GC:97:ASN:C	2.03	0.65
7:LL:124:THR:HG22	7:LL:232:GLN:HG2	1.77	0.65
3:Ff:219:ARG:HD3	3:Ff:233:ARG:HH11	1.60	0.65
8:IM:227:ILE:HD11	3:If:213:GLN:HG2	1.79	0.65
1:HC:69:GLN:HE22	9:HN:114:LEU:HA	1.59	0.65
7:KL:124:THR:HG22	7:KL:232:GLN:HG2	1.77	0.65
5:GH:169:ILE:HD12	5:GH:252:VAL:HG21	1.78	0.65
7:GL:124:THR:HG22	7:GL:232:GLN:HG2	1.77	0.65
5:HH:159:ASN:OD1	5:HH:257:GLN:NE2	2.29	0.65
3:GF:231:THR:O	5:HH:150:LYS:NZ	2.29	0.65
8:FM:306:GLN:CD	3:Ff:251:GLN:HG2	2.20	0.65
6:GK:129:GLU:OE1	6:GK:129:GLU:N	2.30	0.65
6:CK:129:GLU:N	6:CK:129:GLU:OE1	2.30	0.65
8:DM:306:GLN:OE1	3:Df:251:GLN:HG2	1.97	0.65
8:JM:158:VAL:HA	8:JM:178:LYS:HZ3	1.62	0.64
5:KH:311:MET:HE2	5:KH:341:LYS:HA	1.78	0.64
2:AD:98:ILE:HG23	2:AD:128:LEU:HD12	1.79	0.64
6:AK:129:GLU:N	6:AK:129:GLU:OE1	2.30	0.64
8:FM:158:VAL:HA	8:FM:178:LYS:HZ3	1.62	0.64
8:GM:158:VAL:HA	8:GM:178:LYS:HZ3	1.62	0.64
3:IF:229:THR:OG1	5:IH:165:THR:OG1	2.15	0.64
1:KC:97:ASN:C	1:KC:97:ASN:HD22	2.03	0.64
5:KH:111:ALA:O	5:KH:115:MET:HB2	1.97	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:IK:129:GLU:N	6:IK:129:GLU:OE1	2.30	0.64
2:JD:127:SER:OG	2:JD:130:GLU:OE1	2.15	0.64
8:KM:158:VAL:HA	8:KM:178:LYS:HZ3	1.62	0.64
5:LH:189:TRP:CD1	5:LH:230:LEU:HG	2.32	0.64
6:MK:129:GLU:OE1	6:MK:129:GLU:N	2.30	0.64
3:XF:241:TYR:HB3	3:XF:256:THR:HG21	1.80	0.64
6:EK:129:GLU:OE1	6:EK:129:GLU:N	2.30	0.64
8:LM:158:VAL:HA	8:LM:178:LYS:HZ3	1.62	0.64
8:LM:227:ILE:HD11	3:Lf:213:GLN:HG2	1.79	0.64
1:FC:97:ASN:C	1:FC:97:ASN:HD22	2.03	0.64
8:MM:158:VAL:HA	8:MM:178:LYS:HZ3	1.62	0.64
8:DM:158:VAL:HA	8:DM:178:LYS:HZ3	1.62	0.64
3:EF:241:TYR:HB3	3:EF:256:THR:HG21	1.80	0.64
8:HM:158:VAL:HA	8:HM:178:LYS:HZ3	1.63	0.64
3:LF:241:TYR:HB3	3:LF:256:THR:HG21	1.80	0.64
3:WF:241:TYR:HB3	3:WF:256:THR:HG21	1.80	0.64
3:BF:241:TYR:HB3	3:BF:256:THR:HG21	1.80	0.64
8:BM:158:VAL:HA	8:BM:178:LYS:HZ3	1.63	0.64
6:DK:129:GLU:OE1	6:DK:129:GLU:N	2.30	0.64
8:EM:158:VAL:HA	8:EM:178:LYS:HZ3	1.62	0.64
7:HL:187:GLU:OE2	8:HM:268:HIS:ND1	2.30	0.64
6:KK:129:GLU:OE1	6:KK:129:GLU:N	2.30	0.64
5:BH:179:LEU:O	5:BH:212:THR:HA	1.98	0.64
3:FF:241:TYR:HB3	3:FF:256:THR:HG21	1.80	0.64
3:MF:241:TYR:HB3	3:MF:256:THR:HG21	1.80	0.64
8:AM:158:VAL:HA	8:AM:178:LYS:HZ3	1.62	0.64
3:HF:241:TYR:HB3	3:HF:256:THR:HG21	1.80	0.64
5:MH:191:ILE:HG13	5:MH:211:ASN:HA	1.79	0.64
1:AC:35:GLN:HE21	1:MC:177:CYS:HA	1.63	0.63
3:AF:241:TYR:HB3	3:AF:256:THR:HG21	1.80	0.63
1:AC:97:ASN:HD22	1:AC:97:ASN:C	2.03	0.63
3:CF:241:TYR:HB3	3:CF:256:THR:HG21	1.80	0.63
6:HK:129:GLU:OE1	6:HK:129:GLU:N	2.30	0.63
8:IM:58:VAL:HG22	8:IM:106:VAL:HB	1.81	0.63
6:JK:129:GLU:OE1	6:JK:129:GLU:N	2.30	0.63
5:YH:283:VAL:HG21	5:YH:312:TYR:HB3	1.80	0.63
1:DC:66:VAL:HG22	1:EC:34:LEU:HD11	1.80	0.63
8:HM:58:VAL:HG22	8:HM:106:VAL:HB	1.81	0.63
2:LD:107:ARG:NH2	2:LD:155:GLU:OE2	2.32	0.63
5:YH:295:LEU:HD11	5:YH:304:ALA:HB3	1.79	0.63
3:ZF:241:TYR:HB3	3:ZF:256:THR:HG21	1.80	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:DF:241:TYR:HB3	3:DF:256:THR:HG21	1.80	0.63
8:DM:58:VAL:HG22	8:DM:106:VAL:HB	1.81	0.63
7:HL:165:VAL:HG12	7:HL:231:ILE:HG12	1.81	0.63
1:IC:97:ASN:HD22	1:IC:97:ASN:C	2.03	0.63
3:IF:241:TYR:HB3	3:IF:256:THR:HG21	1.80	0.63
3:VF:241:TYR:HB3	3:VF:256:THR:HG21	1.80	0.63
3:YF:241:TYR:HB3	3:YF:256:THR:HG21	1.80	0.63
5:IH:150:LYS:HD2	3:XF:219:ARG:HH22	1.63	0.63
7:IL:165:VAL:HG12	7:IL:231:ILE:HG12	1.81	0.63
8:IM:158:VAL:HA	8:IM:178:LYS:HZ3	1.62	0.63
7:AL:63:ASP:O	7:AL:88:GLY:N	2.32	0.63
8:BM:58:VAL:HG22	8:BM:106:VAL:HB	1.81	0.63
1:CC:180:THR:CG2	1:DC:34:LEU:HD22	2.24	0.63
1:DC:157:LYS:HE2	1:EC:28:THR:OG1	1.99	0.63
8:GM:58:VAL:HG22	8:GM:106:VAL:HB	1.81	0.63
1:BC:177:CYS:HB3	1:CC:35:GLN:HE21	1.61	0.63
7:FL:165:VAL:HG12	7:FL:231:ILE:HG12	1.81	0.63
7:HL:63:ASP:O	7:HL:88:GLY:N	2.32	0.63
7:GL:63:ASP:O	7:GL:88:GLY:N	2.32	0.62
3:KF:241:TYR:HB3	3:KF:256:THR:HG21	1.80	0.62
3:Hf:220:ALA:HB3	3:Hf:232:VAL:HB	1.81	0.62
6:LK:129:GLU:OE1	6:LK:129:GLU:N	2.30	0.62
7:ML:165:VAL:HG12	7:ML:231:ILE:HG12	1.81	0.62
8:AM:58:VAL:HG22	8:AM:106:VAL:HB	1.81	0.62
6:BK:129:GLU:OE1	6:BK:129:GLU:N	2.30	0.62
7:BL:63:ASP:O	7:BL:88:GLY:N	2.32	0.62
7:DL:63:ASP:O	7:DL:88:GLY:N	2.32	0.62
7:EL:63:ASP:O	7:EL:88:GLY:N	2.32	0.62
8:MM:58:VAL:HG22	8:MM:106:VAL:HB	1.81	0.62
7:FL:63:ASP:O	7:FL:88:GLY:N	2.32	0.62
1:GC:177:CYS:CB	1:HC:35:GLN:HE21	2.12	0.62
3:JF:241:TYR:HB3	3:JF:256:THR:HG21	1.80	0.62
8:JM:58:VAL:HG22	8:JM:106:VAL:HB	1.81	0.62
7:KL:63:ASP:O	7:KL:88:GLY:N	2.32	0.62
5:LH:155:SER:OG	5:LH:255:ARG:NH1	2.33	0.62
7:LL:63:ASP:O	7:LL:88:GLY:N	2.32	0.62
5:ZH:223:TYR:HB2	5:ZH:242:ILE:HG22	1.81	0.62
5:ZH:280:LEU:HA	5:ZH:283:VAL:HG22	1.81	0.62
8:AM:40:GLN:HE22	8:AM:42:ARG:HH21	1.48	0.62
3:EF:229:THR:OG1	5:EH:165:THR:OG1	2.15	0.62
7:ML:63:ASP:O	7:ML:88:GLY:N	2.32	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:FK:129:GLU:N	6:FK:129:GLU:OE1	2.30	0.62
3:GF:241:TYR:HB3	3:GF:256:THR:HG21	1.80	0.62
7:IL:63:ASP:O	7:IL:88:GLY:N	2.32	0.62
8:KM:58:VAL:HG22	8:KM:106:VAL:HB	1.81	0.62
4:LG:817:THR:HA	5:LH:216:GLN:HE22	1.65	0.62
7:LL:165:VAL:HG12	7:LL:231:ILE:HG12	1.81	0.62
1:BC:109:VAL:HG13	1:BC:208:GLY:HA3	1.82	0.62
2:Bd:92:GLU:OE2	2:Bd:158:TYR:OH	2.18	0.62
1:CC:177:CYS:HB3	1:DC:35:GLN:NE2	2.15	0.62
7:CL:165:VAL:HG12	7:CL:231:ILE:HG12	1.81	0.62
8:CM:58:VAL:HG22	8:CM:106:VAL:HB	1.81	0.62
8:EM:40:GLN:HE22	8:EM:42:ARG:HH21	1.48	0.62
7:JL:63:ASP:O	7:JL:88:GLY:N	2.32	0.62
5:ZH:283:VAL:HG12	5:ZH:289:PRO:HD3	1.81	0.62
1:KC:109:VAL:HG13	1:KC:208:GLY:HA3	1.82	0.62
7:KL:165:VAL:HG12	7:KL:231:ILE:HG12	1.81	0.62
8:KM:247:ARG:HD2	3:Kf:264:PHE:CE1	2.35	0.62
7:AL:165:VAL:HG12	7:AL:231:ILE:HG12	1.81	0.62
3:ZF:254:ILE:HB	3:ZF:262:ILE:HB	1.82	0.62
5:GH:160:LEU:HD21	5:GH:235:THR:HB	1.80	0.62
2:Hd:134:ASP:OD1	2:Hd:138:GLN:NE2	2.30	0.62
7:JL:165:VAL:HG12	7:JL:231:ILE:HG12	1.81	0.62
1:BC:35:GLN:HA	1:BC:35:GLN:HE21	1.63	0.61
8:FM:58:VAL:HG22	8:FM:106:VAL:HB	1.81	0.61
3:GF:254:ILE:HB	3:GF:262:ILE:HB	1.82	0.61
8:JM:40:GLN:HE22	8:JM:42:ARG:HH21	1.48	0.61
1:LC:109:VAL:HG13	1:LC:208:GLY:HA3	1.82	0.61
8:LM:58:VAL:HG22	8:LM:106:VAL:HB	1.81	0.61
7:BL:165:VAL:HG12	7:BL:231:ILE:HG12	1.81	0.61
8:CM:40:GLN:HE22	8:CM:42:ARG:HH21	1.48	0.61
8:CM:158:VAL:HA	8:CM:178:LYS:HZ3	1.63	0.61
1:EC:69:GLN:HE22	9:EN:114:LEU:HA	1.63	0.61
1:IC:109:VAL:HG13	1:IC:208:GLY:HA3	1.82	0.61
7:ML:187:GLU:OE2	8:MM:268:HIS:ND1	2.29	0.61
1:AC:109:VAL:HG13	1:AC:208:GLY:HA3	1.82	0.61
1:CC:109:VAL:HG13	1:CC:208:GLY:HA3	1.82	0.61
5:DH:225:ASN:C	5:DH:225:ASN:HD22	2.08	0.61
7:EL:165:VAL:HG12	7:EL:231:ILE:HG12	1.81	0.61
5:FH:324:TRP:HB3	5:FH:339:MET:HE2	1.83	0.61
3:HF:254:ILE:HB	3:HF:262:ILE:HB	1.82	0.61
3:AF:254:ILE:HB	3:AF:262:ILE:HB	1.82	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:EC:109:VAL:HG13	1:EC:208:GLY:HA3	1.82	0.61
3:EF:254:ILE:HB	3:EF:262:ILE:HB	1.82	0.61
8:GM:40:GLN:HE22	8:GM:42:ARG:HH21	1.48	0.61
1:HC:109:VAL:HG13	1:HC:208:GLY:HA3	1.82	0.61
8:KM:40:GLN:HE22	8:KM:42:ARG:HH21	1.48	0.61
3:YF:254:ILE:HB	3:YF:262:ILE:HB	1.82	0.61
7:CL:63:ASP:O	7:CL:88:GLY:N	2.32	0.61
7:DL:165:VAL:HG12	7:DL:231:ILE:HG12	1.81	0.61
1:FC:109:VAL:HG13	1:FC:208:GLY:HA3	1.82	0.61
8:LM:40:GLN:HE22	8:LM:42:ARG:HH21	1.48	0.61
5:ZH:230:LEU:HD12	5:ZH:233:LEU:HD12	1.82	0.61
1:DC:69:GLN:HE22	9:DN:114:LEU:HA	1.65	0.61
3:Hf:219:ARG:HH22	3:Hf:233:ARG:HD2	1.65	0.61
5:LH:158:VAL:HG23	5:LH:254:LEU:HD12	1.82	0.61
5:VH:322:PRO:HG2	5:VH:345:LEU:HD23	1.81	0.61
3:CF:254:ILE:HB	3:CF:262:ILE:HB	1.82	0.61
3:DF:254:ILE:HB	3:DF:262:ILE:HB	1.82	0.61
1:GC:109:VAL:HG13	1:GC:208:GLY:HA3	1.82	0.61
5:KH:190:PRO:HB2	5:KH:231:ARG:HG3	1.82	0.61
3:XF:254:ILE:HB	3:XF:262:ILE:HB	1.82	0.61
2:AD:85:VAL:HG12	6:AK:153:ILE:HG23	1.83	0.61
2:CD:24:LYS:O	7:CL:110:GLN:NE2	2.34	0.61
8:DM:40:GLN:HE22	8:DM:42:ARG:HH21	1.48	0.61
8:EM:58:VAL:HG22	8:EM:106:VAL:HB	1.81	0.61
3:KF:254:ILE:HB	3:KF:262:ILE:HB	1.82	0.61
1:BC:180:THR:CG2	1:CC:34:LEU:CD2	2.75	0.60
5:BH:110:LYS:HA	5:BH:113:LYS:HG2	1.83	0.60
8:BM:40:GLN:HE22	8:BM:42:ARG:HH21	1.48	0.60
8:HM:40:GLN:HE22	8:HM:42:ARG:HH21	1.48	0.60
8:JM:306:GLN:HE21	3:Jf:253:ARG:HH21	1.47	0.60
3:IF:254:ILE:HB	3:IF:262:ILE:HB	1.82	0.60
1:JC:109:VAL:HG13	1:JC:208:GLY:HA3	1.82	0.60
3:MF:254:ILE:HB	3:MF:262:ILE:HB	1.82	0.60
5:CH:173:GLN:HA	5:CH:217:ALA:HB3	1.83	0.60
7:GL:165:VAL:HG12	7:GL:231:ILE:HG12	1.81	0.60
8:GM:306:GLN:HE21	3:Gf:253:ARG:HH21	1.48	0.60
4:LG:818:GLN:H	5:LH:216:GLN:HE22	1.49	0.60
3:CF:213:GLN:O	5:CH:170:ARG:NH2	2.33	0.60
1:DC:109:VAL:HG13	1:DC:208:GLY:HA3	1.82	0.60
3:FF:254:ILE:HB	3:FF:262:ILE:HB	1.82	0.60
5:GH:248:VAL:HG13	5:GH:250:TYR:CE2	2.36	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:IM:40:GLN:HE22	8:IM:42:ARG:HH21	1.48	0.60
3:LF:254:ILE:HB	3:LF:262:ILE:HB	1.82	0.60
8:MM:306:GLN:OE1	3:Mf:251:GLN:HG2	2.02	0.60
1:LC:180:THR:HG21	1:MC:34:LEU:HD22	1.84	0.60
3:BF:254:ILE:HB	3:BF:262:ILE:HB	1.82	0.60
3:WF:254:ILE:HB	3:WF:262:ILE:HB	1.82	0.60
3:Hf:233:ARG:HB3	3:Hf:236:SER:HB3	1.84	0.60
1:MC:109:VAL:HG13	1:MC:208:GLY:HA3	1.82	0.60
8:FM:40:GLN:HE22	8:FM:42:ARG:HH21	1.48	0.60
3:VF:254:ILE:HB	3:VF:262:ILE:HB	1.82	0.60
4:CG:818:GLN:NE2	5:CH:205:GLN:OE1	2.35	0.60
4:DG:818:GLN:HG3	5:DH:205:GLN:HE22	1.67	0.60
9:FN:63:GLY:HA3	2:GD:27:LYS:HB2	1.82	0.60
8:MM:40:GLN:HE22	8:MM:42:ARG:HH21	1.48	0.60
5:ZH:205:GLN:HB2	5:ZH:214:MET:HB2	1.83	0.60
7:BL:187:GLU:OE2	8:BM:268:HIS:ND1	2.35	0.59
1:DC:180:THR:CG2	1:EC:34:LEU:HD22	2.28	0.59
2:DD:130:GLU:OE2	2:DD:133:ARG:NH1	2.33	0.59
2:KD:24:LYS:O	7:KL:110:GLN:NE2	2.35	0.59
2:ED:24:LYS:O	7:EL:110:GLN:NE2	2.36	0.59
1:GC:180:THR:CG2	1:HC:34:LEU:HD22	2.32	0.59
5:HH:221:TYR:OH	5:XH:135:GLU:OE2	2.13	0.59
2:AD:29:PRO:HG2	7:AL:225:GLN:HA	1.84	0.59
5:AH:128:VAL:HG12	5:AH:132:GLN:HE22	1.66	0.59
3:JF:254:ILE:HB	3:JF:262:ILE:HB	1.82	0.59
3:Lf:221:TRP:HE1	3:Lf:229:THR:HB	1.68	0.59
5:VH:357:LYS:NZ	5:VH:359:GLU:OE2	2.35	0.59
1:JC:154:ASP:HB2	9:JN:111:GLU:CD	2.27	0.59
5:VH:204:ILE:HD11	5:VH:213:LEU:HD22	1.84	0.59
2:ED:85:VAL:HG12	6:EK:153:ILE:HG23	1.85	0.59
5:BH:115:MET:HE1	4:ZG:802:LEU:HD12	1.84	0.59
4:IG:802:LEU:HA	5:KH:115:MET:HE1	1.85	0.59
5:IH:150:LYS:HD2	3:XF:219:ARG:HH12	1.67	0.59
2:LD:29:PRO:HG2	7:LL:225:GLN:HA	1.85	0.59
2:ED:130:GLU:OE1	2:ED:133:ARG:NH1	2.36	0.59
2:JD:29:PRO:HG2	7:JL:225:GLN:HA	1.85	0.58
1:BC:177:CYS:HB3	1:CC:35:GLN:NE2	2.13	0.58
2:Fd:24:LYS:HE2	2:Fd:25:PHE:HE1	1.68	0.58
2:KD:72:THR:HG23	2:KD:73:ILE:HG13	1.85	0.58
2:AD:55:MET:HE1	2:Ad:55:MET:HE1	1.86	0.58
5:AH:204:ILE:HD11	5:AH:213:LEU:HD22	1.86	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:GF:213:GLN:O	5:GH:170:ARG:NE	2.29	0.58
2:Gd:83:ALA:HA	1:HC:267:ALA:O	2.03	0.58
2:ID:130:GLU:OE2	2:ID:133:ARG:NH1	2.34	0.58
5:HH:204:ILE:HD11	5:HH:213:LEU:HD22	1.85	0.58
5:JH:204:ILE:HD11	5:JH:213:LEU:HD22	1.85	0.58
3:Mf:221:TRP:HE1	3:Mf:229:THR:HB	1.68	0.58
5:VH:230:LEU:HD12	5:VH:233:LEU:HD12	1.86	0.58
5:EH:311:MET:HE2	5:EH:341:LYS:HA	1.86	0.58
5:LH:311:MET:HE3	5:LH:341:LYS:HA	1.85	0.58
5:XH:320:LEU:HD11	5:XH:348:SER:HB3	1.86	0.58
1:AC:126:ARG:HD3	5:AH:275:SER:HA	1.85	0.58
2:Ed:64:PRO:HG2	2:Ed:67:LYS:HZ3	1.68	0.58
1:EC:210:ILE:HD11	2:ED:52:MET:HE1	1.85	0.58
6:GK:71:GLY:O	2:HD:27:LYS:NZ	2.35	0.58
6:JK:71:GLY:O	2:KD:27:LYS:NZ	2.36	0.58
5:XH:295:LEU:HG	5:XH:360:GLY:HA2	1.86	0.58
5:ZH:159:ASN:OD1	5:ZH:257:GLN:NE2	2.37	0.58
2:ID:77:TYR:OH	6:IK:130:ARG:NH1	2.37	0.58
5:WH:179:LEU:O	5:WH:212:THR:HA	2.04	0.58
5:GH:169:ILE:HB	5:GH:241:LEU:HD13	1.86	0.58
2:KD:29:PRO:HG2	7:KL:225:GLN:HA	1.85	0.58
4:MG:802:LEU:O	4:MG:806:THR:OG1	2.22	0.58
4:WG:802:LEU:O	4:WG:806:THR:OG1	2.22	0.58
2:AD:107:ARG:NH1	2:Ad:130:GLU:OE2	2.38	0.57
5:AH:137:SER:O	5:AH:141:LYS:HG2	2.04	0.57
7:AL:187:GLU:OE2	8:AM:268:HIS:ND1	2.34	0.57
1:GC:69:GLN:HE22	9:GN:114:LEU:HA	1.69	0.57
8:IM:306:GLN:OE1	3:If:251:GLN:HG2	2.04	0.57
2:BD:77:TYR:OH	6:BK:130:ARG:NH1	2.35	0.57
5:EH:145:PRO:HB3	5:FH:132:GLN:HE22	1.68	0.57
4:GG:818:GLN:NE2	5:GH:216:GLN:OE1	2.37	0.57
1:IC:180:THR:HG21	1:JC:34:LEU:CD2	2.35	0.57
5:JH:192:ALA:HA	5:JH:231:ARG:HH21	1.69	0.57
8:BM:119:GLU:HG3	8:BM:139:ALA:HB3	1.87	0.57
4:BG:802:LEU:O	4:BG:806:THR:OG1	2.22	0.57
2:GD:141:LYS:NZ	2:Gd:59:GLU:OE2	2.34	0.57
2:Md:64:PRO:HD2	2:Md:67:LYS:HE3	1.85	0.57
2:Dd:112:SER:HA	2:Dd:158:TYR:HE2	1.69	0.57
1:JC:240:ILE:HG12	1:KC:255:LEU:HD13	1.86	0.57
4:KG:802:LEU:O	4:KG:806:THR:OG1	2.22	0.57
6:AK:71:GLY:O	2:BD:27:LYS:NZ	2.37	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:FH:251:ARG:NH2	5:FH:253:ASP:OD1	2.38	0.57
8:LM:119:GLU:HG3	8:LM:139:ALA:HB3	1.87	0.57
5:MH:189:TRP:HE1	5:MH:230:LEU:HB3	1.70	0.57
4:CG:795:GLN:OE1	4:CG:795:GLN:N	2.35	0.57
8:HM:119:GLU:HG3	8:HM:139:ALA:HB3	1.87	0.57
8:KM:247:ARG:CD	3:Kf:213:GLN:O	2.48	0.57
5:VH:190:PRO:HB2	5:VH:231:ARG:HG3	1.86	0.57
1:AC:35:GLN:NE2	1:MC:176:TRP:CE3	2.73	0.57
4:DG:802:LEU:O	4:DG:806:THR:OG1	2.22	0.57
4:FG:802:LEU:O	4:FG:806:THR:OG1	2.22	0.57
6:LK:116:VAL:HG22	6:LK:182:ILE:HG12	1.87	0.57
5:AH:357:LYS:HB2	9:MN:81:LEU:HD21	1.87	0.56
4:CG:802:LEU:O	4:CG:806:THR:OG1	2.22	0.56
6:DK:116:VAL:HG22	6:DK:182:ILE:HG12	1.87	0.56
8:GM:306:GLN:OE1	3:Gf:251:GLN:HG2	2.05	0.56
1:IC:177:CYS:HB3	1:JC:35:GLN:NE2	2.19	0.56
4:JG:802:LEU:O	4:JG:806:THR:OG1	2.22	0.56
8:KM:119:GLU:HG3	8:KM:139:ALA:HB3	1.87	0.56
5:VH:226:LEU:HD23	5:VH:241:LEU:HD11	1.86	0.56
4:AG:802:LEU:O	4:AG:806:THR:OG1	2.22	0.56
5:DH:311:MET:HE2	5:DH:341:LYS:HA	1.87	0.56
9:EN:81:LEU:HD21	5:FH:357:LYS:HB2	1.86	0.56
6:HK:116:VAL:HG22	6:HK:182:ILE:HG12	1.88	0.56
2:ID:85:VAL:HG12	6:IK:153:ILE:HG23	1.87	0.56
2:Id:69:ASN:HD22	2:Id:69:ASN:C	2.13	0.56
1:JC:243:ASP:O	1:KC:128:SER:HA	2.05	0.56
4:XG:802:LEU:O	4:XG:806:THR:OG1	2.22	0.56
8:CM:119:GLU:HG3	8:CM:139:ALA:HB3	1.87	0.56
4:EG:802:LEU:O	4:EG:806:THR:OG1	2.22	0.56
4:GG:802:LEU:O	4:GG:806:THR:OG1	2.22	0.56
4:HG:802:LEU:O	4:HG:806:THR:OG1	2.22	0.56
2:MD:29:PRO:HG2	7:ML:225:GLN:HA	1.87	0.56
3:ZF:214:ALA:HB1	5:ZH:242:ILE:HD11	1.86	0.56
4:DG:795:GLN:OE1	4:DG:795:GLN:N	2.35	0.56
6:DK:71:GLY:O	2:ED:27:LYS:NZ	2.35	0.56
5:EH:169:ILE:HD11	5:EH:239:LEU:HD12	1.87	0.56
6:IK:116:VAL:HG22	6:IK:182:ILE:HG12	1.87	0.56
2:Md:134:ASP:OD1	2:Md:138:GLN:NE2	2.39	0.56
4:YG:802:LEU:O	4:YG:806:THR:OG1	2.22	0.56
5:CH:230:LEU:HD12	5:CH:233:LEU:HD12	1.86	0.56
8:GM:119:GLU:HG3	8:GM:139:ALA:HB3	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:Gd:82:ARG:O	1:HC:268:LYS:HA	2.06	0.56
4:AG:802:LEU:HD12	5:CH:115:MET:HE1	1.88	0.56
6:AK:116:VAL:HG22	6:AK:182:ILE:HG12	1.87	0.56
6:CK:116:VAL:HG22	6:CK:182:ILE:HG12	1.87	0.56
5:DH:250:TYR:HB3	5:VH:238:MET:HE1	1.87	0.56
6:EK:116:VAL:HG22	6:EK:182:ILE:HG12	1.87	0.56
8:EM:119:GLU:HG3	8:EM:139:ALA:HB3	1.87	0.56
2:KD:107:ARG:NH1	2:Kd:130:GLU:OE2	2.38	0.56
6:MK:116:VAL:HG22	6:MK:182:ILE:HG12	1.88	0.56
4:VG:802:LEU:O	4:VG:806:THR:OG1	2.22	0.56
4:VG:807:GLN:OE1	4:VG:810:GLN:NE2	2.37	0.56
8:AM:119:GLU:HG3	8:AM:139:ALA:HB3	1.87	0.56
8:FM:32:LYS:HB3	8:FM:101:TYR:HA	1.88	0.56
6:IK:71:GLY:O	2:JD:27:LYS:NZ	2.38	0.56
8:IM:119:GLU:HG3	8:IM:139:ALA:HB3	1.87	0.56
8:MM:32:LYS:HB3	8:MM:101:TYR:HA	1.88	0.56
5:WH:315:THR:HG22	5:WH:317:LEU:H	1.71	0.56
4:ZG:802:LEU:O	4:ZG:806:THR:OG1	2.22	0.56
3:Af:254:ILE:HB	3:Af:262:ILE:HB	1.87	0.56
2:DD:85:VAL:HG23	6:DK:153:ILE:HG12	1.88	0.56
8:JM:119:GLU:HG3	8:JM:139:ALA:HB3	1.87	0.56
4:BG:807:GLN:OE1	4:BG:810:GLN:NE2	2.37	0.56
5:BH:181:PHE:HD2	5:BH:230:LEU:HD21	1.70	0.56
4:CG:801:MET:HB3	5:DH:115:MET:SD	2.46	0.56
5:GH:179:LEU:O	5:GH:212:THR:HA	2.06	0.56
5:HH:138:GLU:HA	5:HH:141:LYS:HE3	1.87	0.56
5:LH:191:ILE:HA	5:LH:230:LEU:HD13	1.87	0.56
8:AM:32:LYS:HB3	8:AM:101:TYR:HA	1.88	0.55
2:ED:52:MET:HE2	2:Ed:52:MET:HE1	1.88	0.55
8:HM:32:LYS:HB3	8:HM:101:TYR:HA	1.88	0.55
4:IG:802:LEU:O	4:IG:806:THR:OG1	2.22	0.55
8:IM:32:LYS:HB3	8:IM:101:TYR:HA	1.88	0.55
5:MH:169:ILE:HD12	5:MH:252:VAL:HG21	1.86	0.55
2:AD:27:LYS:NZ	6:MK:71:GLY:O	2.37	0.55
1:BC:177:CYS:CA	1:CC:35:GLN:NE2	2.65	0.55
3:Cf:253:ARG:HH21	3:Cf:261:VAL:HG11	1.69	0.55
8:FM:306:GLN:HG3	3:Ff:251:GLN:NE2	2.21	0.55
6:GK:116:VAL:HG22	6:GK:182:ILE:HG12	1.87	0.55
4:JG:807:GLN:OE1	4:JG:810:GLN:NE2	2.37	0.55
5:JH:192:ALA:N	5:JH:229:ARG:O	2.36	0.55
8:JM:32:LYS:HB3	8:JM:101:TYR:HA	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:KG:795:GLN:OE1	4:KG:795:GLN:N	2.35	0.55
8:KM:307:SER:HG	3:Kf:253:ARG:HH22	1.52	0.55
4:LG:802:LEU:O	4:LG:806:THR:OG1	2.22	0.55
2:BD:29:PRO:HG2	7:BL:225:GLN:HA	1.88	0.55
2:Bd:82:ARG:HH21	2:Bd:126:GLU:HA	1.71	0.55
5:EH:238:MET:SD	5:EH:238:MET:N	2.79	0.55
8:EM:32:LYS:HB3	8:EM:101:TYR:HA	1.88	0.55
3:FF:213:GLN:O	5:FH:170:ARG:NH2	2.32	0.55
2:Fd:83:ALA:HA	1:GC:267:ALA:O	2.06	0.55
6:KK:116:VAL:HG22	6:KK:182:ILE:HG12	1.88	0.55
3:AF:229:THR:OG1	5:AH:165:THR:OG1	2.23	0.55
4:GG:807:GLN:OE1	4:GG:810:GLN:NE2	2.37	0.55
8:KM:32:LYS:HB3	8:KM:101:TYR:HA	1.88	0.55
8:KM:308:ASP:OD2	3:Kf:253:ARG:NH2	2.27	0.55
8:MM:119:GLU:HG3	8:MM:139:ALA:HB3	1.87	0.55
5:YH:311:MET:HE1	5:YH:341:LYS:HA	1.88	0.55
2:Ad:87:TRP:CD1	2:Ad:94:LEU:HD12	2.41	0.55
9:GN:63:GLY:HA3	2:HD:27:LYS:HB2	1.89	0.55
5:KH:238:MET:HE1	5:YH:250:TYR:CD2	2.41	0.55
5:XH:205:GLN:HB2	5:XH:214:MET:HB2	1.89	0.55
5:AH:208:LYS:HD2	4:BG:823:GLY:HA2	1.88	0.55
8:BM:32:LYS:HB3	8:BM:101:TYR:HA	1.88	0.55
2:ED:29:PRO:HG2	7:EL:225:GLN:HA	1.87	0.55
2:JD:98:ILE:HG23	2:JD:128:LEU:HD22	1.88	0.55
5:MH:263:ALA:HB1	5:MH:266:MET:HE3	1.89	0.55
8:CM:32:LYS:HB3	8:CM:101:TYR:HA	1.88	0.55
8:DM:119:GLU:HG3	8:DM:139:ALA:HB3	1.87	0.55
5:HH:230:LEU:HD12	5:HH:233:LEU:HD12	1.89	0.55
6:JK:116:VAL:HG22	6:JK:182:ILE:HG12	1.87	0.55
3:XF:229:THR:OG1	5:XH:165:THR:OG1	2.25	0.55
5:BH:205:GLN:HB2	5:BH:214:MET:HB2	1.89	0.55
4:DG:801:MET:HB3	5:FH:115:MET:HE3	1.87	0.55
5:EH:117:ARG:HD3	5:EH:117:ARG:C	2.31	0.55
2:Fd:24:LYS:HG3	2:Fd:25:PHE:HD1	1.71	0.55
1:HC:66:VAL:HG22	1:IC:34:LEU:HD11	1.89	0.55
4:KG:801:MET:HB3	5:LH:115:MET:HE1	1.88	0.55
9:KN:63:GLY:HA3	2:LD:27:LYS:HB2	1.87	0.55
8:GM:32:LYS:HB3	8:GM:101:TYR:HA	1.88	0.55
2:ID:29:PRO:HG2	7:IL:225:GLN:HA	1.88	0.55
2:Id:126:GLU:OE2	2:Id:127:SER:N	2.39	0.55
5:AH:181:PHE:HD2	5:AH:230:LEU:HD21	1.72	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:CH:194:TYR:CE1	5:CH:204:ILE:HD13	2.41	0.55
8:DM:32:LYS:HB3	8:DM:101:TYR:HA	1.88	0.55
5:FH:311:MET:HE2	5:FH:341:LYS:HA	1.89	0.55
8:FM:119:GLU:HG3	8:FM:139:ALA:HB3	1.87	0.55
2:Id:155:GLU:OE2	2:Id:157:ARG:NH2	2.40	0.55
5:KH:155:SER:OG	5:KH:255:ARG:NH1	2.39	0.55
4:AG:802:LEU:HD12	5:CH:115:MET:CE	2.36	0.54
6:BK:116:VAL:HG22	6:BK:182:ILE:HG12	1.87	0.54
6:FK:116:VAL:HG22	6:FK:182:ILE:HG12	1.87	0.54
1:HC:180:THR:HG21	1:IC:34:LEU:CD2	2.37	0.54
4:HG:795:GLN:OE1	4:HG:795:GLN:N	2.35	0.54
7:FL:187:GLU:OE2	8:FM:268:HIS:ND1	2.37	0.54
8:LM:32:LYS:HB3	8:LM:101:TYR:HA	1.88	0.54
1:CC:69:GLN:HE22	9:CN:114:LEU:HA	1.72	0.54
2:Ed:126:GLU:HG3	2:Ed:131:ILE:HG13	1.89	0.54
1:GC:66:VAL:HG22	1:HC:34:LEU:HD11	1.89	0.54
5:IH:202:PHE:HE2	5:IH:241:LEU:HB3	1.71	0.54
2:Id:123:THR:HG21	2:Id:126:GLU:HB2	1.88	0.54
1:JC:157:LYS:NZ	1:KC:28:THR:HA	2.22	0.54
1:JC:247:LEU:HB2	1:KC:125:ILE:HG13	1.89	0.54
4:LG:807:GLN:OE1	4:LG:810:GLN:NE2	2.37	0.54
5:AH:311:MET:HE2	5:AH:341:LYS:HA	1.89	0.54
8:BM:306:GLN:OE1	3:Bf:251:GLN:HG2	2.07	0.54
3:JF:215:VAL:HG12	5:JH:245:GLN:HE22	1.73	0.54
4:BG:795:GLN:OE1	4:BG:795:GLN:N	2.35	0.54
3:DF:229:THR:OG1	5:DH:165:THR:OG1	2.25	0.54
9:IN:63:GLY:HA3	2:JD:27:LYS:HB2	1.88	0.54
4:ZG:807:GLN:OE1	4:ZG:810:GLN:NE2	2.37	0.54
2:DD:55:MET:HE1	2:Dd:55:MET:HE1	1.90	0.54
2:Bd:98:ILE:HG23	2:Bd:128:LEU:HD23	1.89	0.54
5:CH:170:ARG:HH11	5:CH:247:ALA:HB3	1.73	0.54
9:DN:81:LEU:HD21	5:EH:357:LYS:HB2	1.90	0.54
2:ED:23:MET:N	2:ED:23:MET:SD	2.81	0.54
2:Id:146:HIS:HB2	2:Id:155:GLU:HG2	1.90	0.54
4:XG:795:GLN:OE1	4:XG:795:GLN:N	2.35	0.54
4:ZG:795:GLN:OE1	4:ZG:795:GLN:N	2.35	0.54
4:AG:795:GLN:OE1	4:AG:795:GLN:N	2.35	0.54
4:CG:807:GLN:OE1	4:CG:810:GLN:NE2	2.37	0.54
5:DH:158:VAL:HG23	5:DH:254:LEU:HD12	1.89	0.54
5:IH:250:TYR:CG	5:XH:238:MET:HE1	2.43	0.54
2:BD:107:ARG:NH1	2:BD:153:VAL:HG11	2.23	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BC:107:PRO:HG2	1:BC:209:MET:HG2	1.90	0.54
2:Id:113:PRO:HD3	2:Id:158:TYR:HE2	1.71	0.54
7:JL:187:GLU:OE2	8:JM:268:HIS:ND1	2.40	0.54
3:MF:268:ASP:OD1	5:MH:152:THR:HG21	2.08	0.54
9:CN:63:GLY:HA3	2:DD:27:LYS:HB2	1.88	0.53
5:GH:106:VAL:HA	5:GH:109:LYS:HE2	1.89	0.53
2:KD:28:PRO:HG3	7:KL:115:GLN:HB2	1.90	0.53
2:MD:71:LEU:HD12	2:MD:72:THR:HG23	1.90	0.53
1:BC:69:GLN:HE22	9:BN:114:LEU:HA	1.73	0.53
5:EH:251:ARG:NH1	5:EH:253:ASP:OD2	2.42	0.53
1:HC:107:PRO:HG2	1:HC:209:MET:HG2	1.91	0.53
2:Id:82:ARG:NH1	2:Id:125:ASP:OD1	2.33	0.53
5:KH:156:GLN:HB2	5:KH:254:LEU:HD13	1.90	0.53
6:KK:71:GLY:O	2:LD:27:LYS:NZ	2.42	0.53
2:MD:71:LEU:HD13	2:Md:64:PRO:HA	1.90	0.53
5:AH:307:SER:OG	5:AH:308:ASN:N	2.40	0.53
1:CC:126:ARG:HD3	5:CH:275:SER:HA	1.89	0.53
4:GG:818:GLN:HG2	5:GH:216:GLN:NE2	2.23	0.53
1:KC:180:THR:HG21	1:LC:34:LEU:CD1	2.38	0.53
1:LC:157:LYS:HZ3	1:MC:28:THR:HB	1.72	0.53
4:IG:795:GLN:OE1	4:IG:795:GLN:N	2.35	0.53
4:LG:802:LEU:HA	5:ZH:115:MET:HE1	1.90	0.53
5:YH:230:LEU:HD12	5:YH:233:LEU:HD12	1.90	0.53
4:AG:820:TYR:HB3	5:AH:205:GLN:HE22	1.72	0.53
4:GG:795:GLN:OE1	4:GG:795:GLN:N	2.35	0.53
7:GL:57:VAL:HA	7:GL:60:VAL:HG22	1.91	0.53
4:IG:807:GLN:OE1	4:IG:810:GLN:NE2	2.37	0.53
2:KD:92:GLU:OE2	2:KD:158:TYR:OH	2.21	0.53
4:WG:795:GLN:OE1	4:WG:795:GLN:N	2.35	0.53
8:AM:209:SER:OG	8:AM:210:ASP:N	2.42	0.53
8:GM:209:SER:OG	8:GM:210:ASP:N	2.42	0.53
8:HM:209:SER:OG	8:HM:210:ASP:N	2.42	0.53
1:JC:107:PRO:HG2	1:JC:209:MET:HG2	1.91	0.53
8:KM:209:SER:OG	8:KM:210:ASP:N	2.42	0.53
7:LL:57:VAL:HA	7:LL:60:VAL:HG22	1.91	0.53
1:DC:119:LEU:HA	1:DC:125:ILE:HG22	1.91	0.53
4:EG:807:GLN:OE1	4:EG:810:GLN:NE2	2.37	0.53
3:GF:213:GLN:HA	5:GH:170:ARG:HH21	1.73	0.53
1:IC:107:PRO:HG2	1:IC:209:MET:HG2	1.90	0.53
1:KC:107:PRO:HG2	1:KC:209:MET:HG2	1.90	0.53
1:KC:119:LEU:HA	1:KC:125:ILE:HG22	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:KL:57:VAL:HA	7:KL:60:VAL:HG22	1.91	0.53
9:LN:63:GLY:HA3	2:MD:27:LYS:HB2	1.88	0.53
5:WH:190:PRO:HB2	5:WH:231:ARG:HG3	1.89	0.53
5:XH:183:ASP:HA	5:XH:256:VAL:HG12	1.91	0.53
1:AC:107:PRO:HG2	1:AC:209:MET:HG2	1.90	0.53
5:CH:266:MET:HE3	5:CH:267:PRO:HD2	1.90	0.53
7:CL:57:VAL:HA	7:CL:60:VAL:HG22	1.91	0.53
1:DC:107:PRO:HG2	1:DC:209:MET:HG2	1.90	0.53
8:DM:209:SER:OG	8:DM:210:ASP:N	2.42	0.53
2:JD:77:TYR:OH	6:JK:130:ARG:NH1	2.40	0.53
2:Bd:126:GLU:OE2	2:Bd:130:GLU:HB3	2.08	0.53
8:CM:209:SER:OG	8:CM:210:ASP:N	2.42	0.53
1:FC:119:LEU:HA	1:FC:125:ILE:HG22	1.91	0.53
5:YH:280:LEU:HA	5:YH:283:VAL:HG12	1.91	0.53
1:BC:119:LEU:HA	1:BC:125:ILE:HG22	1.91	0.52
5:IH:200:SER:N	4:JG:816:GLU:OE2	2.39	0.52
8:MM:209:SER:OG	8:MM:210:ASP:N	2.42	0.52
1:GC:107:PRO:HG2	1:GC:209:MET:HG2	1.91	0.52
2:JD:85:VAL:HG12	6:JK:153:ILE:HG23	1.92	0.52
3:Jf:247:ILE:HG22	3:Jf:254:ILE:HD12	1.91	0.52
3:LF:245:LYS:HE3	3:LF:257:SER:HA	1.92	0.52
4:MG:795:GLN:OE1	4:MG:795:GLN:N	2.35	0.52
4:XG:807:GLN:OE1	4:XG:810:GLN:NE2	2.37	0.52
5:XH:159:ASN:OD1	5:XH:257:GLN:NE2	2.41	0.52
4:YG:807:GLN:OE1	4:YG:810:GLN:NE2	2.37	0.52
1:AC:119:LEU:HA	1:AC:125:ILE:HG22	1.91	0.52
1:AC:240:ILE:HG12	1:BC:255:LEU:HD13	1.92	0.52
1:BC:184:TRP:CD1	1:CC:31:LEU:HD22	2.44	0.52
8:DM:307:SER:HG	3:Df:253:ARG:HH22	1.58	0.52
1:FC:107:PRO:HG2	1:FC:209:MET:HG2	1.90	0.52
3:FF:245:LYS:HE3	3:FF:257:SER:HA	1.91	0.52
8:JM:209:SER:OG	8:JM:210:ASP:N	2.42	0.52
3:KF:245:LYS:HE3	3:KF:257:SER:HA	1.92	0.52
5:VH:279:LEU:HD11	5:VH:290:PRO:HD2	1.89	0.52
5:XH:319:ILE:HA	5:XH:347:VAL:HG22	1.91	0.52
3:DF:245:LYS:HE3	3:DF:257:SER:HA	1.91	0.52
4:FG:807:GLN:OE1	4:FG:810:GLN:NE2	2.37	0.52
3:HF:245:LYS:HE3	3:HF:257:SER:HA	1.92	0.52
4:LG:795:GLN:OE1	4:LG:795:GLN:N	2.35	0.52
5:CH:223:TYR:HB2	5:CH:242:ILE:HG13	1.91	0.52
1:EC:107:PRO:HG2	1:EC:209:MET:HG2	1.90	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:HC:119:LEU:HA	1:HC:125:ILE:HG22	1.91	0.52
1:IC:119:LEU:HA	1:IC:125:ILE:HG22	1.91	0.52
1:JC:119:LEU:HA	1:JC:125:ILE:HG22	1.91	0.52
2:Kd:64:PRO:HG2	2:Kd:67:LYS:HG2	1.91	0.52
5:MH:311:MET:HE3	5:MH:341:LYS:HA	1.92	0.52
3:WF:245:LYS:HE3	3:WF:257:SER:HA	1.92	0.52
3:AF:245:LYS:HE3	3:AF:257:SER:HA	1.92	0.52
2:CD:82:ARG:NE	2:CD:125:ASP:OD1	2.34	0.52
1:DC:30:SER:OG	1:DC:31:LEU:N	2.43	0.52
6:FK:71:GLY:O	2:GD:27:LYS:NZ	2.42	0.52
4:GG:818:GLN:HG2	5:GH:216:GLN:HE22	1.74	0.52
1:MC:107:PRO:HG2	1:MC:209:MET:HG2	1.91	0.52
3:VF:245:LYS:HE3	3:VF:257:SER:HA	1.92	0.52
1:CC:107:PRO:HG2	1:CC:209:MET:HG2	1.90	0.52
1:CC:176:TRP:HZ3	1:DC:35:GLN:OE1	1.90	0.52
9:JN:63:GLY:HA3	2:KD:27:LYS:HB2	1.90	0.52
2:AD:123:THR:OG1	2:AD:124:LYS:N	2.42	0.52
5:CH:173:GLN:NE2	5:CH:218:THR:O	2.38	0.52
1:GC:119:LEU:HA	1:GC:125:ILE:HG22	1.91	0.52
4:GG:801:MET:C	5:XH:115:MET:HE1	2.35	0.52
8:IM:209:SER:OG	8:IM:210:ASP:N	2.42	0.52
3:JF:245:LYS:HE3	3:JF:257:SER:HA	1.92	0.52
4:JG:801:MET:HB3	5:YH:115:MET:SD	2.50	0.52
5:KH:307:SER:OG	5:KH:308:ASN:N	2.42	0.52
8:CM:212:LEU:HD12	8:CM:214:ILE:HD11	1.92	0.52
7:DL:57:VAL:HA	7:DL:60:VAL:HG22	1.91	0.52
6:EK:71:GLY:O	2:FD:27:LYS:NZ	2.37	0.52
2:LD:28:PRO:HG3	7:LL:115:GLN:HB2	1.91	0.52
6:CK:42:PRO:HD2	9:CN:76:MET:HE3	1.91	0.52
1:GC:225:TYR:HB2	1:GC:251:ALA:HB3	1.92	0.52
4:GG:816:GLU:OE2	5:WH:200:SER:N	2.37	0.52
7:IL:57:VAL:HA	7:IL:60:VAL:HG22	1.91	0.52
7:JL:57:VAL:HA	7:JL:60:VAL:HG22	1.91	0.52
1:LC:107:PRO:HG2	1:LC:209:MET:HG2	1.90	0.52
7:ML:57:VAL:HA	7:ML:60:VAL:HG22	1.91	0.52
4:AG:807:GLN:OE1	4:AG:810:GLN:NE2	2.37	0.51
5:AH:251:ARG:HH22	5:ZH:235:THR:HG23	1.75	0.51
2:BD:55:MET:HE1	2:Bd:55:MET:HE1	1.90	0.51
8:CM:292:MET:HE1	8:CM:298:VAL:H	1.76	0.51
1:DC:225:TYR:HB2	1:DC:251:ALA:HB3	1.93	0.51
5:FH:179:LEU:O	5:FH:212:THR:HA	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:FL:57:VAL:HA	7:FL:60:VAL:HG22	1.91	0.51
5:HH:223:TYR:HB2	5:HH:242:ILE:HG13	1.92	0.51
8:JM:292:MET:HE1	8:JM:298:VAL:H	1.75	0.51
3:MF:245:LYS:HE3	3:MF:257:SER:HA	1.92	0.51
1:CC:119:LEU:HA	1:CC:125:ILE:HG22	1.91	0.51
8:FM:212:LEU:HD12	8:FM:214:ILE:HD11	1.93	0.51
9:FN:56:GLY:HA2	9:FN:71:THR:HG23	1.93	0.51
5:IH:322:PRO:HG2	5:IH:345:LEU:HD23	1.93	0.51
5:JH:192:ALA:HA	5:JH:231:ARG:NH2	2.25	0.51
8:KM:292:MET:HE1	8:KM:298:VAL:H	1.75	0.51
5:LH:238:MET:HE1	5:MH:250:TYR:HD2	1.74	0.51
1:MC:119:LEU:HA	1:MC:125:ILE:HG22	1.91	0.51
8:MM:212:LEU:HD12	8:MM:214:ILE:HD11	1.92	0.51
5:XH:349:TRP:CD1	5:XH:354:MET:HE1	2.44	0.51
7:AL:57:VAL:HA	7:AL:60:VAL:HG22	1.91	0.51
1:BC:177:CYS:HA	1:CC:35:GLN:HE21	1.73	0.51
3:BF:245:LYS:HE3	3:BF:257:SER:HA	1.92	0.51
3:Bf:221:TRP:HE1	3:Bf:229:THR:HB	1.74	0.51
5:EH:230:LEU:HD12	5:EH:233:LEU:HD12	1.92	0.51
2:HD:59:GLU:HA	2:HD:62:ILE:HG22	1.92	0.51
4:HG:807:GLN:OE1	4:HG:810:GLN:NE2	2.37	0.51
8:JM:212:LEU:HD12	8:JM:214:ILE:HD11	1.92	0.51
9:JN:56:GLY:HA2	9:JN:71:THR:HG23	1.92	0.51
1:LC:119:LEU:HA	1:LC:125:ILE:HG22	1.91	0.51
2:MD:82:ARG:NE	2:MD:125:ASP:OD1	2.37	0.51
4:AG:823:GLY:HA2	5:ZH:208:LYS:HE3	1.93	0.51
5:AH:106:VAL:HA	5:AH:109:LYS:HE2	1.91	0.51
8:BM:292:MET:HE1	8:BM:298:VAL:H	1.75	0.51
5:CH:200:SER:N	4:VG:816:GLU:OE2	2.41	0.51
3:EF:245:LYS:HE3	3:EF:257:SER:HA	1.92	0.51
7:EL:57:VAL:HA	7:EL:60:VAL:HG22	1.91	0.51
8:EM:209:SER:OG	8:EM:210:ASP:N	2.42	0.51
8:EM:212:LEU:HD12	8:EM:214:ILE:HD11	1.93	0.51
2:GD:82:ARG:NE	2:GD:125:ASP:OD1	2.35	0.51
8:GM:212:LEU:HD12	8:GM:214:ILE:HD11	1.92	0.51
8:GM:247:ARG:HD3	3:Gf:213:GLN:O	2.10	0.51
1:MC:225:TYR:HB2	1:MC:251:ALA:HB3	1.93	0.51
2:Md:98:ILE:HG23	2:Md:128:LEU:HD23	1.91	0.51
4:VG:795:GLN:OE1	4:VG:795:GLN:N	2.35	0.51
1:CC:184:TRP:CD1	1:DC:31:LEU:HD21	2.45	0.51
4:FG:795:GLN:OE1	4:FG:795:GLN:N	2.35	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:FM:209:SER:OG	8:FM:210:ASP:N	2.42	0.51
8:FM:292:MET:HE1	8:FM:298:VAL:H	1.76	0.51
1:HC:225:TYR:HB2	1:HC:251:ALA:HB3	1.93	0.51
3:IF:245:LYS:HE3	3:IF:257:SER:HA	1.92	0.51
1:JC:225:TYR:HB2	1:JC:251:ALA:HB3	1.93	0.51
5:WH:110:LYS:HA	5:WH:113:LYS:HZ3	1.75	0.51
3:XF:245:LYS:HE3	3:XF:257:SER:HA	1.92	0.51
1:AC:28:THR:OG1	1:MC:157:LYS:NZ	2.40	0.51
1:CC:225:TYR:HB2	1:CC:251:ALA:HB3	1.93	0.51
2:DD:123:THR:OG1	2:DD:124:LYS:N	2.43	0.51
8:DM:292:MET:HE1	8:DM:298:VAL:H	1.76	0.51
7:HL:57:VAL:HA	7:HL:60:VAL:HG22	1.91	0.51
2:KD:77:TYR:OH	6:KK:130:ARG:NH1	2.43	0.51
4:KG:807:GLN:OE1	4:KG:810:GLN:NE2	2.37	0.51
8:KM:306:GLN:OE1	3:Kf:251:GLN:HG2	2.09	0.51
9:LN:56:GLY:HA2	9:LN:71:THR:HG23	1.93	0.51
3:ZF:245:LYS:HE3	3:ZF:257:SER:HA	1.92	0.51
5:ZH:293:ARG:NH2	5:ZH:360:GLY:O	2.40	0.51
5:CH:238:MET:HE1	5:VH:250:TYR:HB3	1.92	0.51
9:DN:56:GLY:HA2	9:DN:71:THR:HG23	1.93	0.51
5:HH:143:ALA:O	5:XH:131:LYS:NZ	2.33	0.51
4:JG:795:GLN:OE1	4:JG:795:GLN:N	2.35	0.51
1:KC:116:THR:HG23	5:KH:328:MET:HG2	1.93	0.51
5:KH:111:ALA:O	5:KH:115:MET:CB	2.59	0.51
5:LH:159:ASN:OD1	5:LH:257:GLN:NE2	2.39	0.51
1:AC:225:TYR:HB2	1:AC:251:ALA:HB3	1.93	0.51
5:BH:190:PRO:HB2	5:BH:231:ARG:HD3	1.92	0.51
7:BL:57:VAL:HA	7:BL:60:VAL:HG22	1.91	0.51
3:CF:245:LYS:HE3	3:CF:257:SER:HA	1.91	0.51
8:EM:292:MET:HE1	8:EM:298:VAL:H	1.76	0.51
8:GM:292:MET:HE1	8:GM:298:VAL:H	1.75	0.51
8:HM:212:LEU:HD12	8:HM:214:ILE:HD11	1.93	0.51
2:LD:123:THR:HG22	2:LD:124:LYS:H	1.75	0.51
8:LM:292:MET:HE1	8:LM:298:VAL:H	1.76	0.51
9:MN:56:GLY:HA2	9:MN:71:THR:HG23	1.93	0.51
8:BM:209:SER:OG	8:BM:210:ASP:N	2.42	0.51
1:DC:157:LYS:NZ	1:EC:28:THR:CB	2.72	0.51
4:DG:807:GLN:OE1	4:DG:810:GLN:NE2	2.37	0.51
1:EC:119:LEU:HA	1:EC:125:ILE:HG22	1.91	0.51
2:Ed:134:ASP:OD1	2:Ed:138:GLN:NE2	2.42	0.51
1:FC:225:TYR:HB2	1:FC:251:ALA:HB3	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:GH:171:LEU:HD11	5:GH:241:LEU:HD12	1.93	0.51
5:GH:322:PRO:HG2	5:GH:345:LEU:HD23	1.92	0.51
2:HD:72:THR:HG23	2:HD:73:ILE:HG12	1.92	0.51
2:Hd:82:ARG:NH1	2:Hd:125:ASP:OD1	2.44	0.51
2:KD:82:ARG:NE	2:KD:125:ASP:OD1	2.30	0.51
3:YF:245:LYS:HE3	3:YF:257:SER:HA	1.91	0.51
1:BC:184:TRP:CG	1:CC:31:LEU:HD22	2.46	0.51
9:HN:63:GLY:HA3	2:ID:27:LYS:HB2	1.93	0.51
9:HN:89:GLY:HA3	9:HN:104:ASN:HB2	1.93	0.51
5:IH:194:TYR:HE2	5:IH:196:LEU:HD22	1.76	0.51
5:IH:208:LYS:HD3	4:JG:823:GLY:HA2	1.93	0.51
9:IN:56:GLY:HA2	9:IN:71:THR:HG23	1.93	0.51
2:JD:52:MET:HE2	2:Jd:52:MET:HE1	1.93	0.51
1:LC:180:THR:HG21	1:MC:34:LEU:CD2	2.41	0.51
1:AC:157:LYS:HZ1	1:BC:28:THR:HB	1.76	0.50
5:BH:115:MET:CE	4:ZG:802:LEU:HD12	2.41	0.50
6:BK:71:GLY:O	2:CD:27:LYS:NZ	2.37	0.50
8:JM:247:ARG:HD2	3:Jf:264:PHE:CE1	2.46	0.50
9:KN:56:GLY:HA2	9:KN:71:THR:HG23	1.93	0.50
8:MM:292:MET:HE1	8:MM:298:VAL:H	1.76	0.50
9:AN:63:GLY:HA3	2:BD:27:LYS:HB2	1.93	0.50
2:Fd:24:LYS:HE2	2:Fd:25:PHE:CE1	2.45	0.50
5:GH:151:PRO:HA	5:GH:250:TYR:CE2	2.46	0.50
9:GN:56:GLY:HA2	9:GN:71:THR:HG23	1.93	0.50
8:IM:212:LEU:HD12	8:IM:214:ILE:HD11	1.93	0.50
8:KM:212:LEU:HD12	8:KM:214:ILE:HD11	1.93	0.50
9:KN:89:GLY:HA3	9:KN:104:ASN:HB2	1.94	0.50
5:LH:190:PRO:HB2	5:LH:231:ARG:HG3	1.91	0.50
8:AM:292:MET:HE1	8:AM:298:VAL:H	1.76	0.50
9:DN:89:GLY:HA3	9:DN:104:ASN:HB2	1.94	0.50
1:EC:225:TYR:HB2	1:EC:251:ALA:HB3	1.93	0.50
9:EN:89:GLY:HA3	9:EN:104:ASN:HB2	1.94	0.50
3:GF:219:ARG:HH12	5:HH:150:LYS:HG2	1.76	0.50
5:GH:230:LEU:HD12	5:GH:233:LEU:HD12	1.92	0.50
7:IL:187:GLU:OE2	8:IM:268:HIS:ND1	2.43	0.50
9:IN:89:GLY:HA3	9:IN:104:ASN:HB2	1.94	0.50
2:Jd:98:ILE:HD12	2:Jd:128:LEU:HD22	1.92	0.50
5:MH:311:MET:HE3	5:MH:342:SER:H	1.75	0.50
9:AN:89:GLY:HA3	9:AN:104:ASN:HB2	1.94	0.50
5:BH:229:ARG:HG2	5:BH:236:PRO:HB3	1.94	0.50
9:BN:56:GLY:HA2	9:BN:71:THR:HG23	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:DD:29:PRO:HG2	7:DL:225:GLN:HA	1.94	0.50
2:JD:59:GLU:HA	2:JD:62:ILE:HG22	1.93	0.50
1:KC:225:TYR:HB2	1:KC:251:ALA:HB3	1.92	0.50
8:LM:31:LYS:O	8:LM:76:ARG:NH2	2.43	0.50
8:LM:209:SER:OG	8:LM:210:ASP:N	2.42	0.50
1:AC:213:ARG:NH2	2:Ad:59:GLU:OE2	2.44	0.50
1:BC:225:TYR:HB2	1:BC:251:ALA:HB3	1.93	0.50
2:CD:23:MET:N	2:CD:23:MET:SD	2.84	0.50
8:FM:31:LYS:O	8:FM:76:ARG:NH2	2.43	0.50
3:GF:245:LYS:HE3	3:GF:257:SER:HA	1.92	0.50
2:Gd:134:ASP:OD1	2:Gd:138:GLN:NE2	2.42	0.50
6:HK:41:LEU:HD22	9:HN:76:MET:HE1	1.94	0.50
2:LD:24:LYS:O	7:LL:110:GLN:NE2	2.45	0.50
5:YH:326:ALA:HB3	5:YH:338:GLU:HG2	1.93	0.50
9:AN:56:GLY:HA2	9:AN:71:THR:HG23	1.93	0.50
2:CD:92:GLU:OE1	2:CD:158:TYR:OH	2.29	0.50
8:DM:212:LEU:HD12	8:DM:214:ILE:HD11	1.93	0.50
1:JC:157:LYS:HZ3	1:KC:28:THR:HA	1.76	0.50
2:LD:85:VAL:HG12	6:LK:153:ILE:HG23	1.93	0.50
6:BK:42:PRO:HD2	9:BN:76:MET:HE3	1.93	0.50
2:Bd:107:ARG:HG3	2:Bd:155:GLU:HG3	1.93	0.50
8:HM:31:LYS:O	8:HM:76:ARG:NH2	2.43	0.50
1:IC:225:TYR:HB2	1:IC:251:ALA:HB3	1.93	0.50
1:MC:35:GLN:NE2	1:MC:35:GLN:HA	2.27	0.50
5:WH:171:LEU:HD11	5:WH:241:LEU:HB3	1.92	0.50
5:ZH:317:LEU:HD11	5:ZH:349:TRP:HB2	1.92	0.50
8:AM:212:LEU:HD12	8:AM:214:ILE:HD11	1.93	0.50
3:Af:238:ILE:HD12	3:Af:239:PRO:HD2	1.93	0.50
6:BK:41:LEU:HD22	9:BN:76:MET:HE1	1.93	0.50
6:EK:41:LEU:HD22	9:EN:76:MET:HE1	1.94	0.50
9:EN:56:GLY:HA2	9:EN:71:THR:HG23	1.93	0.50
9:GN:89:GLY:HA3	9:GN:104:ASN:HB2	1.94	0.50
5:HH:203:ASN:HB3	5:HH:216:GLN:HG2	1.94	0.50
8:HM:292:MET:HE1	8:HM:298:VAL:H	1.76	0.50
5:BH:173:GLN:NE2	5:BH:218:THR:O	2.40	0.50
5:BH:191:ILE:O	5:BH:231:ARG:NH1	2.45	0.50
8:BM:212:LEU:HD12	8:BM:214:ILE:HD11	1.93	0.50
2:CD:59:GLU:HA	2:CD:62:ILE:HG22	1.94	0.50
9:CN:56:GLY:HA2	9:CN:71:THR:HG23	1.93	0.50
8:KM:205:TYR:CZ	3:Kf:219:ARG:NH2	2.79	0.50
1:AC:34:LEU:HD11	1:MC:66:VAL:HG22	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:BH:107:ILE:HD12	4:ZG:794:ILE:HD13	1.94	0.49
2:DD:82:ARG:NE	2:DD:125:ASP:OD1	2.38	0.49
2:Fd:82:ARG:HH22	2:Fd:126:GLU:C	2.19	0.49
8:IM:292:MET:HE1	8:IM:298:VAL:H	1.76	0.49
5:BH:156:GLN:NE2	5:BH:157:PHE:O	2.45	0.49
5:BH:322:PRO:HG2	5:BH:345:LEU:HD23	1.95	0.49
3:GF:219:ARG:HH22	5:HH:150:LYS:HD3	1.77	0.49
5:IH:150:LYS:NZ	3:XF:231:THR:O	2.33	0.49
1:JC:240:ILE:CG1	1:KC:132:TYR:HB2	2.42	0.49
1:LC:225:TYR:HB2	1:LC:251:ALA:HB3	1.93	0.49
9:LN:89:GLY:HA3	9:LN:104:ASN:HB2	1.94	0.49
3:Lf:238:ILE:HD12	3:Lf:239:PRO:HD2	1.93	0.49
9:BN:89:GLY:HA3	9:BN:104:ASN:HB2	1.94	0.49
9:CN:89:GLY:HA3	9:CN:104:ASN:HB2	1.94	0.49
2:DD:77:TYR:OH	6:DK:130:ARG:NH1	2.44	0.49
8:EM:31:LYS:O	8:EM:76:ARG:NH2	2.43	0.49
5:FH:106:VAL:HA	5:FH:109:LYS:HE2	1.94	0.49
8:FM:306:GLN:HG3	3:Ff:251:GLN:CD	2.38	0.49
2:Fd:82:ARG:HH21	2:Fd:83:ALA:N	2.07	0.49
5:GH:208:LYS:NZ	4:HG:822:GLU:O	2.39	0.49
6:GK:166:THR:HG23	6:GK:168:TYR:H	1.78	0.49
5:IH:311:MET:HE3	5:IH:342:SER:H	1.77	0.49
8:JM:274:THR:HA	8:JM:293:ALA:HB3	1.95	0.49
5:ZH:111:ALA:O	5:ZH:115:MET:CB	2.57	0.49
8:BM:31:LYS:O	8:BM:76:ARG:NH2	2.43	0.49
8:EM:274:THR:HA	8:EM:293:ALA:HB3	1.95	0.49
6:FK:166:THR:HG23	6:FK:168:TYR:H	1.78	0.49
6:JK:166:THR:HG23	6:JK:168:TYR:H	1.78	0.49
6:MK:166:THR:HG23	6:MK:168:TYR:H	1.78	0.49
5:WH:348:SER:HA	5:WH:352:LYS:O	2.12	0.49
2:BD:92:GLU:OE2	2:BD:158:TYR:OH	2.30	0.49
7:CL:145:TYR:HA	7:CL:148:LEU:HD12	1.95	0.49
5:DH:183:ASP:HA	5:DH:256:VAL:HG12	1.93	0.49
4:EG:795:GLN:OE1	4:EG:795:GLN:N	2.35	0.49
9:FN:89:GLY:HA3	9:FN:104:ASN:HB2	1.94	0.49
9:HN:56:GLY:HA2	9:HN:71:THR:HG23	1.93	0.49
8:LM:212:LEU:HD12	8:LM:214:ILE:HD11	1.93	0.49
5:AH:110:LYS:HA	5:AH:113:LYS:HE2	1.94	0.49
6:CK:41:LEU:HD22	9:CN:76:MET:HE1	1.94	0.49
7:DL:145:TYR:HA	7:DL:148:LEU:HD12	1.95	0.49
5:IH:225:ASN:C	5:IH:225:ASN:HD22	2.20	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:BM:274:THR:HA	8:BM:293:ALA:HB3	1.95	0.49
5:DH:322:PRO:HG2	5:DH:345:LEU:HD23	1.94	0.49
8:GM:274:THR:HA	8:GM:293:ALA:HB3	1.95	0.49
4:HG:820:TYR:OH	5:HH:207:ASP:OD2	2.11	0.49
6:IK:166:THR:HG23	6:IK:168:TYR:H	1.78	0.49
4:JG:800:ASP:N	4:JG:800:ASP:OD1	2.46	0.49
3:Lf:251:GLN:OE1	3:Lf:253:ARG:NH1	2.46	0.49
2:MD:28:PRO:HG3	7:ML:115:GLN:HB2	1.94	0.49
8:MM:274:THR:HA	8:MM:293:ALA:HB3	1.95	0.49
4:WG:807:GLN:OE1	4:WG:810:GLN:NE2	2.37	0.49
4:CG:800:ASP:OD1	4:CG:800:ASP:N	2.46	0.49
5:CH:322:PRO:HG2	5:CH:345:LEU:HD23	1.95	0.49
6:CK:166:THR:HG23	6:CK:168:TYR:H	1.78	0.49
8:CM:274:THR:HA	8:CM:293:ALA:HB3	1.95	0.49
5:JH:189:TRP:HE1	5:JH:232:GLY:H	1.61	0.49
2:AD:52:MET:HE2	2:AD:52:MET:HB2	1.66	0.49
4:BG:800:ASP:OD1	4:BG:800:ASP:N	2.46	0.49
6:EK:166:THR:HG23	6:EK:168:TYR:H	1.78	0.49
9:JN:89:GLY:HA3	9:JN:104:ASN:HB2	1.94	0.49
5:LH:143:ALA:O	5:MH:131:LYS:NZ	2.42	0.49
5:LH:157:PHE:HD1	5:LH:255:ARG:HB2	1.76	0.49
9:MN:89:GLY:HA3	9:MN:104:ASN:HB2	1.94	0.49
8:FM:274:THR:HA	8:FM:293:ALA:HB3	1.95	0.48
8:HM:274:THR:HA	8:HM:293:ALA:HB3	1.95	0.48
3:Kf:254:ILE:HB	3:Kf:262:ILE:HG13	1.95	0.48
5:LH:183:ASP:HB3	5:LH:189:TRP:CE3	2.48	0.48
4:AG:820:TYR:HB2	5:AH:205:GLN:OE1	2.13	0.48
6:DK:166:THR:HG23	6:DK:168:TYR:H	1.78	0.48
4:EG:801:MET:HB3	5:WH:115:MET:SD	2.53	0.48
5:HH:115:MET:HE1	4:WG:798:THR:HA	1.95	0.48
6:LK:166:THR:HG23	6:LK:168:TYR:H	1.78	0.48
4:YG:795:GLN:OE1	4:YG:795:GLN:N	2.35	0.48
8:DM:274:THR:HA	8:DM:293:ALA:HB3	1.95	0.48
1:EC:104:ASN:OD1	2:ED:142:LYS:NZ	2.47	0.48
4:GG:801:MET:HB3	5:XH:115:MET:SD	2.53	0.48
5:VH:157:PHE:HA	5:VH:255:ARG:HB2	1.94	0.48
6:BK:166:THR:HG23	6:BK:168:TYR:H	1.78	0.48
3:Df:243:MET:HB3	3:Df:257:SER:HB3	1.96	0.48
5:EH:294:ARG:NH2	9:EN:110:GLU:OE2	2.43	0.48
7:HL:145:TYR:HA	7:HL:148:LEU:HD12	1.95	0.48
7:LL:145:TYR:HA	7:LL:148:LEU:HD12	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:VG:800:ASP:OD1	4:VG:800:ASP:N	2.46	0.48
2:Dd:97:ARG:HA	2:Dd:97:ARG:HD3	1.61	0.48
5:GH:184:SER:O	5:GH:255:ARG:NH2	2.46	0.48
6:IK:100:ASN:HD22	6:IK:100:ASN:N	2.11	0.48
5:JH:268:THR:OG1	5:JH:269:GLU:N	2.43	0.48
2:DD:87:TRP:CG	2:DD:94:LEU:HD12	2.48	0.48
5:EH:284:LEU:HD22	5:EH:330:SER:HB3	1.94	0.48
2:Ed:123:THR:HG21	2:Ed:126:GLU:HB2	1.95	0.48
7:GL:145:TYR:HA	7:GL:148:LEU:HD12	1.95	0.48
6:HK:166:THR:HG23	6:HK:168:TYR:H	1.78	0.48
2:Id:126:GLU:HG3	2:Id:131:ILE:HG13	1.95	0.48
1:KC:180:THR:HG21	1:LC:34:LEU:HD11	1.95	0.48
1:KC:240:ILE:HG12	1:LC:255:LEU:HD13	1.95	0.48
3:LF:235:GLY:HA2	3:LF:243:MET:HE1	1.96	0.48
5:XH:297:VAL:HG23	5:XH:358:VAL:HG22	1.94	0.48
5:YH:311:MET:HE3	5:YH:311:MET:HB3	1.66	0.48
8:AM:274:THR:HA	8:AM:293:ALA:HB3	1.95	0.48
1:CC:210:ILE:HD12	1:CC:210:ILE:HA	1.65	0.48
4:CG:818:GLN:HG3	5:CH:205:GLN:HE22	1.78	0.48
4:GG:800:ASP:OD1	4:GG:800:ASP:N	2.46	0.48
5:GH:238:MET:HE1	5:HH:250:TYR:HB3	1.96	0.48
5:IH:250:TYR:HD2	5:XH:166:PRO:HG3	1.79	0.48
5:KH:322:PRO:HG2	5:KH:345:LEU:HD23	1.96	0.48
5:MH:221:TYR:OH	5:ZH:135:GLU:OE2	2.32	0.48
2:AD:28:PRO:HG3	7:AL:115:GLN:HB2	1.95	0.48
2:CD:52:MET:HE2	2:Cd:52:MET:HE1	1.96	0.48
3:CF:269:SER:C	5:CH:247:ALA:HB2	2.38	0.48
5:EH:190:PRO:HG2	5:EH:231:ARG:HH12	1.79	0.48
3:FF:235:GLY:HA2	3:FF:243:MET:HE1	1.96	0.48
4:FG:801:MET:HG3	5:GH:115:MET:HE1	1.96	0.48
5:GH:284:LEU:HD22	5:GH:330:SER:HB3	1.96	0.48
2:ID:67:LYS:HE3	2:Id:64:PRO:HD3	1.95	0.48
5:IH:205:GLN:HB2	5:IH:214:MET:HB2	1.96	0.48
8:IM:31:LYS:O	8:IM:76:ARG:NH2	2.43	0.48
6:LK:71:GLY:O	2:MD:27:LYS:NZ	2.45	0.48
7:LL:47:TYR:OH	7:LL:142:GLU:OE1	2.28	0.48
8:BM:125:THR:HG23	8:BM:145:ARG:HB2	1.96	0.48
2:CD:98:ILE:HG23	2:CD:128:LEU:HD22	1.95	0.48
3:DF:235:GLY:HA2	3:DF:243:MET:HE1	1.96	0.48
5:DH:225:ASN:C	5:DH:225:ASN:ND2	2.72	0.48
6:DK:100:ASN:N	6:DK:100:ASN:HD22	2.11	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:EL:145:TYR:HA	7:EL:148:LEU:HD12	1.95	0.48
2:Ed:126:GLU:OE2	2:Ed:127:SER:N	2.46	0.48
5:FH:129:LYS:HA	5:FH:132:GLN:HG2	1.95	0.48
1:IC:177:CYS:HB3	1:JC:35:GLN:HE22	1.77	0.48
3:JF:235:GLY:HA2	3:JF:243:MET:HE1	1.96	0.48
5:JH:322:PRO:HG2	5:JH:345:LEU:HD23	1.96	0.48
4:KG:794:ILE:HD13	5:LH:107:ILE:HG23	1.94	0.48
1:MC:75:ARG:NH2	2:MD:36:ASP:OD1	2.47	0.48
4:MG:807:GLN:OE1	4:MG:810:GLN:NE2	2.37	0.48
3:VF:235:GLY:HA2	3:VF:243:MET:HE1	1.96	0.48
2:Cd:64:PRO:HG2	2:Cd:67:LYS:HG2	1.96	0.48
6:EK:117:THR:HG22	6:EK:157:GLN:HG3	1.96	0.48
8:EM:125:THR:HG23	8:EM:145:ARG:HB2	1.96	0.48
8:HM:125:THR:HG23	8:HM:145:ARG:HB2	1.96	0.48
7:ML:145:TYR:HA	7:ML:148:LEU:HD12	1.95	0.48
6:AK:166:THR:HG23	6:AK:168:TYR:H	1.78	0.47
7:BL:44:HIS:HB2	7:BL:48:ASN:HA	1.96	0.47
7:BL:145:TYR:HA	7:BL:148:LEU:HD12	1.95	0.47
5:FH:157:PHE:HA	5:FH:255:ARG:HB2	1.96	0.47
2:HD:94:LEU:HD13	2:HD:119:ILE:HD11	1.96	0.47
3:Hf:247:ILE:HG22	3:Hf:254:ILE:HD12	1.95	0.47
8:IM:274:THR:HA	8:IM:293:ALA:HB3	1.95	0.47
7:KL:145:TYR:HA	7:KL:148:LEU:HD12	1.95	0.47
8:KM:308:ASP:CG	3:Kf:253:ARG:HH21	2.19	0.47
8:LM:274:THR:HA	8:LM:293:ALA:HB3	1.95	0.47
5:ZH:296:VAL:HB	5:ZH:359:GLU:HG2	1.96	0.47
7:AL:44:HIS:HB2	7:AL:48:ASN:HA	1.96	0.47
3:Bf:243:MET:HB3	3:Bf:257:SER:HB3	1.96	0.47
3:KF:235:GLY:HA2	3:KF:243:MET:HE1	1.96	0.47
5:LH:284:LEU:HD22	5:LH:330:SER:HB3	1.95	0.47
6:MK:117:THR:HG22	6:MK:157:GLN:HG3	1.96	0.47
5:WH:292:SER:HB2	5:WH:305:TRP:HB3	1.96	0.47
5:XH:317:LEU:HB3	5:XH:347:VAL:HG11	1.95	0.47
7:AL:145:TYR:HA	7:AL:148:LEU:HD12	1.95	0.47
6:BK:117:THR:HG22	6:BK:157:GLN:HG3	1.96	0.47
2:DD:92:GLU:OE2	2:DD:158:TYR:OH	2.33	0.47
8:DM:110:THR:OG1	8:DM:111:LYS:N	2.47	0.47
7:FL:145:TYR:HA	7:FL:148:LEU:HD12	1.95	0.47
8:GM:31:LYS:O	8:GM:76:ARG:NH2	2.43	0.47
5:HH:322:PRO:HG2	5:HH:345:LEU:HD23	1.96	0.47
5:IH:143:ALA:O	5:JH:131:LYS:NZ	2.45	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:JK:117:THR:HG22	6:JK:157:GLN:HG3	1.97	0.47
7:JL:145:TYR:HA	7:JL:148:LEU:HD12	1.95	0.47
6:KK:166:THR:HG23	6:KK:168:TYR:H	1.78	0.47
4:LG:800:ASP:OD1	4:LG:800:ASP:N	2.46	0.47
8:MM:110:THR:OG1	8:MM:111:LYS:N	2.47	0.47
3:BF:235:GLY:HA2	3:BF:243:MET:HE1	1.96	0.47
1:DC:265:ALA:HB2	2:DD:62:ILE:HD11	1.96	0.47
8:FM:125:THR:HG23	8:FM:145:ARG:HB2	1.96	0.47
2:Gd:82:ARG:HB2	1:HC:269:ASP:HB2	1.96	0.47
1:HC:247:LEU:HB2	1:IC:125:ILE:HG13	1.97	0.47
6:HK:117:THR:HG22	6:HK:157:GLN:HG3	1.97	0.47
1:IC:154:ASP:HB2	9:IN:111:GLU:CD	2.39	0.47
5:IH:191:ILE:HA	5:IH:230:LEU:HD23	1.96	0.47
7:IL:145:TYR:HA	7:IL:148:LEU:HD12	1.95	0.47
7:LL:44:HIS:HB2	7:LL:48:ASN:HA	1.96	0.47
8:LM:125:THR:HG23	8:LM:145:ARG:HB2	1.96	0.47
1:MC:210:ILE:HA	1:MC:210:ILE:HD12	1.65	0.47
6:MK:41:LEU:HD22	9:MN:76:MET:HE1	1.96	0.47
3:XF:235:GLY:HA2	3:XF:243:MET:HE1	1.96	0.47
1:EC:210:ILE:HA	1:EC:210:ILE:HD12	1.65	0.47
7:FL:44:HIS:HB2	7:FL:48:ASN:HA	1.96	0.47
9:FN:81:LEU:HD21	5:GH:357:LYS:HB2	1.97	0.47
7:IL:44:HIS:HB2	7:IL:48:ASN:HA	1.96	0.47
5:JH:311:MET:HE3	5:JH:342:SER:H	1.78	0.47
6:JK:41:LEU:HD22	9:JN:76:MET:HE1	1.95	0.47
8:KM:31:LYS:O	8:KM:76:ARG:NH2	2.43	0.47
1:AC:34:LEU:CD2	1:MC:180:THR:HG21	2.44	0.47
9:BN:63:GLY:HA3	2:CD:27:LYS:HB2	1.96	0.47
3:CF:235:GLY:HA2	3:CF:243:MET:HE1	1.96	0.47
7:CL:44:HIS:HB2	7:CL:48:ASN:HA	1.96	0.47
8:CM:319:PHE:HA	8:CM:320:PRO:HD2	1.82	0.47
2:FD:77:TYR:OH	6:FK:130:ARG:NH1	2.45	0.47
4:GG:801:MET:HB3	5:XH:115:MET:HE1	1.97	0.47
7:HL:44:HIS:HB2	7:HL:48:ASN:HA	1.96	0.47
8:HM:110:THR:OG1	8:HM:111:LYS:N	2.47	0.47
1:JC:210:ILE:HA	1:JC:210:ILE:HD12	1.66	0.47
2:MD:71:LEU:HD13	2:Md:65:PRO:HD3	1.96	0.47
7:ML:44:HIS:HB2	7:ML:48:ASN:HA	1.96	0.47
4:ZG:800:ASP:OD1	4:ZG:800:ASP:N	2.46	0.47
3:AF:235:GLY:HA2	3:AF:243:MET:HE1	1.96	0.47
8:CM:125:THR:HG23	8:CM:145:ARG:HB2	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:DC:62:THR:O	1:DC:66:VAL:HG12	2.15	0.47
1:GC:62:THR:O	1:GC:66:VAL:HG12	2.15	0.47
6:GK:117:THR:HG22	6:GK:157:GLN:HG3	1.97	0.47
9:GN:81:LEU:HD21	5:HH:357:LYS:HB2	1.95	0.47
4:HG:818:GLN:HG2	5:HH:216:GLN:NE2	2.30	0.47
5:HH:259:TYR:HD2	5:HH:266:MET:HE1	1.79	0.47
7:HL:224:ALA:HB2	2:Hd:36:ASP:OD1	2.15	0.47
3:IF:235:GLY:HA2	3:IF:243:MET:HE1	1.96	0.47
6:IK:117:THR:HG22	6:IK:157:GLN:HG3	1.96	0.47
8:IM:110:THR:OG1	8:IM:111:LYS:N	2.47	0.47
7:JL:44:HIS:HB2	7:JL:48:ASN:HA	1.96	0.47
8:JM:110:THR:OG1	8:JM:111:LYS:N	2.47	0.47
8:JM:125:THR:HG23	8:JM:145:ARG:HB2	1.96	0.47
8:KM:274:THR:HA	8:KM:293:ALA:HB3	1.95	0.47
6:LK:117:THR:HG22	6:LK:157:GLN:HG3	1.97	0.47
4:MG:800:ASP:OD1	4:MG:800:ASP:N	2.46	0.47
8:MM:125:THR:HG23	8:MM:145:ARG:HB2	1.96	0.47
3:ZF:235:GLY:HA2	3:ZF:243:MET:HE1	1.96	0.47
8:AM:124:VAL:O	8:AM:144:THR:HA	2.15	0.47
1:BC:184:TRP:CD2	1:CC:31:LEU:HB2	2.50	0.47
8:DM:124:VAL:O	8:DM:144:THR:HA	2.15	0.47
5:GH:157:PHE:HA	5:GH:255:ARG:HB2	1.97	0.47
1:IC:62:THR:O	1:IC:66:VAL:HG12	2.15	0.47
6:KK:117:THR:HG22	6:KK:157:GLN:HG3	1.96	0.47
5:MH:207:ASP:OD1	5:MH:208:LYS:N	2.48	0.47
8:MM:31:LYS:O	8:MM:76:ARG:NH2	2.43	0.47
1:BC:210:ILE:HA	1:BC:210:ILE:HD12	1.65	0.47
2:Bd:97:ARG:HA	2:Bd:97:ARG:HD3	1.73	0.47
2:Cd:98:ILE:HG23	2:Cd:128:LEU:HD23	1.97	0.47
3:GF:235:GLY:HA2	3:GF:243:MET:HE1	1.96	0.47
8:GM:124:VAL:O	8:GM:144:THR:HA	2.15	0.47
2:Gd:150:ASN:OD1	2:Gd:151:SER:N	2.46	0.47
1:JC:62:THR:O	1:JC:66:VAL:HG12	2.15	0.47
8:KM:125:THR:HG23	8:KM:145:ARG:HB2	1.96	0.47
1:MC:62:THR:O	1:MC:66:VAL:HG12	2.15	0.47
5:CH:125:GLU:HG2	5:CH:126:GLN:N	2.30	0.47
5:DH:284:LEU:HD22	5:DH:330:SER:HB3	1.97	0.47
8:FM:319:PHE:HA	8:FM:320:PRO:HD2	1.82	0.47
2:Fd:82:ARG:NH2	2:Fd:126:GLU:N	2.63	0.47
3:HF:235:GLY:HA2	3:HF:243:MET:HE1	1.96	0.47
7:IL:47:TYR:OH	7:IL:142:GLU:OE1	2.28	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:JL:47:TYR:OH	7:JL:142:GLU:OE1	2.28	0.47
5:KH:204:ILE:HG12	5:KH:215:ILE:HD12	1.97	0.47
5:KH:238:MET:HE1	5:YH:250:TYR:CG	2.50	0.47
5:MH:263:ALA:HB1	5:MH:266:MET:CE	2.45	0.47
5:VH:168:VAL:HG13	5:VH:242:ILE:HD13	1.97	0.47
5:WH:168:VAL:HG13	5:WH:242:ILE:HD13	1.95	0.47
3:YF:235:GLY:HA2	3:YF:243:MET:HE1	1.96	0.47
6:DK:117:THR:HA	6:DK:157:GLN:O	2.16	0.46
7:DL:44:HIS:HB2	7:DL:48:ASN:HA	1.96	0.46
2:Fd:97:ARG:HD3	2:Fd:97:ARG:HA	1.68	0.46
5:IH:173:GLN:HA	5:IH:217:ALA:HB3	1.96	0.46
6:IK:117:THR:HA	6:IK:157:GLN:O	2.15	0.46
1:JC:242:ILE:HG13	1:KC:130:ARG:HB2	1.97	0.46
1:LC:247:LEU:HB2	1:MC:125:ILE:HG13	1.97	0.46
3:WF:235:GLY:HA2	3:WF:243:MET:HE1	1.96	0.46
5:XH:349:TRP:HD1	5:XH:354:MET:HE1	1.78	0.46
1:AC:62:THR:O	1:AC:66:VAL:HG12	2.15	0.46
5:BH:235:THR:HG23	5:CH:251:ARG:CZ	2.45	0.46
6:BK:117:THR:HA	6:BK:157:GLN:O	2.16	0.46
2:CD:87:TRP:CG	2:CD:94:LEU:HD12	2.49	0.46
6:CK:117:THR:HG22	6:CK:157:GLN:HG3	1.96	0.46
8:DM:31:LYS:O	8:DM:76:ARG:NH2	2.43	0.46
5:EH:196:LEU:HD13	5:EH:226:LEU:HD12	1.98	0.46
7:EL:44:HIS:HB2	7:EL:48:ASN:HA	1.96	0.46
3:GF:229:THR:OG1	5:GH:165:THR:OG1	2.33	0.46
6:HK:117:THR:HA	6:HK:157:GLN:O	2.16	0.46
3:Hf:221:TRP:HE1	3:Hf:229:THR:HB	1.80	0.46
6:IK:41:LEU:HD22	9:IN:76:MET:HE1	1.97	0.46
1:LC:62:THR:O	1:LC:66:VAL:HG12	2.15	0.46
4:LG:817:THR:HA	5:LH:216:GLN:NE2	2.29	0.46
5:MH:106:VAL:HA	5:MH:109:LYS:HE2	1.96	0.46
8:AM:31:LYS:O	8:AM:76:ARG:NH2	2.43	0.46
8:AM:110:THR:OG1	8:AM:111:LYS:N	2.47	0.46
5:CH:238:MET:CE	5:VH:250:TYR:HB3	2.45	0.46
6:CK:117:THR:HA	6:CK:157:GLN:O	2.16	0.46
8:DM:267:ASN:HB3	8:DM:268:HIS:CD2	2.51	0.46
2:ED:157:ARG:HB2	2:Ed:137:TYR:OH	2.14	0.46
8:EM:124:VAL:O	8:EM:144:THR:HA	2.15	0.46
8:GM:125:THR:HG23	8:GM:145:ARG:HB2	1.96	0.46
8:JM:124:VAL:O	8:JM:144:THR:HA	2.15	0.46
1:KC:62:THR:O	1:KC:66:VAL:HG12	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:KH:171:LEU:HD11	5:KH:241:LEU:HB2	1.96	0.46
8:KM:124:VAL:O	8:KM:144:THR:HA	2.15	0.46
6:AK:117:THR:HA	6:AK:157:GLN:O	2.16	0.46
1:BC:157:LYS:HZ3	1:CC:28:THR:HG1	1.49	0.46
8:CM:124:VAL:O	8:CM:144:THR:HA	2.15	0.46
5:EH:354:MET:HE2	5:EH:354:MET:HB3	1.87	0.46
8:EM:110:THR:OG1	8:EM:111:LYS:N	2.47	0.46
1:FC:62:THR:O	1:FC:66:VAL:HG12	2.15	0.46
6:FK:117:THR:HG22	6:FK:157:GLN:HG3	1.96	0.46
6:FK:117:THR:HA	6:FK:157:GLN:O	2.16	0.46
8:FM:110:THR:OG1	8:FM:111:LYS:N	2.47	0.46
5:GH:223:TYR:HB2	5:GH:242:ILE:HG13	1.98	0.46
6:GK:117:THR:HA	6:GK:157:GLN:O	2.16	0.46
7:GL:44:HIS:HB2	7:GL:48:ASN:HA	1.96	0.46
5:HH:279:LEU:HD12	5:HH:289:PRO:HB3	1.97	0.46
8:KM:267:ASN:HB3	8:KM:268:HIS:CD2	2.51	0.46
5:MH:115:MET:HB3	5:MH:115:MET:HE2	1.55	0.46
5:MH:282:HIS:HB3	5:MH:287:VAL:HG13	1.98	0.46
1:AC:69:GLN:HE22	9:AN:114:LEU:HA	1.80	0.46
5:AH:322:PRO:HG2	5:AH:345:LEU:HD23	1.98	0.46
6:JK:117:THR:HA	6:JK:157:GLN:O	2.16	0.46
7:KL:44:HIS:HB2	7:KL:48:ASN:HA	1.96	0.46
7:KL:47:TYR:OH	7:KL:142:GLU:OE1	2.28	0.46
8:KM:110:THR:OG1	8:KM:111:LYS:N	2.47	0.46
6:AK:117:THR:HG22	6:AK:157:GLN:HG3	1.96	0.46
8:BM:267:ASN:HB3	8:BM:268:HIS:CD2	2.51	0.46
8:DM:125:THR:HG23	8:DM:145:ARG:HB2	1.96	0.46
1:HC:62:THR:O	1:HC:66:VAL:HG12	2.15	0.46
1:HC:157:LYS:HZ3	1:IC:28:THR:HB	1.79	0.46
5:HH:311:MET:SD	5:HH:341:LYS:HA	2.56	0.46
7:HL:190:MET:HE3	7:HL:190:MET:HB3	1.82	0.46
8:IM:125:THR:HG23	8:IM:145:ARG:HB2	1.96	0.46
2:KD:59:GLU:HA	2:KD:62:ILE:HG22	1.96	0.46
7:KL:190:MET:HE3	7:KL:190:MET:HB3	1.82	0.46
1:LC:210:ILE:HD11	2:LD:55:MET:HE2	1.97	0.46
5:LH:322:PRO:HG2	5:LH:345:LEU:HD23	1.98	0.46
8:LM:110:THR:OG1	8:LM:111:LYS:N	2.47	0.46
8:LM:124:VAL:O	8:LM:144:THR:HA	2.15	0.46
6:MK:117:THR:HA	6:MK:157:GLN:O	2.16	0.46
5:AH:111:ALA:O	5:AH:115:MET:HB2	2.16	0.46
8:BM:124:VAL:O	8:BM:144:THR:HA	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:CH:204:ILE:HD11	5:CH:213:LEU:HD13	1.96	0.46
6:HK:63:LYS:HB2	6:HK:63:LYS:HE2	1.84	0.46
8:HM:124:VAL:O	8:HM:144:THR:HA	2.15	0.46
2:ID:67:LYS:HB3	2:Id:62:ILE:HD11	1.97	0.46
8:KM:319:PHE:HA	8:KM:320:PRO:HD2	1.82	0.46
2:Kd:126:GLU:HG3	2:Kd:131:ILE:HG13	1.98	0.46
2:AD:24:LYS:O	7:AL:110:GLN:NE2	2.49	0.46
8:AM:125:THR:HG23	8:AM:145:ARG:HB2	1.96	0.46
1:CC:62:THR:O	1:CC:66:VAL:HG12	2.15	0.46
3:EF:235:GLY:HA2	3:EF:243:MET:HE1	1.96	0.46
5:FH:181:PHE:O	5:FH:211:ASN:ND2	2.40	0.46
8:FM:124:VAL:O	8:FM:144:THR:HA	2.15	0.46
8:IM:124:VAL:O	8:IM:144:THR:HA	2.15	0.46
2:KD:87:TRP:CD1	2:KD:89:GLY:H	2.33	0.46
8:MM:124:VAL:O	8:MM:144:THR:HA	2.15	0.46
5:AH:194:TYR:CE2	5:AH:204:ILE:HD13	2.51	0.46
8:AM:217:LYS:HA	8:AM:217:LYS:HD3	1.81	0.46
2:Ad:73:ILE:HA	2:Ad:74:PRO:HD3	1.80	0.46
8:BM:217:LYS:HA	8:BM:217:LYS:HD3	1.80	0.46
8:BM:278:ASP:OD2	8:BM:280:TYR:OH	2.27	0.46
6:DK:117:THR:HG22	6:DK:157:GLN:HG3	1.96	0.46
5:EH:322:PRO:HG2	5:EH:345:LEU:HD23	1.97	0.46
4:FG:797:ARG:O	4:FG:801:MET:HB2	2.16	0.46
7:FL:190:MET:HE3	7:FL:190:MET:HB3	1.82	0.46
2:GD:79:LEU:HA	2:GD:79:LEU:HD23	1.83	0.46
5:IH:208:LYS:NZ	4:JG:822:GLU:O	2.39	0.46
7:IL:166:ALA:HA	7:IL:206:GLU:O	2.16	0.46
2:Id:64:PRO:HG2	2:Id:67:LYS:HG2	1.97	0.46
5:KH:223:TYR:HB2	5:KH:242:ILE:HG13	1.98	0.46
7:LL:190:MET:HE3	7:LL:190:MET:HB3	1.82	0.46
2:MD:52:MET:HE2	2:MD:52:MET:HB2	1.81	0.46
3:MF:235:GLY:HA2	3:MF:243:MET:HE1	1.96	0.46
8:MM:267:ASN:HB3	8:MM:268:HIS:CD2	2.51	0.46
4:VG:797:ARG:O	4:VG:801:MET:HB2	2.16	0.46
4:ZG:797:ARG:O	4:ZG:801:MET:HB2	2.16	0.46
8:AM:54:GLY:O	8:AM:118:TYR:OH	2.33	0.46
2:Dd:55:MET:HB2	2:Dd:55:MET:HE3	1.64	0.46
4:EG:797:ARG:O	4:EG:801:MET:HB2	2.16	0.46
1:FC:210:ILE:HD12	1:FC:210:ILE:HA	1.65	0.46
7:HL:166:ALA:HA	7:HL:206:GLU:O	2.16	0.46
4:JG:797:ARG:O	4:JG:801:MET:HB2	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:JM:267:ASN:HB3	8:JM:268:HIS:CD2	2.51	0.46
6:KK:117:THR:HA	6:KK:157:GLN:O	2.16	0.46
8:LM:267:ASN:HB3	8:LM:268:HIS:CD2	2.51	0.46
4:WG:797:ARG:O	4:WG:801:MET:HB2	2.16	0.46
5:YH:191:ILE:HG13	5:YH:211:ASN:HA	1.97	0.46
2:CD:55:MET:HE1	2:Cd:55:MET:HE1	1.99	0.45
4:DG:797:ARG:O	4:DG:801:MET:HB2	2.16	0.45
6:EK:117:THR:HA	6:EK:157:GLN:O	2.16	0.45
2:GD:28:PRO:HG3	7:GL:115:GLN:HB2	1.98	0.45
2:GD:29:PRO:HG2	7:GL:225:GLN:HA	1.98	0.45
7:GL:166:ALA:HA	7:GL:206:GLU:O	2.16	0.45
7:GL:190:MET:HE3	7:GL:190:MET:HB3	1.82	0.45
8:GM:267:ASN:HB3	8:GM:268:HIS:CD2	2.51	0.45
5:HH:129:LYS:HA	5:HH:132:GLN:OE1	2.16	0.45
1:KC:88:ARG:HA	1:KC:88:ARG:HD2	1.82	0.45
5:WH:283:VAL:HG22	5:WH:289:PRO:HD3	1.97	0.45
5:WH:349:TRP:CD1	5:WH:354:MET:HE1	2.50	0.45
1:AC:88:ARG:HA	1:AC:88:ARG:HD2	1.82	0.45
4:CG:797:ARG:O	4:CG:801:MET:HB2	2.16	0.45
5:CH:245:GLN:HE21	5:CH:246:LYS:N	2.11	0.45
8:FM:267:ASN:HB3	8:FM:268:HIS:CD2	2.51	0.45
4:MG:797:ARG:O	4:MG:801:MET:HB2	2.16	0.45
4:WG:800:ASP:N	4:WG:800:ASP:OD1	2.46	0.45
5:YH:106:VAL:HA	5:YH:109:LYS:HE2	1.97	0.45
1:AC:97:ASN:C	1:AC:97:ASN:ND2	2.74	0.45
5:AH:141:LYS:HB3	5:AH:141:LYS:HE2	1.73	0.45
1:BC:62:THR:O	1:BC:66:VAL:HG12	2.15	0.45
1:DC:157:LYS:HZ1	1:EC:28:THR:CB	2.26	0.45
8:EM:54:GLY:O	8:EM:118:TYR:OH	2.33	0.45
8:IM:267:ASN:HB3	8:IM:268:HIS:CD2	2.51	0.45
2:Id:97:ARG:HD3	2:Id:97:ARG:HA	1.80	0.45
5:MH:134:TYR:OH	5:MH:138:GLU:OE2	2.28	0.45
7:ML:166:ALA:HA	7:ML:206:GLU:O	2.16	0.45
5:AH:190:PRO:HB2	5:AH:231:ARG:HG3	1.99	0.45
9:BN:81:LEU:HD21	5:CH:357:LYS:HB2	1.98	0.45
8:CM:267:ASN:HB3	8:CM:268:HIS:CD2	2.51	0.45
8:EM:267:ASN:HB3	8:EM:268:HIS:CD2	2.51	0.45
7:FL:47:TYR:OH	7:FL:142:GLU:OE1	2.28	0.45
4:GG:797:ARG:O	4:GG:801:MET:HB2	2.16	0.45
5:GH:220:LEU:HA	5:GH:220:LEU:HD23	1.80	0.45
4:LG:797:ARG:O	4:LG:801:MET:HB2	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:LL:166:ALA:HA	7:LL:206:GLU:O	2.16	0.45
2:Md:126:GLU:OE2	2:Md:130:GLU:HG3	2.16	0.45
4:YG:797:ARG:O	4:YG:801:MET:HB2	2.16	0.45
5:YH:168:VAL:HG13	5:YH:242:ILE:HD13	1.99	0.45
5:ZH:339:MET:HB2	5:ZH:339:MET:HE2	1.66	0.45
7:AL:166:ALA:HA	7:AL:206:GLU:O	2.16	0.45
1:EC:62:THR:O	1:EC:66:VAL:HG12	2.15	0.45
3:FF:219:ARG:HG2	5:WH:148:PRO:HD2	1.98	0.45
4:FG:800:ASP:OD1	4:FG:800:ASP:N	2.46	0.45
7:GL:47:TYR:OH	7:GL:142:GLU:OE1	2.28	0.45
1:HC:210:ILE:HD12	1:HC:210:ILE:HA	1.65	0.45
8:HM:319:PHE:HA	8:HM:320:PRO:HD2	1.82	0.45
2:JD:87:TRP:CD1	2:JD:89:GLY:H	2.34	0.45
3:Jf:256:THR:HG23	3:Jf:258:SER:H	1.81	0.45
2:KD:27:LYS:HA	2:KD:28:PRO:HD3	1.81	0.45
4:KG:797:ARG:O	4:KG:801:MET:HB2	2.16	0.45
6:LK:117:THR:HA	6:LK:157:GLN:O	2.16	0.45
4:AG:797:ARG:O	4:AG:801:MET:HB2	2.16	0.45
8:AM:267:ASN:HB3	8:AM:268:HIS:CD2	2.51	0.45
4:BG:797:ARG:O	4:BG:801:MET:HB2	2.16	0.45
5:EH:106:VAL:HA	5:EH:109:LYS:HE2	1.99	0.45
7:EL:166:ALA:HA	7:EL:206:GLU:O	2.16	0.45
6:GK:41:LEU:HD22	9:GN:76:MET:HE1	1.98	0.45
7:HL:47:TYR:OH	7:HL:142:GLU:OE1	2.28	0.45
4:AG:800:ASP:N	4:AG:800:ASP:OD1	2.46	0.45
6:CK:42:PRO:HD2	9:CN:76:MET:CE	2.46	0.45
8:CM:31:LYS:O	8:CM:76:ARG:NH2	2.43	0.45
7:EL:47:TYR:OH	7:EL:142:GLU:OE1	2.28	0.45
5:FH:182:LEU:HD12	5:FH:255:ARG:NH2	2.31	0.45
2:HD:79:LEU:HD23	2:HD:79:LEU:HA	1.84	0.45
2:ID:87:TRP:CD1	2:ID:89:GLY:H	2.35	0.45
4:IG:797:ARG:O	4:IG:801:MET:HB2	2.16	0.45
5:IH:250:TYR:CD2	5:XH:166:PRO:HG3	2.52	0.45
8:JM:31:LYS:O	8:JM:76:ARG:NH2	2.43	0.45
5:KH:263:ALA:O	5:KH:264:LYS:NZ	2.39	0.45
8:KM:217:LYS:HD3	8:KM:217:LYS:HA	1.81	0.45
5:LH:251:ARG:HB2	5:YH:238:MET:HE1	1.99	0.45
8:MM:217:LYS:HA	8:MM:217:LYS:HD3	1.81	0.45
2:HD:87:TRP:CG	2:HD:94:LEU:HD12	2.51	0.45
8:HM:267:ASN:HB3	8:HM:268:HIS:CD2	2.51	0.45
2:Id:69:ASN:O	2:Id:69:ASN:ND2	2.38	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:JH:173:GLN:NE2	5:JH:217:ALA:O	2.50	0.45
7:JL:166:ALA:HA	7:JL:206:GLU:O	2.16	0.45
2:LD:87:TRP:CG	2:LD:94:LEU:HD12	2.51	0.45
7:BL:166:ALA:HA	7:BL:206:GLU:O	2.16	0.45
2:CD:123:THR:OG1	2:CD:124:LYS:N	2.50	0.45
8:CM:121:ALA:HA	8:CM:141:GLU:O	2.17	0.45
5:FH:190:PRO:HB2	5:FH:231:ARG:HD2	1.98	0.45
5:JH:223:TYR:HB2	5:JH:242:ILE:HD13	1.99	0.45
6:KK:63:LYS:HB2	6:KK:63:LYS:HE2	1.84	0.45
3:Lf:212:ILE:HD11	3:Lf:215:VAL:HB	1.99	0.45
4:XG:797:ARG:O	4:XG:801:MET:HB2	2.16	0.45
5:ZH:160:LEU:HD13	5:ZH:235:THR:HG21	1.98	0.45
8:CM:110:THR:OG1	8:CM:111:LYS:N	2.47	0.45
2:Cd:130:GLU:H	2:Cd:130:GLU:CD	2.25	0.45
7:FL:166:ALA:HA	7:FL:206:GLU:O	2.16	0.45
1:GC:97:ASN:C	1:GC:97:ASN:ND2	2.74	0.45
4:HG:797:ARG:O	4:HG:801:MET:HB2	2.16	0.45
5:IH:238:MET:HE3	5:JH:251:ARG:HB2	1.98	0.45
5:LH:228:VAL:HG12	5:LH:230:LEU:HD22	1.99	0.45
8:LM:319:PHE:HA	8:LM:320:PRO:HD2	1.82	0.45
1:AC:210:ILE:HD12	1:AC:210:ILE:HA	1.65	0.44
4:AG:801:MET:CB	5:CH:115:MET:HE3	2.43	0.44
5:AH:128:VAL:O	5:AH:132:GLN:NE2	2.50	0.44
8:AM:279:ILE:O	8:AM:298:VAL:HA	2.18	0.44
2:CD:29:PRO:HG2	7:CL:225:GLN:HA	2.00	0.44
4:DG:797:ARG:CZ	5:EH:117:ARG:NH2	2.74	0.44
8:DM:279:ILE:O	8:DM:298:VAL:HA	2.17	0.44
2:Dd:24:LYS:HG3	2:Dd:25:PHE:HD2	1.82	0.44
8:EM:279:ILE:O	8:EM:298:VAL:HA	2.17	0.44
2:FD:62:ILE:HG22	2:FD:63:THR:HG23	1.99	0.44
8:FM:121:ALA:HA	8:FM:141:GLU:O	2.17	0.44
2:GD:130:GLU:OE1	2:GD:133:ARG:HD2	2.17	0.44
2:HD:107:ARG:HD2	2:Hd:130:GLU:CD	2.42	0.44
8:HM:121:ALA:HA	8:HM:141:GLU:O	2.17	0.44
1:JC:249:ILE:HG13	1:KC:123:GLN:OE1	2.17	0.44
5:JH:106:VAL:HA	5:JH:109:LYS:HE2	1.98	0.44
3:Jf:221:TRP:HE1	3:Jf:229:THR:HB	1.82	0.44
1:LC:210:ILE:HD12	1:LC:210:ILE:HA	1.66	0.44
8:MM:205:TYR:CD2	3:Mf:219:ARG:NH2	2.84	0.44
4:YG:800:ASP:OD1	4:YG:800:ASP:N	2.46	0.44
1:BC:35:GLN:HA	1:BC:35:GLN:NE2	2.31	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:DM:250:ARG:HD2	8:DM:268:HIS:HB2	2.00	0.44
9:DN:86:LEU:HD23	5:EH:344:VAL:HG21	1.98	0.44
4:GG:794:ILE:HD13	5:XH:107:ILE:HG23	1.99	0.44
3:Gf:247:ILE:HG22	3:Gf:254:ILE:HD12	1.99	0.44
8:IM:121:ALA:HA	8:IM:141:GLU:O	2.17	0.44
5:JH:189:TRP:HE1	5:JH:231:ARG:HB2	1.82	0.44
8:JM:121:ALA:HA	8:JM:141:GLU:O	2.17	0.44
7:KL:166:ALA:HA	7:KL:206:GLU:O	2.16	0.44
8:KM:279:ILE:O	8:KM:298:VAL:HA	2.18	0.44
2:LD:107:ARG:HB3	2:LD:155:GLU:HG2	1.99	0.44
2:LD:161:ILE:HD12	2:LD:161:ILE:HA	1.87	0.44
5:AH:179:LEU:O	5:AH:212:THR:HA	2.17	0.44
8:AM:121:ALA:HA	8:AM:141:GLU:O	2.17	0.44
5:CH:180:VAL:HG11	5:CH:251:ARG:HH22	1.82	0.44
6:CK:71:GLY:O	2:DD:27:LYS:NZ	2.47	0.44
8:DM:121:ALA:HA	8:DM:141:GLU:O	2.17	0.44
3:Df:210:TYR:HB3	3:Df:222:LEU:HD12	1.98	0.44
3:Df:247:ILE:HG22	3:Df:254:ILE:HG12	1.99	0.44
5:EH:113:LYS:O	5:EH:117:ARG:HB2	2.16	0.44
8:GM:279:ILE:O	8:GM:298:VAL:HA	2.18	0.44
4:HG:797:ARG:NH2	5:IH:108:ASP:OD1	2.50	0.44
8:HM:284:LEU:HD12	8:HM:284:LEU:HA	1.85	0.44
1:IC:177:CYS:CB	1:JC:35:GLN:NE2	2.80	0.44
5:KH:181:PHE:HE2	5:KH:228:VAL:HG11	1.82	0.44
2:LD:24:LYS:HE2	2:LD:24:LYS:HB2	1.85	0.44
2:Ad:64:PRO:HG2	2:Ad:67:LYS:HG2	1.98	0.44
5:BH:223:TYR:HB3	5:BH:242:ILE:HD12	1.98	0.44
6:BK:42:PRO:HD2	9:BN:76:MET:CE	2.47	0.44
2:CD:68:ASP:OD2	2:CD:70:THR:OG1	2.31	0.44
8:CM:58:VAL:HA	8:CM:106:VAL:O	2.18	0.44
7:DL:166:ALA:HA	7:DL:206:GLU:O	2.16	0.44
2:FD:79:LEU:HD23	2:FD:79:LEU:HA	1.87	0.44
1:GC:180:THR:CG2	1:HC:34:LEU:CD2	2.90	0.44
8:GM:121:ALA:HA	8:GM:141:GLU:O	2.17	0.44
8:GM:284:LEU:HD12	8:GM:284:LEU:HA	1.85	0.44
5:IH:150:LYS:HD2	3:XF:219:ARG:NH1	2.33	0.44
5:IH:151:PRO:HD2	3:XF:231:THR:HG21	2.00	0.44
2:Id:150:ASN:OD1	2:Id:151:SER:N	2.47	0.44
5:JH:208:LYS:HA	5:JH:208:LYS:HD3	1.62	0.44
2:Kd:24:LYS:HB3	2:Kd:24:LYS:HE2	1.87	0.44
8:MM:58:VAL:HA	8:MM:106:VAL:O	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:CM:76:ARG:HH11	8:CM:76:ARG:HD3	1.63	0.44
7:DL:190:MET:HE3	7:DL:190:MET:HB3	1.82	0.44
8:FM:279:ILE:O	8:FM:298:VAL:HA	2.17	0.44
5:GH:125:GLU:HG2	5:GH:126:GLN:N	2.32	0.44
8:GM:58:VAL:HA	8:GM:106:VAL:O	2.18	0.44
8:KM:250:ARG:HD2	8:KM:268:HIS:HB2	2.00	0.44
9:LN:62:ASN:OD1	9:LN:62:ASN:N	2.51	0.44
1:BC:184:TRP:CE2	1:CC:31:LEU:HB2	2.53	0.44
5:DH:220:LEU:HD23	5:DH:220:LEU:HA	1.84	0.44
5:FH:223:TYR:HB2	5:FH:242:ILE:HG13	2.00	0.44
8:IM:58:VAL:HA	8:IM:106:VAL:O	2.18	0.44
1:JC:240:ILE:HG13	1:KC:132:TYR:HB2	2.00	0.44
5:JH:266:MET:HE2	5:JH:268:THR:HA	1.99	0.44
8:JM:250:ARG:HD2	8:JM:268:HIS:HB2	2.00	0.44
8:JM:279:ILE:O	8:JM:298:VAL:HA	2.18	0.44
8:LM:58:VAL:HA	8:LM:106:VAL:O	2.18	0.44
8:MM:121:ALA:HA	8:MM:141:GLU:O	2.17	0.44
8:AM:279:ILE:HB	8:AM:298:VAL:HG22	2.00	0.44
7:BL:47:TYR:OH	7:BL:142:GLU:OE1	2.28	0.44
5:CH:248:VAL:HG13	5:CH:250:TYR:CE1	2.53	0.44
8:CM:70:MET:O	8:CM:106:VAL:HA	2.18	0.44
4:DG:800:ASP:OD1	4:DG:800:ASP:N	2.46	0.44
8:DM:58:VAL:HA	8:DM:106:VAL:O	2.18	0.44
5:EH:196:LEU:HD22	5:EH:204:ILE:HD13	2.00	0.44
8:FM:58:VAL:HA	8:FM:106:VAL:O	2.18	0.44
8:FM:250:ARG:HD2	8:FM:268:HIS:HB2	2.00	0.44
8:FM:279:ILE:HB	8:FM:298:VAL:HG22	2.00	0.44
5:HH:196:LEU:HD11	5:HH:202:PHE:HD2	1.83	0.44
5:IH:150:LYS:HD2	3:XF:219:ARG:NH2	2.30	0.44
8:JM:279:ILE:HB	8:JM:298:VAL:HG22	2.00	0.44
2:MD:150:ASN:OD1	2:MD:150:ASN:N	2.50	0.44
5:MH:166:PRO:HA	5:MH:167:PRO:HD3	1.89	0.44
8:MM:279:ILE:O	8:MM:298:VAL:HA	2.18	0.44
8:MM:284:LEU:HD12	8:MM:284:LEU:HA	1.85	0.44
5:BH:307:SER:OG	5:BH:308:ASN:N	2.51	0.44
8:BM:121:ALA:HA	8:BM:141:GLU:O	2.17	0.44
2:CD:87:TRP:CD1	2:CD:89:GLY:H	2.35	0.44
7:CL:166:ALA:HA	7:CL:206:GLU:O	2.16	0.44
8:CM:279:ILE:HB	8:CM:298:VAL:HG22	2.00	0.44
8:DM:279:ILE:HB	8:DM:298:VAL:HG22	2.00	0.44
2:ED:130:GLU:OE1	2:ED:133:ARG:HD2	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:EL:190:MET:HE3	7:EL:190:MET:HB3	1.82	0.44
8:EM:121:ALA:HA	8:EM:141:GLU:O	2.17	0.44
7:FL:124:THR:HA	7:FL:231:ILE:O	2.18	0.44
8:GM:279:ILE:HB	8:GM:298:VAL:HG22	2.00	0.44
8:IM:250:ARG:HD2	8:IM:268:HIS:HB2	2.00	0.44
8:JM:217:LYS:HD3	8:JM:217:LYS:HA	1.81	0.44
7:LL:124:THR:HA	7:LL:231:ILE:O	2.18	0.44
8:LM:279:ILE:HB	8:LM:298:VAL:HG22	2.00	0.44
8:MM:70:MET:O	8:MM:106:VAL:HA	2.18	0.44
8:MM:279:ILE:HB	8:MM:298:VAL:HG22	2.00	0.44
8:BM:58:VAL:HA	8:BM:106:VAL:O	2.18	0.44
2:Bd:150:ASN:OD1	2:Bd:151:SER:N	2.46	0.44
2:CD:77:TYR:OH	6:CK:130:ARG:NH1	2.47	0.44
8:CM:279:ILE:O	8:CM:298:VAL:HA	2.17	0.44
9:EN:62:ASN:OD1	9:EN:62:ASN:N	2.51	0.44
2:Ed:150:ASN:OD1	2:Ed:151:SER:N	2.44	0.44
6:FK:144:LEU:HD23	6:FK:144:LEU:HA	1.85	0.44
8:FM:217:LYS:HA	8:FM:217:LYS:HD3	1.80	0.44
2:Fd:150:ASN:OD1	2:Fd:151:SER:N	2.44	0.44
1:KC:97:ASN:C	1:KC:97:ASN:ND2	2.74	0.44
8:KM:61:HIS:HB2	8:KM:109:LYS:HG2	2.00	0.44
3:Kf:210:TYR:HB3	3:Kf:222:LEU:HD11	2.00	0.44
8:LM:70:MET:O	8:LM:106:VAL:HA	2.18	0.44
8:LM:279:ILE:O	8:LM:298:VAL:HA	2.18	0.44
5:MH:193:ALA:HB1	4:ZG:821:THR:O	2.17	0.44
8:AM:58:VAL:HA	8:AM:106:VAL:O	2.18	0.43
5:BH:213:LEU:C	5:BH:214:MET:HE2	2.43	0.43
3:Cf:251:GLN:HB3	3:Cf:253:ARG:HG3	1.99	0.43
1:DC:75:ARG:NH2	2:DD:36:ASP:OD1	2.51	0.43
1:DC:157:LYS:CE	1:EC:28:THR:HG1	2.27	0.43
8:FM:61:HIS:HB2	8:FM:109:LYS:HG2	2.00	0.43
2:GD:87:TRP:CG	2:GD:94:LEU:HD12	2.53	0.43
1:HC:240:ILE:HG12	1:IC:255:LEU:HD13	2.00	0.43
2:ID:79:LEU:HD23	2:ID:79:LEU:HA	1.85	0.43
6:IK:144:LEU:HD23	6:IK:144:LEU:HA	1.85	0.43
3:JF:219:ARG:HH22	5:KH:150:LYS:HG3	1.83	0.43
8:JM:58:VAL:HA	8:JM:106:VAL:O	2.18	0.43
8:JM:84:ILE:HG22	8:JM:91:TYR:HB2	2.00	0.43
7:KL:98:TYR:HE1	7:KL:104:LYS:HE3	1.83	0.43
5:LH:307:SER:OG	5:LH:308:ASN:N	2.51	0.43
8:LM:121:ALA:HA	8:LM:141:GLU:O	2.17	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:MH:130:LEU:HA	5:MH:133:ILE:HG22	1.99	0.43
5:XH:157:PHE:CZ	5:XH:257:GLN:HG2	2.53	0.43
1:AC:228:HIS:HB3	1:AC:247:LEU:HG	2.00	0.43
2:AD:128:LEU:HA	2:AD:128:LEU:HD13	1.82	0.43
5:BH:110:LYS:HG3	4:ZG:794:ILE:HD12	2.01	0.43
7:CL:47:TYR:OH	7:CL:142:GLU:OE1	2.28	0.43
7:CL:98:TYR:HE1	7:CL:104:LYS:HE3	1.83	0.43
8:CM:275:ASP:HB3	8:CM:294:PHE:HB2	2.01	0.43
5:DH:171:LEU:HD11	5:DH:241:LEU:HD23	1.99	0.43
8:DM:284:LEU:HD12	8:DM:284:LEU:HA	1.85	0.43
2:Dd:150:ASN:OD1	2:Dd:151:SER:N	2.48	0.43
2:ED:57:LYS:O	2:ED:60:LYS:HG3	2.18	0.43
8:EM:306:GLN:CD	3:Ef:251:GLN:HG2	2.43	0.43
2:Ed:97:ARG:HD3	2:Ed:97:ARG:HA	1.77	0.43
5:FH:284:LEU:HD22	5:FH:330:SER:HB3	1.99	0.43
7:FL:98:TYR:HE1	7:FL:104:LYS:HE3	1.83	0.43
8:FM:84:ILE:HG22	8:FM:91:TYR:HB2	2.00	0.43
1:GC:213:ARG:HD2	1:GC:213:ARG:HA	1.83	0.43
8:GM:84:ILE:HG22	8:GM:91:TYR:HB2	2.00	0.43
2:Gd:111:LYS:HB3	2:Gd:111:LYS:HE3	1.76	0.43
2:HD:87:TRP:CD1	2:HD:89:GLY:H	2.37	0.43
7:HL:124:THR:HA	7:HL:231:ILE:O	2.18	0.43
7:IL:124:THR:HA	7:IL:231:ILE:O	2.18	0.43
8:IM:279:ILE:O	8:IM:298:VAL:HA	2.18	0.43
2:Id:73:ILE:HA	2:Id:74:PRO:HD3	1.82	0.43
8:KM:121:ALA:HA	8:KM:141:GLU:O	2.17	0.43
8:KM:307:SER:OG	3:Kf:253:ARG:NH2	2.35	0.43
2:LD:87:TRP:CD1	2:LD:89:GLY:H	2.36	0.43
1:MC:228:HIS:HB3	1:MC:247:LEU:HG	2.00	0.43
7:ML:98:TYR:HE1	7:ML:104:LYS:HE3	1.83	0.43
7:ML:124:THR:HA	7:ML:231:ILE:O	2.18	0.43
7:AL:47:TYR:OH	7:AL:142:GLU:OE1	2.28	0.43
8:EM:319:PHE:HA	8:EM:320:PRO:HD2	1.82	0.43
2:FD:29:PRO:HG2	7:FL:225:GLN:HA	1.99	0.43
8:FM:70:MET:O	8:FM:106:VAL:HA	2.18	0.43
5:GH:250:TYR:HB3	5:WH:238:MET:CE	2.48	0.43
5:HH:166:PRO:HA	5:HH:167:PRO:HD3	1.87	0.43
8:HM:279:ILE:O	8:HM:298:VAL:HA	2.18	0.43
1:IC:97:ASN:C	1:IC:97:ASN:ND2	2.74	0.43
4:JG:808:LEU:HD13	4:JG:808:LEU:HA	1.88	0.43
5:JH:115:MET:HE1	4:XG:802:LEU:HA	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:Jd:150:ASN:OD1	2:Jd:151:SER:N	2.45	0.43
4:KG:797:ARG:HH12	5:YH:121:PRO:HA	1.84	0.43
8:KM:245:TYR:CD1	3:Kf:215:VAL:O	2.71	0.43
3:LF:265:SER:O	3:LF:269:SER:HB3	2.19	0.43
5:LH:184:SER:N	5:LH:256:VAL:O	2.52	0.43
9:LN:81:LEU:HD21	5:MH:357:LYS:HB2	2.00	0.43
5:AH:166:PRO:HA	5:AH:167:PRO:HD3	1.90	0.43
7:AL:98:TYR:HE1	7:AL:104:LYS:HE3	1.83	0.43
8:BM:250:ARG:HD2	8:BM:268:HIS:HB2	2.00	0.43
8:BM:279:ILE:O	8:BM:298:VAL:HA	2.18	0.43
1:DC:210:ILE:HA	1:DC:210:ILE:HD12	1.65	0.43
7:DL:98:TYR:HE1	7:DL:104:LYS:HE3	1.83	0.43
8:DM:70:MET:O	8:DM:106:VAL:HA	2.18	0.43
8:EM:58:VAL:HA	8:EM:106:VAL:O	2.18	0.43
8:FM:275:ASP:HB3	8:FM:294:PHE:HB2	2.01	0.43
2:GD:87:TRP:CD1	2:GD:89:GLY:H	2.36	0.43
2:HD:27:LYS:HA	2:HD:28:PRO:HD3	1.84	0.43
5:HH:169:ILE:HD11	5:HH:239:LEU:HD12	2.01	0.43
8:HM:279:ILE:HB	8:HM:298:VAL:HG22	2.00	0.43
4:IG:819:VAL:HG13	5:XH:199:PRO:HG2	2.00	0.43
8:IM:279:ILE:HB	8:IM:298:VAL:HG22	2.00	0.43
1:JC:118:ASN:ND2	5:JH:274:PRO:O	2.47	0.43
3:JF:265:SER:O	3:JF:269:SER:HB3	2.19	0.43
4:JG:815:VAL:HG21	5:JH:173:GLN:NE2	2.33	0.43
8:JM:70:MET:O	8:JM:106:VAL:HA	2.18	0.43
5:LH:130:LEU:HA	5:LH:133:ILE:HG22	1.99	0.43
8:MM:275:ASP:HB3	8:MM:294:PHE:HB2	2.01	0.43
3:Mf:210:TYR:HB3	3:Mf:222:LEU:HD23	2.01	0.43
3:VF:265:SER:O	3:VF:269:SER:HB3	2.19	0.43
5:XH:230:LEU:HD12	5:XH:233:LEU:HD12	2.01	0.43
3:BF:265:SER:O	3:BF:269:SER:HB3	2.19	0.43
7:BL:124:THR:HA	7:BL:231:ILE:O	2.18	0.43
5:CH:179:LEU:O	5:CH:212:THR:HA	2.18	0.43
8:DM:127:ASN:HB3	8:DM:147:ALA:HB3	2.01	0.43
2:Dd:83:ALA:HA	1:EC:267:ALA:O	2.18	0.43
3:FF:265:SER:O	3:FF:269:SER:HB3	2.19	0.43
5:FH:169:ILE:HD11	5:FH:239:LEU:HD12	2.00	0.43
2:Fd:24:LYS:HG3	2:Fd:25:PHE:CD1	2.51	0.43
6:GK:63:LYS:HB2	6:GK:63:LYS:HE2	1.84	0.43
8:IM:70:MET:O	8:IM:106:VAL:HA	2.18	0.43
8:IM:284:LEU:HD12	8:IM:284:LEU:HA	1.85	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:KC:213:ARG:HD2	1:KC:213:ARG:HA	1.82	0.43
6:KK:144:LEU:HD23	6:KK:144:LEU:HA	1.85	0.43
4:LG:815:VAL:HG11	5:LH:218:THR:O	2.18	0.43
8:LM:84:ILE:HG22	8:LM:91:TYR:HB2	2.00	0.43
8:LM:284:LEU:HD12	8:LM:284:LEU:HA	1.85	0.43
5:MH:122:LEU:HD12	5:MH:126:GLN:HB3	2.00	0.43
5:WH:279:LEU:HD12	5:WH:280:LEU:HD22	2.00	0.43
5:XH:169:ILE:HB	5:XH:241:LEU:HD23	2.01	0.43
6:AK:63:LYS:HB2	6:AK:63:LYS:HE2	1.84	0.43
5:EH:200:SER:N	4:FG:816:GLU:OE2	2.48	0.43
8:EM:250:ARG:HD2	8:EM:268:HIS:HB2	2.00	0.43
8:FM:127:ASN:HB3	8:FM:147:ALA:HB3	2.01	0.43
7:GL:124:THR:HA	7:GL:231:ILE:O	2.18	0.43
2:HD:162:TYR:HB3	1:IC:127:ILE:HD13	2.00	0.43
8:HM:58:VAL:HA	8:HM:106:VAL:O	2.18	0.43
8:IM:84:ILE:HG22	8:IM:91:TYR:HB2	2.00	0.43
5:JH:194:TYR:CE2	5:JH:204:ILE:HD13	2.53	0.43
1:KC:58:ALA:O	1:KC:62:THR:HG23	2.19	0.43
2:KD:67:LYS:HB3	2:Kd:62:ILE:HD11	1.99	0.43
8:KM:58:VAL:HA	8:KM:106:VAL:O	2.18	0.43
3:Kf:210:TYR:CG	3:Kf:222:LEU:HD21	2.54	0.43
1:MC:265:ALA:HB2	2:MD:62:ILE:HD11	2.00	0.43
5:YH:171:LEU:HD11	5:YH:241:LEU:HB3	2.01	0.43
8:AM:61:HIS:HB2	8:AM:109:LYS:HG2	2.00	0.43
8:AM:70:MET:O	8:AM:106:VAL:HA	2.18	0.43
8:AM:84:ILE:HG22	8:AM:91:TYR:HB2	2.00	0.43
2:BD:60:LYS:HD2	2:Bd:54:GLU:OE2	2.19	0.43
8:BM:70:MET:O	8:BM:106:VAL:HA	2.18	0.43
2:CD:27:LYS:HA	2:CD:28:PRO:HD3	1.84	0.43
5:CH:191:ILE:HG13	5:CH:211:ASN:HA	2.01	0.43
8:DM:61:HIS:HB2	8:DM:109:LYS:HG2	2.00	0.43
8:DM:217:LYS:HA	8:DM:217:LYS:HD3	1.81	0.43
8:DM:275:ASP:HB3	8:DM:294:PHE:HB2	2.01	0.43
5:EH:145:PRO:HB3	5:FH:132:GLN:NE2	2.33	0.43
7:EL:124:THR:HA	7:EL:231:ILE:O	2.18	0.43
8:EM:84:ILE:HG22	8:EM:91:TYR:HB2	2.00	0.43
8:EM:217:LYS:HD3	8:EM:217:LYS:HA	1.80	0.43
5:GH:307:SER:OG	5:GH:308:ASN:N	2.52	0.43
1:HC:58:ALA:O	1:HC:62:THR:HG23	2.19	0.43
8:HM:61:HIS:HB2	8:HM:109:LYS:HG2	2.00	0.43
3:Hf:243:MET:HB3	3:Hf:257:SER:HB3	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:JD:27:LYS:HA	2:JD:28:PRO:HD3	1.85	0.43
2:Kd:91:ILE:O	2:Kd:95:THR:HG22	2.18	0.43
2:LD:77:TYR:OH	6:LK:130:ARG:NH1	2.47	0.43
8:MM:250:ARG:HD2	8:MM:268:HIS:HB2	2.00	0.43
3:XF:265:SER:O	3:XF:269:SER:HB3	2.19	0.43
5:ZH:324:TRP:HA	5:ZH:339:MET:HE1	2.00	0.43
8:CM:127:ASN:HB3	8:CM:147:ALA:HB3	2.01	0.43
8:CM:217:LYS:HA	8:CM:217:LYS:HD3	1.81	0.43
2:DD:27:LYS:HA	2:DD:28:PRO:HD3	1.88	0.43
6:EK:144:LEU:HD23	6:EK:144:LEU:HA	1.85	0.43
7:EL:98:TYR:HE1	7:EL:104:LYS:HE3	1.83	0.43
8:EM:257:ASP:OD1	8:EM:257:ASP:N	2.52	0.43
1:FC:213:ARG:HD2	1:FC:213:ARG:HA	1.83	0.43
2:Fd:134:ASP:OD1	2:Fd:138:GLN:NE2	2.48	0.43
8:GM:70:MET:O	8:GM:106:VAL:HA	2.18	0.43
8:GM:275:ASP:HB3	8:GM:294:PHE:HB2	2.01	0.43
2:Gd:24:LYS:HG3	2:Gd:25:PHE:HD2	1.83	0.43
2:Gd:157:ARG:HE	2:Gd:157:ARG:HB2	1.61	0.43
8:IM:275:ASP:HB3	8:IM:294:PHE:HB2	2.01	0.43
3:If:254:ILE:HG13	3:If:262:ILE:HB	2.00	0.43
1:JC:228:HIS:HB3	1:JC:247:LEU:HG	2.00	0.43
5:JH:156:GLN:HB2	5:JH:254:LEU:HD13	1.99	0.43
8:KM:279:ILE:HB	8:KM:298:VAL:HG22	2.00	0.43
2:Kd:73:ILE:HA	2:Kd:74:PRO:HD3	1.88	0.43
2:Kd:150:ASN:OD1	2:Kd:151:SER:N	2.46	0.43
7:LL:98:TYR:HE1	7:LL:104:LYS:HE3	1.83	0.43
8:LM:257:ASP:OD1	8:LM:257:ASP:N	2.52	0.43
5:WH:166:PRO:HA	5:WH:167:PRO:HD3	1.89	0.43
1:AC:35:GLN:CD	1:MC:176:TRP:CZ3	2.97	0.43
1:BC:58:ALA:O	1:BC:62:THR:HG23	2.19	0.43
8:BM:61:HIS:HB2	8:BM:109:LYS:HG2	2.00	0.43
3:Bf:219:ARG:NH2	3:Bf:233:ARG:HH22	2.16	0.43
1:CC:58:ALA:O	1:CC:62:THR:HG23	2.19	0.43
2:CD:155:GLU:OE1	2:Cd:133:ARG:NH2	2.51	0.43
5:CH:307:SER:OG	5:CH:308:ASN:N	2.51	0.43
1:DC:228:HIS:HB3	1:DC:247:LEU:HG	2.01	0.43
1:EC:58:ALA:O	1:EC:62:THR:HG23	2.19	0.43
6:EK:63:LYS:HB2	6:EK:63:LYS:HE2	1.84	0.43
5:HH:132:GLN:HA	5:HH:135:GLU:CD	2.44	0.43
6:HK:71:GLY:O	2:ID:27:LYS:NZ	2.37	0.43
7:KL:124:THR:HA	7:KL:231:ILE:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:KM:257:ASP:OD1	8:KM:257:ASP:N	2.52	0.43
1:LC:69:GLN:HE22	9:LN:114:LEU:HA	1.82	0.43
5:LH:223:TYR:HB2	5:LH:242:ILE:HD13	2.01	0.43
3:MF:265:SER:O	3:MF:269:SER:HB3	2.19	0.43
5:MH:208:LYS:HE3	5:MH:208:LYS:HB3	1.71	0.43
8:MM:127:ASN:HB3	8:MM:147:ALA:HB3	2.01	0.43
5:YH:311:MET:SD	5:YH:311:MET:N	2.90	0.43
5:YH:335:HIS:HB3	5:YH:337:TYR:CE1	2.53	0.43
4:AG:808:LEU:CD1	5:AH:141:LYS:HE3	2.45	0.43
8:AM:275:ASP:HB3	8:AM:294:PHE:HB2	2.01	0.43
7:CL:111:LYS:HB2	7:CL:111:LYS:HE3	1.86	0.43
8:CM:84:ILE:HG22	8:CM:91:TYR:HB2	2.00	0.43
9:CN:62:ASN:OD1	9:CN:62:ASN:N	2.51	0.43
6:DK:65:ILE:HD12	6:DK:78:ILE:HG23	2.01	0.43
7:DL:124:THR:HA	7:DL:231:ILE:O	2.18	0.43
1:EC:116:THR:HG23	5:EH:328:MET:HG2	2.01	0.43
2:ED:87:TRP:CD1	2:ED:89:GLY:H	2.36	0.43
7:EL:111:LYS:HE3	7:EL:111:LYS:HB2	1.86	0.43
8:FM:278:ASP:OD2	8:FM:280:TYR:OH	2.27	0.43
1:GC:210:ILE:HA	1:GC:210:ILE:HD12	1.65	0.43
1:IC:58:ALA:O	1:IC:62:THR:HG23	2.19	0.43
8:MM:84:ILE:HG22	8:MM:91:TYR:HB2	2.00	0.43
8:AM:250:ARG:HD2	8:AM:268:HIS:HB2	2.00	0.42
2:Ad:94:LEU:HD22	2:Ad:135:ILE:HD11	2.00	0.42
1:CC:228:HIS:HB3	1:CC:247:LEU:HG	2.00	0.42
5:CH:199:PRO:HG2	4:VG:819:VAL:HG13	2.00	0.42
8:CM:250:ARG:HD2	8:CM:268:HIS:HB2	2.00	0.42
8:CM:284:LEU:HD12	8:CM:284:LEU:HA	1.85	0.42
3:DF:265:SER:O	3:DF:269:SER:HB3	2.19	0.42
5:DH:191:ILE:HG13	5:DH:211:ASN:HA	2.01	0.42
8:DM:84:ILE:HG22	8:DM:91:TYR:HB2	2.00	0.42
2:ED:87:TRP:CG	2:ED:94:LEU:HD12	2.54	0.42
1:FC:154:ASP:HB3	9:FN:84:CYS:HB3	2.01	0.42
2:FD:161:ILE:HD12	2:FD:161:ILE:HA	1.93	0.42
8:FM:71:LEU:HD22	8:FM:81:VAL:HG11	2.02	0.42
6:GK:65:ILE:HD12	6:GK:78:ILE:HG23	2.01	0.42
8:GM:49:GLN:HG3	8:GM:115:ARG:HB3	2.01	0.42
8:GM:71:LEU:HD22	8:GM:81:VAL:HG11	2.01	0.42
8:GM:250:ARG:HD2	8:GM:268:HIS:HB2	2.00	0.42
2:HD:29:PRO:HG2	7:HL:225:GLN:HA	2.01	0.42
2:Hd:97:ARG:HA	2:Hd:97:ARG:HD3	1.74	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:IM:61:HIS:HB2	8:IM:109:LYS:HG2	2.00	0.42
8:IM:127:ASN:HB3	8:IM:147:ALA:HB3	2.01	0.42
8:JM:61:HIS:HB2	8:JM:109:LYS:HG2	2.00	0.42
5:KH:181:PHE:HD2	5:KH:230:LEU:HD21	1.84	0.42
1:LC:228:HIS:HB3	1:LC:247:LEU:HG	2.01	0.42
2:MD:87:TRP:CD1	2:MD:89:GLY:H	2.37	0.42
4:MG:797:ARG:HH12	5:ZH:121:PRO:HA	1.83	0.42
2:AD:27:LYS:HA	2:AD:28:PRO:HD3	1.82	0.42
2:AD:161:ILE:HD12	2:AD:161:ILE:HA	1.88	0.42
4:AG:808:LEU:HD13	4:AG:808:LEU:HA	1.88	0.42
6:AK:65:ILE:HD12	6:AK:78:ILE:HG23	2.01	0.42
5:BH:166:PRO:HA	5:BH:167:PRO:HD3	1.94	0.42
7:CL:124:THR:HA	7:CL:231:ILE:O	2.18	0.42
8:EM:70:MET:O	8:EM:106:VAL:HA	2.18	0.42
3:GF:219:ARG:HH22	5:HH:150:LYS:CD	2.32	0.42
5:GH:131:LYS:NZ	5:WH:143:ALA:O	2.51	0.42
6:HK:65:ILE:HD12	6:HK:78:ILE:HG23	2.01	0.42
2:ID:28:PRO:HG3	7:IL:115:GLN:HB2	1.99	0.42
2:Id:113:PRO:HD3	2:Id:158:TYR:CE2	2.52	0.42
1:JC:157:LYS:HZ3	1:KC:28:THR:CA	2.32	0.42
8:JM:275:ASP:HB3	8:JM:294:PHE:HB2	2.01	0.42
8:KM:70:MET:O	8:KM:106:VAL:HA	2.18	0.42
8:LM:250:ARG:HD2	8:LM:268:HIS:HB2	2.00	0.42
8:MM:61:HIS:HB2	8:MM:109:LYS:HG2	2.00	0.42
5:ZH:319:ILE:HA	5:ZH:347:VAL:HG22	2.00	0.42
7:AL:124:THR:HA	7:AL:231:ILE:O	2.18	0.42
8:AM:127:ASN:HB3	8:AM:147:ALA:HB3	2.01	0.42
1:BC:88:ARG:HA	1:BC:88:ARG:HD2	1.82	0.42
1:BC:228:HIS:HB3	1:BC:247:LEU:HG	2.01	0.42
5:CH:106:VAL:HA	5:CH:109:LYS:HE2	2.01	0.42
2:Cd:50:ASP:O	2:Cd:54:GLU:HG2	2.19	0.42
6:EK:65:ILE:HD12	6:EK:78:ILE:HG23	2.01	0.42
3:Ef:219:ARG:HG2	3:Ef:233:ARG:HD2	2.00	0.42
3:GF:265:SER:O	3:GF:269:SER:HB3	2.19	0.42
5:GH:219:LYS:HE2	5:GH:219:LYS:HB2	1.81	0.42
1:IC:228:HIS:HB3	1:IC:247:LEU:HG	2.00	0.42
3:IF:265:SER:O	3:IF:269:SER:HB3	2.19	0.42
4:IG:800:ASP:N	4:IG:800:ASP:OD1	2.46	0.42
1:MC:58:ALA:O	1:MC:62:THR:HG23	2.19	0.42
5:WH:203:ASN:OD1	5:WH:216:GLN:HB3	2.20	0.42
4:XG:800:ASP:N	4:XG:800:ASP:OD1	2.46	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AC:35:GLN:NE2	1:MC:176:TRP:HE3	2.14	0.42
6:BK:144:LEU:HD23	6:BK:144:LEU:HA	1.85	0.42
8:BM:84:ILE:HG22	8:BM:91:TYR:HB2	2.00	0.42
3:CF:265:SER:O	3:CF:269:SER:HB3	2.19	0.42
4:DG:797:ARG:NH1	5:EH:117:ARG:HH12	2.16	0.42
4:DG:819:VAL:HG13	5:VH:199:PRO:HG2	2.01	0.42
7:DL:47:TYR:OH	7:DL:142:GLU:OE1	2.28	0.42
1:EC:228:HIS:HB3	1:EC:247:LEU:HG	2.00	0.42
5:EH:122:LEU:HD13	5:EH:130:LEU:HD21	2.00	0.42
5:EH:127:VAL:HA	5:EH:130:LEU:HG	2.01	0.42
5:EH:289:PRO:O	5:EH:292:SER:OG	2.35	0.42
1:FC:58:ALA:O	1:FC:62:THR:HG23	2.19	0.42
1:FC:88:ARG:HA	1:FC:88:ARG:HD2	1.82	0.42
5:IH:179:LEU:O	5:IH:212:THR:HA	2.19	0.42
7:JL:124:THR:HA	7:JL:231:ILE:O	2.18	0.42
7:JL:190:MET:HE3	7:JL:190:MET:HB3	1.82	0.42
2:KD:55:MET:HE2	2:KD:55:MET:HB2	1.95	0.42
5:KH:196:LEU:HD13	5:KH:226:LEU:HD12	2.01	0.42
8:LM:127:ASN:HB3	8:LM:147:ALA:HB3	2.01	0.42
5:MH:183:ASP:O	5:MH:255:ARG:HD3	2.20	0.42
5:MH:307:SER:OG	5:MH:308:ASN:N	2.51	0.42
8:MM:257:ASP:OD1	8:MM:257:ASP:N	2.52	0.42
5:YH:179:LEU:O	5:YH:212:THR:HA	2.19	0.42
3:ZF:265:SER:O	3:ZF:269:SER:HB3	2.19	0.42
2:BD:87:TRP:CD1	2:BD:89:GLY:H	2.38	0.42
6:BK:65:ILE:HD12	6:BK:78:ILE:HG23	2.01	0.42
8:BM:279:ILE:HB	8:BM:298:VAL:HG22	2.00	0.42
3:Bf:212:ILE:HG22	3:Bf:222:LEU:HG	2.00	0.42
5:CH:151:PRO:HB3	5:CH:250:TYR:CE2	2.54	0.42
8:CM:257:ASP:OD1	8:CM:257:ASP:N	2.52	0.42
2:DD:87:TRP:CD1	2:DD:89:GLY:H	2.37	0.42
3:EF:265:SER:O	3:EF:269:SER:HB3	2.19	0.42
2:Fd:73:ILE:HA	2:Fd:74:PRO:HD3	1.81	0.42
1:GC:58:ALA:O	1:GC:62:THR:HG23	2.19	0.42
8:HM:70:MET:O	8:HM:106:VAL:HA	2.18	0.42
1:IC:265:ALA:HB2	2:ID:62:ILE:HD11	2.01	0.42
5:IH:166:PRO:HA	5:IH:167:PRO:HD3	1.92	0.42
5:IH:213:LEU:C	5:IH:214:MET:HE2	2.44	0.42
7:IL:98:TYR:HE1	7:IL:104:LYS:HE3	1.83	0.42
5:JH:173:GLN:HE22	5:JH:218:THR:HA	1.84	0.42
4:KG:801:MET:C	5:LH:115:MET:HE1	2.44	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:MM:49:GLN:HG3	8:MM:115:ARG:HB3	2.02	0.42
3:WF:265:SER:O	3:WF:269:SER:HB3	2.19	0.42
3:YF:265:SER:O	3:YF:269:SER:HB3	2.19	0.42
3:AF:265:SER:O	3:AF:269:SER:HB3	2.19	0.42
1:BC:114:ARG:HG2	1:BC:130:ARG:HG2	2.02	0.42
8:CM:61:HIS:HB2	8:CM:109:LYS:HG2	2.00	0.42
1:DC:58:ALA:O	1:DC:62:THR:HG23	2.19	0.42
5:EH:183:ASP:O	5:EH:255:ARG:HD3	2.20	0.42
8:EM:279:ILE:HB	8:EM:298:VAL:HG22	2.00	0.42
1:FC:228:HIS:HB3	1:FC:247:LEU:HG	2.00	0.42
2:FD:82:ARG:HH22	8:FM:311:ASP:CG	2.28	0.42
5:FH:169:ILE:HB	5:FH:241:LEU:HD23	2.01	0.42
1:GC:229:THR:HG22	5:XH:323:GLY:HA3	2.01	0.42
2:GD:55:MET:HE3	2:GD:55:MET:HB3	1.78	0.42
1:HC:114:ARG:HG2	1:HC:130:ARG:HG2	2.02	0.42
1:HC:228:HIS:HB3	1:HC:247:LEU:HG	2.01	0.42
3:HF:265:SER:O	3:HF:269:SER:HB3	2.19	0.42
5:IH:181:PHE:HD2	5:IH:230:LEU:HD11	1.83	0.42
5:IH:199:PRO:HG2	4:JG:819:VAL:HG13	2.01	0.42
8:JM:284:LEU:HD12	8:JM:284:LEU:HA	1.85	0.42
2:Jd:97:ARG:HD3	2:Jd:97:ARG:HA	1.68	0.42
5:KH:179:LEU:O	5:KH:212:THR:HA	2.19	0.42
6:KK:41:LEU:HD22	9:KN:76:MET:HE1	2.01	0.42
8:KM:306:GLN:CD	3:Kf:251:GLN:HB3	2.45	0.42
2:LD:35:ASP:OD1	2:LD:37:ALA:N	2.43	0.42
8:LM:61:HIS:HB2	8:LM:109:LYS:HG2	2.00	0.42
8:LM:247:ARG:HD3	3:Lf:213:GLN:O	2.19	0.42
8:LM:275:ASP:HB3	8:LM:294:PHE:HB2	2.01	0.42
8:BM:71:LEU:HD22	8:BM:81:VAL:HG11	2.01	0.42
5:CH:122:LEU:HA	5:CH:122:LEU:HD23	1.86	0.42
1:EC:114:ARG:HG2	1:EC:130:ARG:HG2	2.02	0.42
5:FH:170:ARG:HH11	5:FH:170:ARG:HD2	1.71	0.42
2:Hd:132:LEU:HD23	2:Hd:132:LEU:HA	1.87	0.42
6:IK:65:ILE:HD12	6:IK:78:ILE:HG23	2.01	0.42
7:IL:190:MET:HE3	7:IL:190:MET:HB3	1.82	0.42
8:KM:84:ILE:HG22	8:KM:91:TYR:HB2	2.00	0.42
2:Kd:97:ARG:HD3	2:Kd:97:ARG:HA	1.81	0.42
1:LC:58:ALA:O	1:LC:62:THR:HG23	2.19	0.42
5:LH:229:ARG:HG2	5:LH:236:PRO:HB3	2.01	0.42
6:LK:65:ILE:HD12	6:LK:78:ILE:HG23	2.01	0.42
1:AC:58:ALA:O	1:AC:62:THR:HG23	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:AM:49:GLN:HG3	8:AM:115:ARG:HB3	2.02	0.42
5:CH:158:VAL:HG23	5:CH:254:LEU:HD12	2.01	0.42
5:DH:273:PRO:HA	5:DH:274:PRO:HD3	1.89	0.42
5:DH:307:SER:OG	5:DH:308:ASN:N	2.52	0.42
6:DK:63:LYS:HB2	6:DK:63:LYS:HE2	1.84	0.42
8:DM:69:VAL:O	8:DM:70:MET:HE2	2.20	0.42
5:EH:166:PRO:HA	5:EH:167:PRO:HD3	1.92	0.42
8:EM:69:VAL:O	8:EM:70:MET:HE2	2.20	0.42
8:EM:274:THR:HG23	8:EM:275:ASP:OD1	2.19	0.42
8:EM:275:ASP:HB3	8:EM:294:PHE:HB2	2.01	0.42
1:FC:114:ARG:HG2	1:FC:130:ARG:HG2	2.02	0.42
5:FH:267:PRO:HB2	5:WH:328:MET:HE2	2.01	0.42
8:FM:257:ASP:OD1	8:FM:257:ASP:N	2.52	0.42
8:GM:76:ARG:HH11	8:GM:76:ARG:HD3	1.63	0.42
8:HM:49:GLN:HG3	8:HM:115:ARG:HB3	2.02	0.42
8:HM:127:ASN:HB3	8:HM:147:ALA:HB3	2.01	0.42
8:HM:275:ASP:HB3	8:HM:294:PHE:HB2	2.01	0.42
1:IC:69:GLN:HE22	9:IN:114:LEU:HA	1.84	0.42
2:ID:87:TRP:CG	2:ID:94:LEU:HD12	2.54	0.42
5:IH:195:ASP:HB2	4:JG:820:TYR:HD1	1.85	0.42
2:JD:123:THR:OG1	2:JD:124:LYS:N	2.53	0.42
8:JM:49:GLN:HG3	8:JM:115:ARG:HB3	2.02	0.42
1:KC:210:ILE:HA	1:KC:210:ILE:HD12	1.65	0.42
8:KM:71:LEU:HD22	8:KM:81:VAL:HG11	2.01	0.42
2:MD:57:LYS:HE2	2:MD:57:LYS:HB3	1.80	0.42
7:BL:91:VAL:HA	7:BL:94:VAL:HG22	2.02	0.42
5:CH:169:ILE:HD11	5:CH:239:LEU:HD23	2.02	0.42
2:Cd:73:ILE:HA	2:Cd:74:PRO:HD3	1.76	0.42
3:Cf:238:ILE:HA	3:Cf:239:PRO:HD3	1.95	0.42
3:Cf:247:ILE:HA	3:Cf:254:ILE:HG22	2.01	0.42
5:DH:179:LEU:O	5:DH:212:THR:HA	2.20	0.42
6:FK:65:ILE:HD12	6:FK:78:ILE:HG23	2.01	0.42
6:GK:144:LEU:HD23	6:GK:144:LEU:HA	1.85	0.42
8:GM:110:THR:OG1	8:GM:111:LYS:N	2.47	0.42
8:HM:69:VAL:O	8:HM:70:MET:HE2	2.20	0.42
8:HM:250:ARG:HD2	8:HM:268:HIS:HB2	2.00	0.42
1:IC:114:ARG:HG2	1:IC:130:ARG:HG2	2.02	0.42
5:IH:194:TYR:CE1	4:JG:821:THR:HB	2.55	0.42
5:IH:284:LEU:HD22	5:IH:330:SER:HB3	2.02	0.42
8:IM:69:VAL:O	8:IM:70:MET:HE2	2.20	0.42
8:IM:274:THR:HG23	8:IM:275:ASP:OD1	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:JK:65:ILE:HD12	6:JK:78:ILE:HG23	2.01	0.42
8:JM:257:ASP:N	8:JM:257:ASP:OD1	2.52	0.42
1:KC:228:HIS:HB3	1:KC:247:LEU:HG	2.00	0.42
3:KF:265:SER:O	3:KF:269:SER:HB3	2.19	0.42
7:KL:91:VAL:HA	7:KL:94:VAL:HG22	2.02	0.42
1:LC:114:ARG:HG2	1:LC:130:ARG:HG2	2.02	0.42
4:MG:808:LEU:HD13	4:MG:808:LEU:HA	1.87	0.42
9:AN:62:ASN:OD1	9:AN:62:ASN:N	2.51	0.42
2:Ad:150:ASN:OD1	2:Ad:151:SER:N	2.47	0.42
3:Af:243:MET:HB3	3:Af:257:SER:HB3	2.01	0.42
5:BH:200:SER:N	4:CG:816:GLU:OE2	2.51	0.42
5:CH:289:PRO:O	5:CH:292:SER:OG	2.37	0.42
5:DH:195:ASP:HB3	5:DH:227:ALA:HB3	2.02	0.42
9:DN:76:MET:H	9:DN:76:MET:HG3	1.73	0.42
7:FL:91:VAL:HA	7:FL:94:VAL:HG22	2.02	0.42
8:FM:69:VAL:O	8:FM:70:MET:HE2	2.20	0.42
1:GC:228:HIS:HB3	1:GC:247:LEU:HG	2.01	0.42
8:GM:61:HIS:HB2	8:GM:109:LYS:HG2	2.00	0.42
8:GM:69:VAL:O	8:GM:70:MET:HE2	2.20	0.42
2:Gd:132:LEU:HD23	2:Gd:132:LEU:HA	1.89	0.42
3:HF:269:SER:O	3:HF:269:SER:OG	2.37	0.42
7:HL:98:TYR:HE1	7:HL:104:LYS:HE3	1.83	0.42
8:HM:84:ILE:HG22	8:HM:91:TYR:HB2	2.00	0.42
9:HN:81:LEU:HD21	5:IH:357:LYS:HB2	2.02	0.42
1:IC:213:ARG:HA	1:IC:213:ARG:HD2	1.83	0.42
9:IN:81:LEU:HD21	5:JH:357:LYS:HB2	2.01	0.42
8:JM:127:ASN:HB3	8:JM:147:ALA:HB3	2.01	0.42
8:KM:284:LEU:HD12	8:KM:284:LEU:HA	1.85	0.42
2:Ld:131:ILE:O	2:Ld:135:ILE:HG12	2.20	0.42
1:MC:114:ARG:HG2	1:MC:130:ARG:HG2	2.02	0.42
3:ZF:267:GLU:O	5:ZH:150:LYS:NZ	2.36	0.42
1:AC:114:ARG:HG2	1:AC:130:ARG:HG2	2.02	0.41
5:AH:191:ILE:HG13	5:AH:211:ASN:HA	2.01	0.41
2:BD:155:GLU:OE2	2:Bd:133:ARG:NH2	2.49	0.41
5:BH:122:LEU:HD13	5:BH:130:LEU:HD21	2.01	0.41
2:Cd:83:ALA:HA	1:DC:267:ALA:O	2.20	0.41
2:DD:57:LYS:HB3	2:DD:57:LYS:HE2	1.85	0.41
8:DM:49:GLN:HG3	8:DM:115:ARG:HB3	2.02	0.41
8:DM:274:THR:HG23	8:DM:275:ASP:OD1	2.20	0.41
8:EM:61:HIS:HB2	8:EM:109:LYS:HG2	2.00	0.41
3:FF:254:ILE:O	3:FF:261:VAL:HA	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:FH:307:SER:OG	5:FH:308:ASN:N	2.53	0.41
8:FM:49:GLN:HG3	8:FM:115:ARG:HB3	2.02	0.41
8:GM:127:ASN:HB3	8:GM:147:ALA:HB3	2.01	0.41
5:HH:183:ASP:HB3	5:HH:189:TRP:HE3	1.84	0.41
5:IH:150:LYS:CD	3:XF:219:ARG:HH22	2.32	0.41
8:IM:247:ARG:HD3	3:If:213:GLN:O	2.20	0.41
1:JC:58:ALA:O	1:JC:62:THR:HG23	2.19	0.41
1:JC:215:LEU:HD12	1:JC:220:MET:HB2	2.02	0.41
7:JL:91:VAL:HA	7:JL:94:VAL:HG22	2.02	0.41
2:Jd:79:LEU:O	2:Jd:127:SER:OG	2.38	0.41
3:LF:254:ILE:O	3:LF:261:VAL:HA	2.20	0.41
1:MC:116:THR:HG23	5:MH:328:MET:HG2	2.02	0.41
8:MM:274:THR:HG23	8:MM:275:ASP:OD1	2.19	0.41
2:Md:97:ARG:HA	2:Md:97:ARG:HD3	1.83	0.41
5:AH:205:GLN:HG2	5:AH:206:TRP:N	2.29	0.41
8:AM:257:ASP:N	8:AM:257:ASP:OD1	2.52	0.41
8:AM:274:THR:HG23	8:AM:275:ASP:OD1	2.20	0.41
1:BC:215:LEU:HD12	1:BC:220:MET:HB2	2.02	0.41
5:BH:238:MET:HE2	5:BH:238:MET:HB2	1.73	0.41
2:Bd:26:LYS:HD3	2:Bd:26:LYS:HA	1.90	0.41
1:EC:154:ASP:HB3	9:EN:84:CYS:HB3	2.02	0.41
3:EF:254:ILE:O	3:EF:261:VAL:HA	2.20	0.41
1:FC:215:LEU:HD12	1:FC:220:MET:HB2	2.02	0.41
2:FD:136:ASP:OD1	2:FD:145:ILE:N	2.51	0.41
8:HM:71:LEU:HD22	8:HM:81:VAL:HG11	2.01	0.41
5:JH:208:LYS:HD2	4:KG:823:GLY:CA	2.48	0.41
7:JL:98:TYR:HE1	7:JL:104:LYS:HE3	1.83	0.41
8:JM:71:LEU:HD22	8:JM:81:VAL:HG11	2.02	0.41
1:KC:215:LEU:HD12	1:KC:220:MET:HB2	2.02	0.41
6:KK:65:ILE:HD12	6:KK:78:ILE:HG23	2.01	0.41
1:LC:207:GLY:HA2	1:LC:210:ILE:HG22	2.03	0.41
7:LL:111:LYS:HE3	7:LL:111:LYS:HB2	1.86	0.41
7:ML:91:VAL:HA	7:ML:94:VAL:HG22	2.02	0.41
3:WF:254:ILE:O	3:WF:261:VAL:HA	2.20	0.41
5:WH:320:LEU:HD13	5:WH:348:SER:H	1.85	0.41
3:YF:254:ILE:O	3:YF:261:VAL:HA	2.20	0.41
1:AC:215:LEU:HD12	1:AC:220:MET:HB2	2.02	0.41
8:BM:127:ASN:HB3	8:BM:147:ALA:HB3	2.01	0.41
2:Cd:131:ILE:O	2:Cd:135:ILE:HG12	2.20	0.41
1:EC:215:LEU:HD12	1:EC:220:MET:HB2	2.02	0.41
1:EC:259:SER:HA	1:EC:262:TRP:CE2	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:EM:71:LEU:HD22	8:EM:81:VAL:HG11	2.01	0.41
1:FC:259:SER:HA	1:FC:262:TRP:CE2	2.55	0.41
5:FH:319:ILE:HG21	5:FH:339:MET:HE1	2.02	0.41
2:Fd:131:ILE:O	2:Fd:135:ILE:HG12	2.20	0.41
5:GH:208:LYS:HE3	5:GH:208:LYS:HB3	1.79	0.41
5:HH:199:PRO:HG2	4:XG:819:VAL:HG13	2.01	0.41
8:HM:274:THR:HG23	8:HM:275:ASP:OD1	2.19	0.41
1:KC:114:ARG:HG2	1:KC:130:ARG:HG2	2.02	0.41
8:LM:49:GLN:HG3	8:LM:115:ARG:HB3	2.01	0.41
1:MC:259:SER:HA	1:MC:262:TRP:CE2	2.55	0.41
2:MD:140:GLY:O	2:Md:60:LYS:NZ	2.53	0.41
3:MF:254:ILE:O	3:MF:261:VAL:HA	2.20	0.41
3:XF:212:ILE:HD11	3:XF:254:ILE:HG13	2.03	0.41
3:ZF:214:ALA:CB	5:ZH:242:ILE:HD11	2.51	0.41
1:AC:207:GLY:HA2	1:AC:210:ILE:HG22	2.03	0.41
1:AC:259:SER:HA	1:AC:262:TRP:CE2	2.56	0.41
5:AH:284:LEU:HD22	5:AH:330:SER:HB3	2.01	0.41
8:AM:71:LEU:HD22	8:AM:81:VAL:HG11	2.01	0.41
3:BF:212:ILE:HD11	3:BF:254:ILE:HG13	2.03	0.41
8:BM:69:VAL:O	8:BM:70:MET:HE2	2.20	0.41
8:BM:76:ARG:HH11	8:BM:76:ARG:HD3	1.63	0.41
6:CK:65:ILE:HD12	6:CK:78:ILE:HG23	2.01	0.41
8:CM:71:LEU:HD22	8:CM:81:VAL:HG11	2.01	0.41
7:DL:111:LYS:HB2	7:DL:111:LYS:HE3	1.86	0.41
3:EF:212:ILE:HD11	3:EF:254:ILE:HG13	2.03	0.41
4:EG:800:ASP:OD1	4:EG:800:ASP:N	2.46	0.41
5:EH:130:LEU:HA	5:EH:133:ILE:HG12	2.02	0.41
8:EM:127:ASN:HB3	8:EM:147:ALA:HB3	2.01	0.41
1:FC:180:THR:CG2	1:GC:34:LEU:HD23	2.50	0.41
8:FM:90:LEU:HG	8:FM:92:VAL:HG23	2.03	0.41
2:Fd:33:PRO:HB2	2:Fd:39:ILE:HD13	2.01	0.41
7:GL:98:TYR:HE1	7:GL:104:LYS:HE3	1.83	0.41
8:GM:274:THR:HG23	8:GM:275:ASP:OD1	2.20	0.41
8:HM:257:ASP:N	8:HM:257:ASP:OD1	2.52	0.41
2:ID:155:GLU:OE2	2:ID:157:ARG:NH2	2.53	0.41
3:JF:254:ILE:O	3:JF:261:VAL:HA	2.20	0.41
6:JK:63:LYS:HB2	6:JK:63:LYS:HE2	1.84	0.41
4:KG:800:ASP:N	4:KG:800:ASP:OD1	2.46	0.41
5:LH:251:ARG:NH1	5:LH:253:ASP:OD2	2.53	0.41
8:LM:69:VAL:O	8:LM:70:MET:HE2	2.20	0.41
1:MC:88:ARG:HA	1:MC:88:ARG:HD2	1.82	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:ML:190:MET:HE3	7:ML:190:MET:HB3	1.82	0.41
3:VF:212:ILE:HD11	3:VF:254:ILE:HG13	2.03	0.41
5:XH:330:SER:OG	5:XH:332:ASP:OD1	2.30	0.41
3:YF:212:ILE:HD11	3:YF:254:ILE:HG13	2.03	0.41
2:AD:79:LEU:HD23	2:AD:79:LEU:HA	1.84	0.41
4:AG:812:TRP:HZ3	5:BH:130:LEU:HD12	1.86	0.41
8:AM:69:VAL:O	8:AM:70:MET:HE2	2.20	0.41
8:BM:90:LEU:HG	8:BM:92:VAL:HG23	2.03	0.41
8:BM:275:ASP:HB3	8:BM:294:PHE:HB2	2.01	0.41
1:DC:114:ARG:HG2	1:DC:130:ARG:HG2	2.02	0.41
5:EH:115:MET:CE	4:VG:801:MET:HB3	2.51	0.41
5:EH:307:SER:OG	5:EH:308:ASN:N	2.53	0.41
9:EN:76:MET:H	9:EN:76:MET:HG3	1.73	0.41
3:GF:254:ILE:O	3:GF:261:VAL:HA	2.20	0.41
1:HC:259:SER:HA	1:HC:262:TRP:CE2	2.56	0.41
3:HF:254:ILE:O	3:HF:261:VAL:HA	2.20	0.41
8:HM:217:LYS:HA	8:HM:217:LYS:HD3	1.80	0.41
1:IC:259:SER:HA	1:IC:262:TRP:CE2	2.55	0.41
1:JC:259:SER:HA	1:JC:262:TRP:CE2	2.56	0.41
4:JG:815:VAL:HG21	5:JH:173:GLN:HE21	1.86	0.41
6:JK:144:LEU:HD23	6:JK:144:LEU:HA	1.85	0.41
8:KM:127:ASN:HB3	8:KM:147:ALA:HB3	2.01	0.41
8:KM:274:THR:HG23	8:KM:275:ASP:OD1	2.19	0.41
8:LM:306:GLN:OE1	3:Lf:251:GLN:HG2	2.20	0.41
2:MD:71:LEU:HD11	2:Md:62:ILE:HD12	2.01	0.41
7:ML:47:TYR:OH	7:ML:142:GLU:OE1	2.28	0.41
9:MN:76:MET:H	9:MN:76:MET:HG3	1.73	0.41
3:ZF:254:ILE:O	3:ZF:261:VAL:HA	2.20	0.41
5:ZH:190:PRO:HB2	5:ZH:231:ARG:HG3	2.02	0.41
3:AF:212:ILE:HD11	3:AF:254:ILE:HG13	2.03	0.41
2:BD:136:ASP:OD1	2:Bd:60:LYS:HE2	2.19	0.41
8:BM:274:THR:HG23	8:BM:275:ASP:OD1	2.19	0.41
1:CC:114:ARG:HG2	1:CC:130:ARG:HG2	2.02	0.41
1:DC:207:GLY:HA2	1:DC:210:ILE:HG22	2.03	0.41
5:EH:173:GLN:NE2	5:EH:218:THR:O	2.43	0.41
5:EH:273:PRO:HA	5:EH:274:PRO:HD3	1.92	0.41
8:EM:49:GLN:HG3	8:EM:115:ARG:HB3	2.02	0.41
8:EM:310:LYS:HB3	8:EM:310:LYS:HE3	1.88	0.41
5:FH:289:PRO:O	5:FH:292:SER:OG	2.35	0.41
8:FM:274:THR:HG23	8:FM:275:ASP:OD1	2.19	0.41
8:HM:90:LEU:HG	8:HM:92:VAL:HG23	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:ID:107:ARG:CZ	2:ID:153:VAL:HG11	2.51	0.41
8:IM:71:LEU:HD22	8:IM:81:VAL:HG11	2.01	0.41
2:Id:157:ARG:HE	2:Id:157:ARG:HB2	1.69	0.41
3:JF:212:ILE:HD11	3:JF:254:ILE:HG13	2.03	0.41
7:JL:111:LYS:HE3	7:JL:111:LYS:HB2	1.86	0.41
8:JM:274:THR:HG23	8:JM:275:ASP:OD1	2.20	0.41
1:KC:259:SER:HA	1:KC:262:TRP:CE2	2.56	0.41
3:KF:254:ILE:O	3:KF:261:VAL:HA	2.20	0.41
8:KM:69:VAL:O	8:KM:70:MET:HE2	2.20	0.41
5:ZH:242:ILE:HG21	5:ZH:242:ILE:HD13	1.75	0.41
3:BF:254:ILE:O	3:BF:261:VAL:HA	2.20	0.41
5:BH:133:ILE:HD12	5:BH:133:ILE:HA	1.82	0.41
5:BH:220:LEU:HB3	5:BH:221:TYR:CD1	2.55	0.41
5:BH:273:PRO:HA	5:BH:274:PRO:HD3	1.94	0.41
7:BL:98:TYR:HE1	7:BL:104:LYS:HE3	1.83	0.41
7:BL:193:LEU:HD23	7:BL:193:LEU:HA	1.91	0.41
8:CM:49:GLN:HG3	8:CM:115:ARG:HB3	2.02	0.41
8:CM:69:VAL:O	8:CM:70:MET:HE2	2.20	0.41
8:CM:90:LEU:HG	8:CM:92:VAL:HG23	2.03	0.41
8:CM:274:THR:HG23	8:CM:275:ASP:OD1	2.20	0.41
2:Cd:97:ARG:HD3	2:Cd:97:ARG:HA	1.83	0.41
7:DL:193:LEU:HD23	7:DL:193:LEU:HA	1.91	0.41
8:DM:54:GLY:O	8:DM:118:TYR:OH	2.33	0.41
6:EK:66:LYS:HE2	7:FL:118:GLU:O	2.20	0.41
1:GC:114:ARG:HG2	1:GC:130:ARG:HG2	2.02	0.41
1:GC:215:LEU:HD12	1:GC:220:MET:HB2	2.02	0.41
3:HF:212:ILE:HD11	3:HF:254:ILE:HG13	2.03	0.41
7:HL:91:VAL:HA	7:HL:94:VAL:HG22	2.02	0.41
2:Hd:73:ILE:HA	2:Hd:74:PRO:HD3	1.86	0.41
1:IC:207:GLY:HA2	1:IC:210:ILE:HG22	2.03	0.41
8:IM:70:MET:HB2	8:IM:107:ASP:HB3	2.03	0.41
8:LM:274:THR:HG23	8:LM:275:ASP:OD1	2.20	0.41
5:MH:208:LYS:NZ	4:ZG:822:GLU:O	2.51	0.41
8:MM:69:VAL:O	8:MM:70:MET:HE2	2.20	0.41
8:MM:71:LEU:HD22	8:MM:81:VAL:HG11	2.01	0.41
5:ZH:166:PRO:HA	5:ZH:167:PRO:HD3	1.92	0.41
1:AC:154:ASP:HB3	9:AN:84:CYS:HB3	2.01	0.41
2:AD:87:TRP:CD1	2:AD:89:GLY:H	2.39	0.41
3:AF:254:ILE:O	3:AF:261:VAL:HA	2.21	0.41
8:BM:49:GLN:HG3	8:BM:115:ARG:HB3	2.02	0.41
2:Bd:73:ILE:HA	2:Bd:74:PRO:HD3	1.84	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CC:215:LEU:HD12	1:CC:220:MET:HB2	2.02	0.41
1:DC:213:ARG:HD2	1:DC:213:ARG:HA	1.83	0.41
3:DF:254:ILE:O	3:DF:261:VAL:HA	2.20	0.41
8:DM:257:ASP:OD1	8:DM:257:ASP:N	2.52	0.41
5:EH:264:LYS:HA	5:EH:264:LYS:HD3	1.78	0.41
1:GC:93:ILE:HA	2:Gd:48:VAL:HG11	2.02	0.41
5:GH:184:SER:OG	5:GH:257:GLN:O	2.34	0.41
9:GN:58:LEU:HD12	9:GN:68:CYS:HB2	2.03	0.41
7:IL:44:HIS:HA	7:IL:45:PRO:HD3	1.95	0.41
7:IL:111:LYS:HE3	7:IL:111:LYS:HB2	1.86	0.41
1:JC:114:ARG:HG2	1:JC:130:ARG:HG2	2.02	0.41
2:JD:161:ILE:HD12	2:JD:161:ILE:HA	1.88	0.41
7:JL:94:VAL:HA	7:JL:97:ILE:HG12	2.03	0.41
8:JM:69:VAL:O	8:JM:70:MET:HE2	2.20	0.41
1:KC:207:GLY:HA2	1:KC:210:ILE:HG22	2.03	0.41
8:KM:49:GLN:HG3	8:KM:115:ARG:HB3	2.02	0.41
8:LM:71:LEU:HD22	8:LM:81:VAL:HG11	2.01	0.41
2:Ld:73:ILE:HA	2:Ld:74:PRO:HD3	1.82	0.41
5:XH:166:PRO:HA	5:XH:167:PRO:HD3	1.90	0.41
5:ZH:112:PHE:O	5:ZH:115:MET:HB3	2.20	0.41
5:AH:254:LEU:HD13	5:AH:254:LEU:HA	1.86	0.41
6:BK:63:LYS:HB2	6:BK:63:LYS:HE2	1.84	0.41
7:BL:94:VAL:HA	7:BL:97:ILE:HG12	2.03	0.41
8:BM:110:THR:OG1	8:BM:111:LYS:N	2.47	0.41
1:CC:157:LYS:CE	1:DC:28:THR:OG1	2.68	0.41
5:CH:238:MET:HB3	5:CH:238:MET:HE2	1.42	0.41
7:CL:94:VAL:HA	7:CL:97:ILE:HG12	2.03	0.41
8:CM:70:MET:HB2	8:CM:107:ASP:HB3	2.03	0.41
8:DM:71:LEU:HD22	8:DM:81:VAL:HG11	2.01	0.41
1:EC:207:GLY:HA2	1:EC:210:ILE:HG22	2.03	0.41
4:EG:808:LEU:HA	4:EG:808:LEU:HD13	1.88	0.41
7:EL:91:VAL:HA	7:EL:94:VAL:HG22	2.02	0.41
8:EM:70:MET:HB2	8:EM:107:ASP:HB3	2.03	0.41
8:EM:90:LEU:HG	8:EM:92:VAL:HG23	2.03	0.41
2:Ed:105:ARG:HB2	2:Ed:153:VAL:HG23	2.03	0.41
1:FC:207:GLY:HA2	1:FC:210:ILE:HG22	2.03	0.41
3:FF:212:ILE:HD11	3:FF:254:ILE:HG13	2.03	0.41
7:FL:94:VAL:HA	7:FL:97:ILE:HG12	2.03	0.41
8:FM:70:MET:HB2	8:FM:107:ASP:HB3	2.03	0.41
1:GC:259:SER:HA	1:GC:262:TRP:CE2	2.56	0.41
5:GH:113:LYS:HB3	5:GH:113:LYS:HE3	1.88	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:GM:310:LYS:HB3	8:GM:310:LYS:HE3	1.88	0.41
1:HC:207:GLY:HA2	1:HC:210:ILE:HG22	2.03	0.41
7:HL:125:LEU:HD23	7:HL:125:LEU:HA	1.87	0.41
1:IC:215:LEU:HD12	1:IC:220:MET:HB2	2.02	0.41
5:IH:169:ILE:HB	5:IH:241:LEU:HD23	2.02	0.41
7:IL:125:LEU:HD23	7:IL:125:LEU:HA	1.87	0.41
8:IM:49:GLN:HG3	8:IM:115:ARG:HB3	2.02	0.41
8:IM:217:LYS:HA	8:IM:217:LYS:HD3	1.81	0.41
8:IM:307:SER:HG	3:If:253:ARG:HH22	1.67	0.41
9:JN:58:LEU:HD12	9:JN:68:CYS:HB2	2.03	0.41
3:KF:212:ILE:HD11	3:KF:254:ILE:HG13	2.03	0.41
8:KM:275:ASP:HB3	8:KM:294:PHE:HB2	2.01	0.41
9:KN:58:LEU:HD12	9:KN:68:CYS:HB2	2.03	0.41
1:LC:259:SER:HA	1:LC:262:TRP:CE2	2.56	0.41
1:MC:215:LEU:HD12	1:MC:220:MET:HB2	2.02	0.41
6:MK:65:ILE:HD12	6:MK:78:ILE:HG23	2.01	0.41
7:ML:44:HIS:HA	7:ML:45:PRO:HD3	1.95	0.41
7:ML:125:LEU:HD23	7:ML:125:LEU:HA	1.87	0.41
8:MM:65:ASN:HD21	8:MM:109:LYS:HD3	1.86	0.41
8:MM:90:LEU:HG	8:MM:92:VAL:HG23	2.03	0.41
3:VF:254:ILE:O	3:VF:261:VAL:HA	2.20	0.41
4:WG:808:LEU:HA	4:WG:808:LEU:HD13	1.87	0.41
3:ZF:212:ILE:HD11	3:ZF:254:ILE:HG13	2.03	0.41
1:AC:177:CYS:HB2	1:BC:35:GLN:OE1	2.20	0.41
1:BC:207:GLY:HA2	1:BC:210:ILE:HG22	2.03	0.41
1:BC:259:SER:HA	1:BC:262:TRP:CE2	2.55	0.41
5:BH:107:ILE:HD12	4:ZG:794:ILE:CD1	2.50	0.41
2:Bd:84:SER:OG	2:CD:57:LYS:NZ	2.50	0.41
2:Bd:124:LYS:HB3	2:Bd:124:LYS:HE2	1.83	0.41
3:CF:254:ILE:O	3:CF:261:VAL:HA	2.20	0.41
5:CH:190:PRO:O	5:CH:231:ARG:HG2	2.21	0.41
5:CH:208:LYS:HA	5:CH:208:LYS:HD2	1.92	0.41
8:CM:56:LEU:HD13	8:CM:58:VAL:HG23	2.03	0.41
1:DC:215:LEU:HD12	1:DC:220:MET:HB2	2.02	0.41
1:EC:88:ARG:HA	1:EC:88:ARG:HD2	1.82	0.41
4:EG:801:MET:C	5:WH:115:MET:HE1	2.46	0.41
1:GC:207:GLY:HA2	1:GC:210:ILE:HG22	2.03	0.41
8:IM:90:LEU:HG	8:IM:92:VAL:HG23	2.03	0.41
2:Id:126:GLU:OE2	2:Id:130:GLU:HB2	2.21	0.41
8:KM:218:LYS:HA	8:KM:218:LYS:HD3	1.95	0.41
5:LH:109:LYS:O	5:LH:113:LYS:HG3	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:MH:143:ALA:O	5:ZH:131:LYS:NZ	2.37	0.41
8:MM:54:GLY:O	8:MM:118:TYR:OH	2.33	0.41
8:MM:70:MET:HB2	8:MM:107:ASP:HB3	2.03	0.41
5:ZH:126:GLN:O	5:ZH:130:LEU:HD23	2.21	0.41
9:AN:46:GLY:HA3	9:AN:61:ASN:HB2	2.04	0.40
2:BD:24:LYS:O	7:BL:110:GLN:NE2	2.54	0.40
7:BL:125:LEU:HD23	7:BL:125:LEU:HA	1.87	0.40
1:CC:207:GLY:HA2	1:CC:210:ILE:HG22	2.03	0.40
8:CM:65:ASN:HD21	8:CM:109:LYS:HD3	1.86	0.40
1:DC:83:LEU:HD23	1:DC:83:LEU:HA	1.95	0.40
7:DL:91:VAL:HA	7:DL:94:VAL:HG22	2.02	0.40
8:DM:56:LEU:HD13	8:DM:58:VAL:HG23	2.04	0.40
5:EH:156:GLN:HB3	5:EH:254:LEU:HD13	2.03	0.40
1:GC:240:ILE:HG12	1:HC:255:LEU:HD13	2.03	0.40
3:GF:212:ILE:HD11	3:GF:254:ILE:HG13	2.03	0.40
5:IH:307:SER:OG	5:IH:308:ASN:N	2.54	0.40
8:IM:319:PHE:HA	8:IM:320:PRO:HD2	1.82	0.40
7:JL:125:LEU:HD23	7:JL:125:LEU:HA	1.87	0.40
8:JM:70:MET:HB2	8:JM:107:ASP:HB3	2.03	0.40
8:JM:90:LEU:HG	8:JM:92:VAL:HG23	2.03	0.40
2:KD:79:LEU:HA	2:KD:79:LEU:HD23	1.81	0.40
7:KL:125:LEU:HD23	7:KL:125:LEU:HA	1.87	0.40
3:LF:212:ILE:HD11	3:LF:254:ILE:HG13	2.03	0.40
8:LM:90:LEU:HG	8:LM:92:VAL:HG23	2.03	0.40
9:MN:58:LEU:HD12	9:MN:68:CYS:HB2	2.03	0.40
4:WG:822:GLU:OE2	5:WH:207:ASP:OD2	2.39	0.40
1:AC:213:ARG:HA	1:AC:213:ARG:HD2	1.83	0.40
5:AH:208:LYS:CD	4:BG:823:GLY:HA2	2.52	0.40
1:BC:35:GLN:HE21	1:BC:35:GLN:CA	2.26	0.40
5:BH:233:LEU:HD23	5:BH:233:LEU:HA	1.94	0.40
8:BM:54:GLY:O	8:BM:118:TYR:OH	2.33	0.40
7:CL:91:VAL:HA	7:CL:94:VAL:HG22	2.02	0.40
4:DG:797:ARG:NE	5:EH:117:ARG:HH22	2.19	0.40
5:DH:219:LYS:HA	5:DH:219:LYS:HD2	1.90	0.40
8:DM:271:SER:OG	8:DM:290:ASP:OD1	2.40	0.40
2:ED:79:LEU:HD23	2:ED:79:LEU:HA	1.83	0.40
5:EH:144:THR:HA	5:EH:145:PRO:HD3	1.94	0.40
8:EM:56:LEU:HD13	8:EM:58:VAL:HG23	2.04	0.40
8:FM:76:ARG:HH11	8:FM:76:ARG:HD3	1.63	0.40
2:GD:136:ASP:OD1	2:Gd:60:LYS:HE2	2.21	0.40
3:GF:219:ARG:HE	3:GF:219:ARG:HB3	1.75	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:HD:77:TYR:OH	6:HK:130:ARG:NH1	2.52	0.40
5:HH:307:SER:OG	5:HH:308:ASN:N	2.52	0.40
3:IF:212:ILE:HD11	3:IF:254:ILE:HG13	2.03	0.40
8:IM:56:LEU:HD13	8:IM:58:VAL:HG23	2.04	0.40
8:IM:257:ASP:N	8:IM:257:ASP:OD1	2.52	0.40
1:JC:213:ARG:HD2	1:JC:213:ARG:HA	1.83	0.40
5:JH:354:MET:HB2	5:JH:354:MET:HE3	1.72	0.40
2:KD:52:MET:HE2	2:Kd:52:MET:HE1	2.03	0.40
7:LL:91:VAL:HA	7:LL:94:VAL:HG22	2.02	0.40
2:Ld:97:ARG:HA	2:Ld:97:ARG:HD3	1.78	0.40
8:MM:301:MET:HE3	8:MM:301:MET:HB3	2.00	0.40
5:WH:354:MET:HE2	5:WH:354:MET:HB2	1.91	0.40
8:AM:319:PHE:HA	8:AM:320:PRO:HD2	1.82	0.40
9:AN:58:LEU:HD12	9:AN:68:CYS:HB2	2.03	0.40
3:BF:219:ARG:NH2	5:CH:151:PRO:HD2	2.36	0.40
9:BN:46:GLY:HA3	9:BN:61:ASN:HB2	2.04	0.40
5:CH:208:LYS:NZ	4:VG:822:GLU:O	2.41	0.40
4:DG:818:GLN:HG3	5:DH:205:GLN:NE2	2.35	0.40
8:EM:65:ASN:HD21	8:EM:109:LYS:HD3	1.86	0.40
9:FN:58:LEU:HD12	9:FN:68:CYS:HB2	2.03	0.40
5:GH:149:PRO:HB2	5:GH:248:VAL:HG12	2.03	0.40
7:GL:94:VAL:HA	7:GL:97:ILE:HG12	2.03	0.40
8:HM:218:LYS:HA	8:HM:218:LYS:HD3	1.95	0.40
9:HN:58:LEU:HD12	9:HN:68:CYS:HB2	2.03	0.40
2:ID:55:MET:HB3	2:ID:55:MET:HE3	1.91	0.40
5:IH:289:PRO:O	5:IH:292:SER:OG	2.35	0.40
5:JH:191:ILE:HG23	5:JH:228:VAL:HG13	2.03	0.40
5:LH:289:PRO:O	5:LH:292:SER:OG	2.35	0.40
7:LL:94:VAL:HA	7:LL:97:ILE:HG12	2.03	0.40
6:MK:138:ARG:HH11	6:MK:138:ARG:HD3	1.75	0.40
3:WF:212:ILE:HD11	3:WF:254:ILE:HG13	2.03	0.40
5:YH:348:SER:HA	5:YH:352:LYS:O	2.21	0.40
5:AH:311:MET:HE3	5:AH:311:MET:HB2	1.81	0.40
7:AL:91:VAL:HA	7:AL:94:VAL:HG22	2.02	0.40
3:Af:238:ILE:HG23	3:Af:241:TYR:HB2	2.03	0.40
3:Af:254:ILE:O	3:Af:261:VAL:HA	2.22	0.40
1:DC:259:SER:HA	1:DC:262:TRP:CE2	2.56	0.40
5:DH:280:LEU:HB2	5:DH:328:MET:HE2	2.04	0.40
8:FM:56:LEU:HD13	8:FM:58:VAL:HG23	2.04	0.40
5:GH:204:ILE:HG12	5:GH:215:ILE:HD12	2.02	0.40
7:GL:193:LEU:HD23	7:GL:193:LEU:HA	1.91	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:HC:215:LEU:HD12	1:HC:220:MET:HB2	2.02	0.40
4:HG:818:GLN:H	5:HH:216:GLN:HE22	1.70	0.40
5:HH:115:MET:SD	4:WG:801:MET:HB3	2.60	0.40
5:HH:184:SER:OG	5:HH:257:GLN:O	2.36	0.40
3:IF:254:ILE:O	3:IF:261:VAL:HA	2.20	0.40
5:IH:133:ILE:HD12	5:IH:133:ILE:HA	1.95	0.40
1:LC:215:LEU:HD12	1:LC:220:MET:HB2	2.02	0.40
2:Ld:133:ARG:HH11	2:Ld:133:ARG:HD3	1.74	0.40
2:MD:87:TRP:CG	2:MD:94:LEU:HD12	2.56	0.40
5:MH:169:ILE:HB	5:MH:241:LEU:HD23	2.03	0.40
7:ML:111:LYS:HE3	7:ML:111:LYS:HB2	1.86	0.40
8:AM:76:ARG:HH11	8:AM:76:ARG:HD3	1.63	0.40
5:BH:254:LEU:HD23	5:BH:254:LEU:HA	1.86	0.40
8:BM:56:LEU:HD13	8:BM:58:VAL:HG23	2.04	0.40
2:Cd:123:THR:HG21	2:Cd:126:GLU:HB2	2.02	0.40
9:DN:58:LEU:HD12	9:DN:68:CYS:HB2	2.03	0.40
1:FC:204:GLU:HA	5:FH:285:GLU:OE2	2.21	0.40
9:FN:46:GLY:HA3	9:FN:61:ASN:HB2	2.04	0.40
5:GH:170:ARG:HG2	5:GH:245:GLN:HG3	2.03	0.40
1:HC:213:ARG:HD2	1:HC:213:ARG:HA	1.83	0.40
8:HM:56:LEU:HD13	8:HM:58:VAL:HG23	2.04	0.40
5:IH:189:TRP:CZ2	5:IH:260:GLY:HA2	2.57	0.40
7:IL:91:VAL:HA	7:IL:94:VAL:HG22	2.02	0.40
4:KG:801:MET:HB3	5:LH:115:MET:CE	2.51	0.40
9:KN:81:LEU:HD21	5:LH:357:LYS:HB2	2.03	0.40
1:LC:177:CYS:HA	1:MC:35:GLN:OE1	2.21	0.40
8:LM:70:MET:HB2	8:LM:107:ASP:HB3	2.03	0.40
2:Md:112:SER:HA	2:Md:113:PRO:HD3	1.96	0.40
3:WF:219:ARG:HE	3:WF:219:ARG:HB3	1.75	0.40
3:XF:254:ILE:O	3:XF:261:VAL:HA	2.20	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	AC	207/303 (68%)	202 (98%)	5 (2%)	0	100	100
1	BC	207/303 (68%)	202 (98%)	5 (2%)	0	100	100
1	CC	207/303 (68%)	202 (98%)	5 (2%)	0	100	100
1	DC	207/303 (68%)	201 (97%)	6 (3%)	0	100	100
1	EC	207/303 (68%)	202 (98%)	5 (2%)	0	100	100
1	FC	207/303 (68%)	202 (98%)	5 (2%)	0	100	100
1	GC	207/303 (68%)	202 (98%)	5 (2%)	0	100	100
1	HC	207/303 (68%)	202 (98%)	5 (2%)	0	100	100
1	IC	207/303 (68%)	202 (98%)	5 (2%)	0	100	100
1	JC	207/303 (68%)	202 (98%)	5 (2%)	0	100	100
1	KC	207/303 (68%)	202 (98%)	5 (2%)	0	100	100
1	LC	207/303 (68%)	202 (98%)	5 (2%)	0	100	100
1	MC	207/303 (68%)	202 (98%)	5 (2%)	0	100	100
2	AD	138/163 (85%)	133 (96%)	5 (4%)	0	100	100
2	Ad	135/163 (83%)	130 (96%)	5 (4%)	0	100	100
2	BD	138/163 (85%)	134 (97%)	4 (3%)	0	100	100
2	Bd	135/163 (83%)	129 (96%)	6 (4%)	0	100	100
2	CD	138/163 (85%)	134 (97%)	4 (3%)	0	100	100
2	Cd	135/163 (83%)	129 (96%)	6 (4%)	0	100	100
2	DD	138/163 (85%)	134 (97%)	4 (3%)	0	100	100
2	Dd	135/163 (83%)	130 (96%)	5 (4%)	0	100	100
2	ED	138/163 (85%)	134 (97%)	4 (3%)	0	100	100
2	Ed	135/163 (83%)	129 (96%)	6 (4%)	0	100	100
2	FD	138/163 (85%)	134 (97%)	4 (3%)	0	100	100
2	Fd	135/163 (83%)	129 (96%)	6 (4%)	0	100	100
2	GD	138/163 (85%)	133 (96%)	5 (4%)	0	100	100
2	Gd	135/163 (83%)	129 (96%)	6 (4%)	0	100	100
2	HD	138/163 (85%)	134 (97%)	4 (3%)	0	100	100
2	Hd	135/163 (83%)	130 (96%)	5 (4%)	0	100	100
2	ID	138/163 (85%)	134 (97%)	4 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	Id	135/163 (83%)	130 (96%)	5 (4%)	0	100	100
2	JD	138/163 (85%)	134 (97%)	4 (3%)	0	100	100
2	Jd	135/163 (83%)	129 (96%)	6 (4%)	0	100	100
2	KD	138/163 (85%)	134 (97%)	4 (3%)	0	100	100
2	Kd	135/163 (83%)	129 (96%)	6 (4%)	0	100	100
2	LD	138/163 (85%)	133 (96%)	5 (4%)	0	100	100
2	Ld	135/163 (83%)	130 (96%)	5 (4%)	0	100	100
2	MD	138/163 (85%)	134 (97%)	4 (3%)	0	100	100
2	Md	135/163 (83%)	129 (96%)	6 (4%)	0	100	100
3	AF	61/269 (23%)	57 (93%)	4 (7%)	0	100	100
3	Af	57/269 (21%)	53 (93%)	4 (7%)	0	100	100
3	BF	61/269 (23%)	57 (93%)	4 (7%)	0	100	100
3	Bf	57/269 (21%)	55 (96%)	2 (4%)	0	100	100
3	CF	61/269 (23%)	57 (93%)	4 (7%)	0	100	100
3	Cf	57/269 (21%)	53 (93%)	4 (7%)	0	100	100
3	DF	61/269 (23%)	57 (93%)	4 (7%)	0	100	100
3	Df	57/269 (21%)	55 (96%)	2 (4%)	0	100	100
3	EF	61/269 (23%)	57 (93%)	4 (7%)	0	100	100
3	Ef	57/269 (21%)	52 (91%)	5 (9%)	0	100	100
3	FF	61/269 (23%)	57 (93%)	4 (7%)	0	100	100
3	Ff	57/269 (21%)	55 (96%)	2 (4%)	0	100	100
3	GF	61/269 (23%)	57 (93%)	4 (7%)	0	100	100
3	Gf	57/269 (21%)	56 (98%)	1 (2%)	0	100	100
3	HF	61/269 (23%)	57 (93%)	4 (7%)	0	100	100
3	Hf	57/269 (21%)	54 (95%)	3 (5%)	0	100	100
3	IF	61/269 (23%)	57 (93%)	4 (7%)	0	100	100
3	If	57/269 (21%)	54 (95%)	3 (5%)	0	100	100
3	JF	61/269 (23%)	57 (93%)	4 (7%)	0	100	100
3	Jf	57/269 (21%)	55 (96%)	2 (4%)	0	100	100
3	KF	61/269 (23%)	57 (93%)	4 (7%)	0	100	100
3	Kf	57/269 (21%)	53 (93%)	4 (7%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	LF	61/269 (23%)	57 (93%)	4 (7%)	0	100	100
3	Lf	57/269 (21%)	55 (96%)	2 (4%)	0	100	100
3	MF	61/269 (23%)	57 (93%)	4 (7%)	0	100	100
3	Mf	57/269 (21%)	54 (95%)	3 (5%)	0	100	100
3	VF	61/269 (23%)	57 (93%)	4 (7%)	0	100	100
3	WF	61/269 (23%)	57 (93%)	4 (7%)	0	100	100
3	XF	61/269 (23%)	57 (93%)	4 (7%)	0	100	100
3	YF	61/269 (23%)	57 (93%)	4 (7%)	0	100	100
3	ZF	61/269 (23%)	57 (93%)	4 (7%)	0	100	100
4	AG	32/1048 (3%)	32 (100%)	0	0	100	100
4	BG	32/1048 (3%)	32 (100%)	0	0	100	100
4	CG	32/1048 (3%)	32 (100%)	0	0	100	100
4	DG	32/1048 (3%)	32 (100%)	0	0	100	100
4	EG	32/1048 (3%)	32 (100%)	0	0	100	100
4	FG	32/1048 (3%)	32 (100%)	0	0	100	100
4	GG	32/1048 (3%)	32 (100%)	0	0	100	100
4	HG	32/1048 (3%)	32 (100%)	0	0	100	100
4	IG	32/1048 (3%)	32 (100%)	0	0	100	100
4	JG	32/1048 (3%)	32 (100%)	0	0	100	100
4	KG	32/1048 (3%)	32 (100%)	0	0	100	100
4	LG	32/1048 (3%)	32 (100%)	0	0	100	100
4	MG	32/1048 (3%)	32 (100%)	0	0	100	100
4	VG	32/1048 (3%)	32 (100%)	0	0	100	100
4	WG	32/1048 (3%)	32 (100%)	0	0	100	100
4	XG	32/1048 (3%)	32 (100%)	0	0	100	100
4	YG	32/1048 (3%)	32 (100%)	0	0	100	100
4	ZG	32/1048 (3%)	32 (100%)	0	0	100	100
5	AH	256/361 (71%)	243 (95%)	13 (5%)	0	100	100
5	BH	256/361 (71%)	243 (95%)	13 (5%)	0	100	100
5	CH	256/361 (71%)	245 (96%)	11 (4%)	0	100	100
5	DH	256/361 (71%)	243 (95%)	13 (5%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
5	EH	256/361 (71%)	245 (96%)	11 (4%)	0	100	100
5	FH	256/361 (71%)	242 (94%)	14 (6%)	0	100	100
5	GH	256/361 (71%)	241 (94%)	15 (6%)	0	100	100
5	HH	256/361 (71%)	245 (96%)	11 (4%)	0	100	100
5	IH	256/361 (71%)	240 (94%)	16 (6%)	0	100	100
5	JH	256/361 (71%)	242 (94%)	14 (6%)	0	100	100
5	KH	256/361 (71%)	242 (94%)	14 (6%)	0	100	100
5	LH	256/361 (71%)	241 (94%)	15 (6%)	0	100	100
5	MH	256/361 (71%)	243 (95%)	13 (5%)	0	100	100
5	VH	239/361 (66%)	231 (97%)	8 (3%)	0	100	100
5	WH	239/361 (66%)	231 (97%)	8 (3%)	0	100	100
5	XH	239/361 (66%)	230 (96%)	9 (4%)	0	100	100
5	YH	239/361 (66%)	232 (97%)	7 (3%)	0	100	100
5	ZH	239/361 (66%)	232 (97%)	7 (3%)	0	100	100
6	AK	149/189 (79%)	144 (97%)	5 (3%)	0	100	100
6	BK	149/189 (79%)	144 (97%)	5 (3%)	0	100	100
6	CK	149/189 (79%)	144 (97%)	5 (3%)	0	100	100
6	DK	149/189 (79%)	144 (97%)	5 (3%)	0	100	100
6	EK	149/189 (79%)	144 (97%)	5 (3%)	0	100	100
6	FK	149/189 (79%)	144 (97%)	5 (3%)	0	100	100
6	GK	149/189 (79%)	144 (97%)	5 (3%)	0	100	100
6	HK	149/189 (79%)	144 (97%)	5 (3%)	0	100	100
6	IK	149/189 (79%)	144 (97%)	5 (3%)	0	100	100
6	JK	149/189 (79%)	144 (97%)	5 (3%)	0	100	100
6	KK	149/189 (79%)	144 (97%)	5 (3%)	0	100	100
6	LK	149/189 (79%)	144 (97%)	5 (3%)	0	100	100
6	MK	149/189 (79%)	144 (97%)	5 (3%)	0	100	100
7	AL	169/249 (68%)	159 (94%)	10 (6%)	0	100	100
7	BL	169/249 (68%)	159 (94%)	10 (6%)	0	100	100
7	CL	169/249 (68%)	159 (94%)	10 (6%)	0	100	100
7	DL	169/249 (68%)	159 (94%)	10 (6%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
7	EL	169/249 (68%)	159 (94%)	10 (6%)	0	100	100
7	FL	169/249 (68%)	159 (94%)	10 (6%)	0	100	100
7	GL	169/249 (68%)	159 (94%)	10 (6%)	0	100	100
7	HL	169/249 (68%)	159 (94%)	10 (6%)	0	100	100
7	IL	169/249 (68%)	159 (94%)	10 (6%)	0	100	100
7	JL	169/249 (68%)	159 (94%)	10 (6%)	0	100	100
7	KL	169/249 (68%)	159 (94%)	10 (6%)	0	100	100
7	LL	169/249 (68%)	159 (94%)	10 (6%)	0	100	100
7	ML	169/249 (68%)	159 (94%)	10 (6%)	0	100	100
8	AM	290/320 (91%)	265 (91%)	25 (9%)	0	100	100
8	BM	290/320 (91%)	265 (91%)	25 (9%)	0	100	100
8	CM	290/320 (91%)	265 (91%)	25 (9%)	0	100	100
8	DM	290/320 (91%)	265 (91%)	25 (9%)	0	100	100
8	EM	290/320 (91%)	265 (91%)	25 (9%)	0	100	100
8	FM	290/320 (91%)	265 (91%)	25 (9%)	0	100	100
8	GM	290/320 (91%)	265 (91%)	25 (9%)	0	100	100
8	HM	290/320 (91%)	265 (91%)	25 (9%)	0	100	100
8	IM	290/320 (91%)	266 (92%)	24 (8%)	0	100	100
8	JM	290/320 (91%)	265 (91%)	25 (9%)	0	100	100
8	KM	290/320 (91%)	266 (92%)	24 (8%)	0	100	100
8	LM	290/320 (91%)	265 (91%)	25 (9%)	0	100	100
8	MM	290/320 (91%)	265 (91%)	25 (9%)	0	100	100
9	AN	76/124 (61%)	68 (90%)	8 (10%)	0	100	100
9	BN	76/124 (61%)	68 (90%)	8 (10%)	0	100	100
9	CN	76/124 (61%)	68 (90%)	8 (10%)	0	100	100
9	DN	76/124 (61%)	68 (90%)	8 (10%)	0	100	100
9	EN	76/124 (61%)	68 (90%)	8 (10%)	0	100	100
9	FN	76/124 (61%)	68 (90%)	8 (10%)	0	100	100
9	GN	76/124 (61%)	68 (90%)	8 (10%)	0	100	100
9	HN	76/124 (61%)	68 (90%)	8 (10%)	0	100	100
9	IN	76/124 (61%)	68 (90%)	8 (10%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
9	JN	76/124 (61%)	68 (90%)	8 (10%)	0	100	100
9	KN	76/124 (61%)	68 (90%)	8 (10%)	0	100	100
9	LN	76/124 (61%)	68 (90%)	8 (10%)	0	100	100
9	MN	76/124 (61%)	68 (90%)	8 (10%)	0	100	100
All	All	22070/53344 (41%)	20933 (95%)	1137 (5%)	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	AC	178/257 (69%)	166 (93%)	12 (7%)	15	41
1	BC	178/257 (69%)	166 (93%)	12 (7%)	15	41
1	CC	178/257 (69%)	167 (94%)	11 (6%)	16	42
1	DC	178/257 (69%)	167 (94%)	11 (6%)	16	42
1	EC	178/257 (69%)	167 (94%)	11 (6%)	16	42
1	FC	178/257 (69%)	166 (93%)	12 (7%)	15	41
1	GC	178/257 (69%)	166 (93%)	12 (7%)	15	41
1	HC	178/257 (69%)	167 (94%)	11 (6%)	16	42
1	IC	178/257 (69%)	166 (93%)	12 (7%)	15	41
1	JC	178/257 (69%)	167 (94%)	11 (6%)	16	42
1	KC	178/257 (69%)	166 (93%)	12 (7%)	15	41
1	LC	178/257 (69%)	167 (94%)	11 (6%)	16	42
1	MC	178/257 (69%)	166 (93%)	12 (7%)	15	41
2	AD	121/139 (87%)	121 (100%)	0	100	100
2	Ad	118/139 (85%)	118 (100%)	0	100	100
2	BD	121/139 (87%)	121 (100%)	0	100	100
2	Bd	118/139 (85%)	118 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
2	CD	121/139 (87%)	121 (100%)	0	100 100
2	Cd	118/139 (85%)	118 (100%)	0	100 100
2	DD	121/139 (87%)	121 (100%)	0	100 100
2	Dd	118/139 (85%)	118 (100%)	0	100 100
2	ED	121/139 (87%)	121 (100%)	0	100 100
2	Ed	118/139 (85%)	118 (100%)	0	100 100
2	FD	121/139 (87%)	121 (100%)	0	100 100
2	Fd	118/139 (85%)	118 (100%)	0	100 100
2	GD	121/139 (87%)	121 (100%)	0	100 100
2	Gd	118/139 (85%)	118 (100%)	0	100 100
2	HD	121/139 (87%)	121 (100%)	0	100 100
2	Hd	118/139 (85%)	118 (100%)	0	100 100
2	ID	121/139 (87%)	121 (100%)	0	100 100
2	Id	118/139 (85%)	117 (99%)	1 (1%)	73 76
2	JD	121/139 (87%)	121 (100%)	0	100 100
2	Jd	118/139 (85%)	118 (100%)	0	100 100
2	KD	121/139 (87%)	121 (100%)	0	100 100
2	Kd	118/139 (85%)	118 (100%)	0	100 100
2	LD	121/139 (87%)	121 (100%)	0	100 100
2	Ld	118/139 (85%)	118 (100%)	0	100 100
2	MD	121/139 (87%)	121 (100%)	0	100 100
2	Md	118/139 (85%)	118 (100%)	0	100 100
3	AF	53/237 (22%)	46 (87%)	7 (13%)	4 19
3	Af	49/237 (21%)	49 (100%)	0	100 100
3	BF	53/237 (22%)	46 (87%)	7 (13%)	4 19
3	Bf	49/237 (21%)	49 (100%)	0	100 100
3	CF	53/237 (22%)	46 (87%)	7 (13%)	4 19
3	Cf	49/237 (21%)	49 (100%)	0	100 100
3	DF	53/237 (22%)	46 (87%)	7 (13%)	4 19
3	Df	49/237 (21%)	49 (100%)	0	100 100
3	EF	53/237 (22%)	46 (87%)	7 (13%)	4 19

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
3	Ef	49/237 (21%)	49 (100%)	0	100 100
3	FF	53/237 (22%)	46 (87%)	7 (13%)	4 19
3	Ff	49/237 (21%)	49 (100%)	0	100 100
3	GF	53/237 (22%)	46 (87%)	7 (13%)	4 19
3	Gf	49/237 (21%)	49 (100%)	0	100 100
3	HF	53/237 (22%)	46 (87%)	7 (13%)	4 19
3	Hf	49/237 (21%)	49 (100%)	0	100 100
3	IF	53/237 (22%)	46 (87%)	7 (13%)	4 19
3	If	49/237 (21%)	49 (100%)	0	100 100
3	JF	53/237 (22%)	46 (87%)	7 (13%)	4 19
3	Jf	49/237 (21%)	49 (100%)	0	100 100
3	KF	53/237 (22%)	46 (87%)	7 (13%)	4 19
3	Kf	49/237 (21%)	49 (100%)	0	100 100
3	LF	53/237 (22%)	46 (87%)	7 (13%)	4 19
3	Lf	49/237 (21%)	49 (100%)	0	100 100
3	MF	53/237 (22%)	46 (87%)	7 (13%)	4 19
3	Mf	49/237 (21%)	49 (100%)	0	100 100
3	VF	53/237 (22%)	46 (87%)	7 (13%)	4 19
3	WF	53/237 (22%)	46 (87%)	7 (13%)	4 19
3	XF	53/237 (22%)	46 (87%)	7 (13%)	4 19
3	YF	53/237 (22%)	46 (87%)	7 (13%)	4 19
3	ZF	53/237 (22%)	46 (87%)	7 (13%)	4 19
4	AG	31/765 (4%)	21 (68%)	10 (32%)	0 1
4	BG	31/765 (4%)	21 (68%)	10 (32%)	0 1
4	CG	31/765 (4%)	21 (68%)	10 (32%)	0 1
4	DG	31/765 (4%)	21 (68%)	10 (32%)	0 1
4	EG	31/765 (4%)	21 (68%)	10 (32%)	0 1
4	FG	31/765 (4%)	21 (68%)	10 (32%)	0 1
4	GG	31/765 (4%)	21 (68%)	10 (32%)	0 1
4	HG	31/765 (4%)	21 (68%)	10 (32%)	0 1
4	IG	31/765 (4%)	21 (68%)	10 (32%)	0 1

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
4	JG	31/765 (4%)	21 (68%)	10 (32%)	0 1
4	KG	31/765 (4%)	21 (68%)	10 (32%)	0 1
4	LG	31/765 (4%)	21 (68%)	10 (32%)	0 1
4	MG	31/765 (4%)	21 (68%)	10 (32%)	0 1
4	VG	31/765 (4%)	21 (68%)	10 (32%)	0 1
4	WG	31/765 (4%)	21 (68%)	10 (32%)	0 1
4	XG	31/765 (4%)	21 (68%)	10 (32%)	0 1
4	YG	31/765 (4%)	21 (68%)	10 (32%)	0 1
4	ZG	31/765 (4%)	21 (68%)	10 (32%)	0 1
5	AH	220/300 (73%)	220 (100%)	0	100 100
5	BH	220/300 (73%)	220 (100%)	0	100 100
5	CH	220/300 (73%)	219 (100%)	1 (0%)	81 81
5	DH	220/300 (73%)	219 (100%)	1 (0%)	81 81
5	EH	220/300 (73%)	220 (100%)	0	100 100
5	FH	220/300 (73%)	220 (100%)	0	100 100
5	GH	220/300 (73%)	220 (100%)	0	100 100
5	HH	220/300 (73%)	220 (100%)	0	100 100
5	IH	220/300 (73%)	219 (100%)	1 (0%)	81 81
5	JH	220/300 (73%)	220 (100%)	0	100 100
5	KH	220/300 (73%)	220 (100%)	0	100 100
5	LH	220/300 (73%)	220 (100%)	0	100 100
5	MH	220/300 (73%)	220 (100%)	0	100 100
5	VH	207/300 (69%)	207 (100%)	0	100 100
5	WH	207/300 (69%)	207 (100%)	0	100 100
5	XH	207/300 (69%)	207 (100%)	0	100 100
5	YH	207/300 (69%)	207 (100%)	0	100 100
5	ZH	207/300 (69%)	207 (100%)	0	100 100
6	AK	129/163 (79%)	119 (92%)	10 (8%)	11 36
6	BK	129/163 (79%)	119 (92%)	10 (8%)	11 36
6	CK	129/163 (79%)	119 (92%)	10 (8%)	11 36
6	DK	129/163 (79%)	118 (92%)	11 (8%)	10 33

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
6	EK	129/163 (79%)	119 (92%)	10 (8%)	11	36
6	FK	129/163 (79%)	119 (92%)	10 (8%)	11	36
6	GK	129/163 (79%)	119 (92%)	10 (8%)	11	36
6	HK	129/163 (79%)	119 (92%)	10 (8%)	11	36
6	IK	129/163 (79%)	118 (92%)	11 (8%)	10	33
6	JK	129/163 (79%)	119 (92%)	10 (8%)	11	36
6	KK	129/163 (79%)	119 (92%)	10 (8%)	11	36
6	LK	129/163 (79%)	119 (92%)	10 (8%)	11	36
6	MK	129/163 (79%)	119 (92%)	10 (8%)	11	36
7	AL	148/203 (73%)	139 (94%)	9 (6%)	17	43
7	BL	148/203 (73%)	139 (94%)	9 (6%)	17	43
7	CL	148/203 (73%)	139 (94%)	9 (6%)	17	43
7	DL	148/203 (73%)	139 (94%)	9 (6%)	17	43
7	EL	148/203 (73%)	139 (94%)	9 (6%)	17	43
7	FL	148/203 (73%)	139 (94%)	9 (6%)	17	43
7	GL	148/203 (73%)	139 (94%)	9 (6%)	17	43
7	HL	148/203 (73%)	139 (94%)	9 (6%)	17	43
7	IL	148/203 (73%)	139 (94%)	9 (6%)	17	43
7	JL	148/203 (73%)	139 (94%)	9 (6%)	17	43
7	KL	148/203 (73%)	139 (94%)	9 (6%)	17	43
7	LL	148/203 (73%)	139 (94%)	9 (6%)	17	43
7	ML	148/203 (73%)	139 (94%)	9 (6%)	17	43
8	AM	251/276 (91%)	241 (96%)	10 (4%)	28	52
8	BM	251/276 (91%)	241 (96%)	10 (4%)	28	52
8	CM	251/276 (91%)	241 (96%)	10 (4%)	28	52
8	DM	251/276 (91%)	241 (96%)	10 (4%)	28	52
8	EM	251/276 (91%)	241 (96%)	10 (4%)	28	52
8	FM	251/276 (91%)	241 (96%)	10 (4%)	28	52
8	GM	251/276 (91%)	241 (96%)	10 (4%)	28	52
8	HM	251/276 (91%)	241 (96%)	10 (4%)	28	52
8	IM	251/276 (91%)	241 (96%)	10 (4%)	28	52

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
8	JM	251/276 (91%)	241 (96%)	10 (4%)	28	52
8	KM	251/276 (91%)	241 (96%)	10 (4%)	28	52
8	LM	251/276 (91%)	241 (96%)	10 (4%)	28	52
8	MM	251/276 (91%)	241 (96%)	10 (4%)	28	52
9	AN	66/107 (62%)	60 (91%)	6 (9%)	9	32
9	BN	66/107 (62%)	60 (91%)	6 (9%)	9	32
9	CN	66/107 (62%)	60 (91%)	6 (9%)	9	32
9	DN	66/107 (62%)	60 (91%)	6 (9%)	9	32
9	EN	66/107 (62%)	60 (91%)	6 (9%)	9	32
9	FN	66/107 (62%)	60 (91%)	6 (9%)	9	32
9	GN	66/107 (62%)	60 (91%)	6 (9%)	9	32
9	HN	66/107 (62%)	60 (91%)	6 (9%)	9	32
9	IN	66/107 (62%)	60 (91%)	6 (9%)	9	32
9	JN	66/107 (62%)	60 (91%)	6 (9%)	9	32
9	KN	66/107 (62%)	60 (91%)	6 (9%)	9	32
9	LN	66/107 (62%)	60 (91%)	6 (9%)	9	32
9	MN	66/107 (62%)	60 (91%)	6 (9%)	9	32
All	All	19187/43209 (44%)	18270 (95%)	917 (5%)	24	47

All (917) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	AC	57	MET
1	AC	59	LEU
1	AC	60	LYS
1	AC	61	GLU
1	AC	66	VAL
1	AC	69	GLN
1	AC	97	ASN
1	AC	123	GLN
1	AC	137	GLN
1	AC	157	LYS
1	AC	182	ARG
1	AC	185	LYS
3	AF	207	ARG
3	AF	229	THR

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Mol	Chain	Res	Type
3	AF	238	ILE
3	AF	243	MET
3	AF	250	LEU
3	AF	263	LYS
3	AF	267	GLU
4	AG	795	GLN
4	AG	796	GLN
4	AG	802	LEU
4	AG	806	THR
4	AG	808	LEU
4	AG	810	GLN
4	AG	815	VAL
4	AG	819	VAL
4	AG	822	GLU
4	AG	824	THR
6	AK	43	CYS
6	AK	62	LYS
6	AK	63	LYS
6	AK	70	VAL
6	AK	101	GLU
6	AK	122	LYS
6	AK	129	GLU
6	AK	156	THR
6	AK	166	THR
6	AK	181	GLU
7	AL	104	LYS
7	AL	110	GLN
7	AL	130	ASP
7	AL	154	LEU
7	AL	155	LEU
7	AL	160	GLN
7	AL	190	MET
7	AL	201	LYS
7	AL	211	LYS
8	AM	69	VAL
8	AM	70	MET
8	AM	79	VAL
8	AM	171	ARG
8	AM	194	LEU
8	AM	196	MET
8	AM	262	GLU
8	AM	301	MET

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Mol	Chain	Res	Type
8	AM	303	GLU
8	AM	310	LYS
9	AN	39	ASP
9	AN	67	THR
9	AN	71	THR
9	AN	76	MET
9	AN	113	SER
9	AN	114	LEU
1	BC	35	GLN
1	BC	57	MET
1	BC	59	LEU
1	BC	60	LYS
1	BC	61	GLU
1	BC	66	VAL
1	BC	69	GLN
1	BC	123	GLN
1	BC	137	GLN
1	BC	157	LYS
1	BC	182	ARG
1	BC	185	LYS
3	BF	207	ARG
3	BF	229	THR
3	BF	238	ILE
3	BF	243	MET
3	BF	250	LEU
3	BF	263	LYS
3	BF	267	GLU
4	BG	795	GLN
4	BG	796	GLN
4	BG	802	LEU
4	BG	806	THR
4	BG	808	LEU
4	BG	810	GLN
4	BG	815	VAL
4	BG	819	VAL
4	BG	822	GLU
4	BG	824	THR
6	BK	43	CYS
6	BK	62	LYS
6	BK	63	LYS
6	BK	70	VAL
6	BK	101	GLU

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Mol	Chain	Res	Type
6	BK	122	LYS
6	BK	129	GLU
6	BK	156	THR
6	BK	166	THR
6	BK	181	GLU
7	BL	104	LYS
7	BL	110	GLN
7	BL	130	ASP
7	BL	154	LEU
7	BL	155	LEU
7	BL	160	GLN
7	BL	190	MET
7	BL	201	LYS
7	BL	211	LYS
8	BM	69	VAL
8	BM	70	MET
8	BM	79	VAL
8	BM	171	ARG
8	BM	194	LEU
8	BM	196	MET
8	BM	262	GLU
8	BM	301	MET
8	BM	303	GLU
8	BM	310	LYS
9	BN	39	ASP
9	BN	67	THR
9	BN	71	THR
9	BN	76	MET
9	BN	113	SER
9	BN	114	LEU
1	CC	57	MET
1	CC	59	LEU
1	CC	60	LYS
1	CC	61	GLU
1	CC	66	VAL
1	CC	69	GLN
1	CC	123	GLN
1	CC	137	GLN
1	CC	157	LYS
1	CC	182	ARG
1	CC	185	LYS
3	CF	207	ARG

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Mol	Chain	Res	Type
3	CF	229	THR
3	CF	238	ILE
3	CF	243	MET
3	CF	250	LEU
3	CF	263	LYS
3	CF	267	GLU
4	CG	795	GLN
4	CG	796	GLN
4	CG	802	LEU
4	CG	806	THR
4	CG	808	LEU
4	CG	810	GLN
4	CG	815	VAL
4	CG	819	VAL
4	CG	822	GLU
4	CG	824	THR
5	CH	245	GLN
6	CK	43	CYS
6	CK	62	LYS
6	CK	63	LYS
6	CK	70	VAL
6	CK	101	GLU
6	CK	122	LYS
6	CK	129	GLU
6	CK	156	THR
6	CK	166	THR
6	CK	181	GLU
7	CL	104	LYS
7	CL	110	GLN
7	CL	130	ASP
7	CL	154	LEU
7	CL	155	LEU
7	CL	160	GLN
7	CL	190	MET
7	CL	201	LYS
7	CL	211	LYS
8	CM	69	VAL
8	CM	70	MET
8	CM	79	VAL
8	CM	171	ARG
8	CM	194	LEU
8	CM	196	MET

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Mol	Chain	Res	Type
8	CM	262	GLU
8	CM	301	MET
8	CM	303	GLU
8	CM	310	LYS
9	CN	39	ASP
9	CN	67	THR
9	CN	71	THR
9	CN	76	MET
9	CN	113	SER
9	CN	114	LEU
1	DC	57	MET
1	DC	59	LEU
1	DC	60	LYS
1	DC	61	GLU
1	DC	66	VAL
1	DC	69	GLN
1	DC	123	GLN
1	DC	137	GLN
1	DC	157	LYS
1	DC	182	ARG
1	DC	185	LYS
3	DF	207	ARG
3	DF	229	THR
3	DF	238	ILE
3	DF	243	MET
3	DF	250	LEU
3	DF	263	LYS
3	DF	267	GLU
4	DG	795	GLN
4	DG	796	GLN
4	DG	802	LEU
4	DG	806	THR
4	DG	808	LEU
4	DG	810	GLN
4	DG	815	VAL
4	DG	819	VAL
4	DG	822	GLU
4	DG	824	THR
5	DH	225	ASN
6	DK	43	CYS
6	DK	62	LYS
6	DK	63	LYS

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Mol	Chain	Res	Type
6	DK	70	VAL
6	DK	100	ASN
6	DK	101	GLU
6	DK	122	LYS
6	DK	129	GLU
6	DK	156	THR
6	DK	166	THR
6	DK	181	GLU
7	DL	104	LYS
7	DL	110	GLN
7	DL	130	ASP
7	DL	154	LEU
7	DL	155	LEU
7	DL	160	GLN
7	DL	190	MET
7	DL	201	LYS
7	DL	211	LYS
8	DM	69	VAL
8	DM	70	MET
8	DM	79	VAL
8	DM	171	ARG
8	DM	194	LEU
8	DM	196	MET
8	DM	262	GLU
8	DM	301	MET
8	DM	303	GLU
8	DM	310	LYS
9	DN	39	ASP
9	DN	67	THR
9	DN	71	THR
9	DN	76	MET
9	DN	113	SER
9	DN	114	LEU
1	EC	57	MET
1	EC	59	LEU
1	EC	60	LYS
1	EC	61	GLU
1	EC	66	VAL
1	EC	69	GLN
1	EC	123	GLN
1	EC	137	GLN
1	EC	157	LYS

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Mol	Chain	Res	Type
1	EC	182	ARG
1	EC	185	LYS
3	EF	207	ARG
3	EF	229	THR
3	EF	238	ILE
3	EF	243	MET
3	EF	250	LEU
3	EF	263	LYS
3	EF	267	GLU
4	EG	795	GLN
4	EG	796	GLN
4	EG	802	LEU
4	EG	806	THR
4	EG	808	LEU
4	EG	810	GLN
4	EG	815	VAL
4	EG	819	VAL
4	EG	822	GLU
4	EG	824	THR
6	EK	43	CYS
6	EK	62	LYS
6	EK	63	LYS
6	EK	70	VAL
6	EK	101	GLU
6	EK	122	LYS
6	EK	129	GLU
6	EK	156	THR
6	EK	166	THR
6	EK	181	GLU
7	EL	104	LYS
7	EL	110	GLN
7	EL	130	ASP
7	EL	154	LEU
7	EL	155	LEU
7	EL	160	GLN
7	EL	190	MET
7	EL	201	LYS
7	EL	211	LYS
8	EM	69	VAL
8	EM	70	MET
8	EM	79	VAL
8	EM	171	ARG

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Mol	Chain	Res	Type
8	EM	194	LEU
8	EM	196	MET
8	EM	262	GLU
8	EM	301	MET
8	EM	303	GLU
8	EM	310	LYS
9	EN	39	ASP
9	EN	67	THR
9	EN	71	THR
9	EN	76	MET
9	EN	113	SER
9	EN	114	LEU
1	FC	57	MET
1	FC	59	LEU
1	FC	60	LYS
1	FC	61	GLU
1	FC	66	VAL
1	FC	69	GLN
1	FC	97	ASN
1	FC	123	GLN
1	FC	137	GLN
1	FC	157	LYS
1	FC	182	ARG
1	FC	185	LYS
3	FF	207	ARG
3	FF	229	THR
3	FF	238	ILE
3	FF	243	MET
3	FF	250	LEU
3	FF	263	LYS
3	FF	267	GLU
4	FG	795	GLN
4	FG	796	GLN
4	FG	802	LEU
4	FG	806	THR
4	FG	808	LEU
4	FG	810	GLN
4	FG	815	VAL
4	FG	819	VAL
4	FG	822	GLU
4	FG	824	THR
6	FK	43	CYS

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Mol	Chain	Res	Type
6	FK	62	LYS
6	FK	63	LYS
6	FK	70	VAL
6	FK	101	GLU
6	FK	122	LYS
6	FK	129	GLU
6	FK	156	THR
6	FK	166	THR
6	FK	181	GLU
7	FL	104	LYS
7	FL	110	GLN
7	FL	130	ASP
7	FL	154	LEU
7	FL	155	LEU
7	FL	160	GLN
7	FL	190	MET
7	FL	201	LYS
7	FL	211	LYS
8	FM	69	VAL
8	FM	70	MET
8	FM	79	VAL
8	FM	171	ARG
8	FM	194	LEU
8	FM	196	MET
8	FM	262	GLU
8	FM	301	MET
8	FM	303	GLU
8	FM	310	LYS
9	FN	39	ASP
9	FN	67	THR
9	FN	71	THR
9	FN	76	MET
9	FN	113	SER
9	FN	114	LEU
1	GC	57	MET
1	GC	59	LEU
1	GC	60	LYS
1	GC	61	GLU
1	GC	66	VAL
1	GC	69	GLN
1	GC	97	ASN
1	GC	123	GLN

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Mol	Chain	Res	Type
1	GC	137	GLN
1	GC	157	LYS
1	GC	182	ARG
1	GC	185	LYS
3	GF	207	ARG
3	GF	229	THR
3	GF	238	ILE
3	GF	243	MET
3	GF	250	LEU
3	GF	263	LYS
3	GF	267	GLU
4	GG	795	GLN
4	GG	796	GLN
4	GG	802	LEU
4	GG	806	THR
4	GG	808	LEU
4	GG	810	GLN
4	GG	815	VAL
4	GG	819	VAL
4	GG	822	GLU
4	GG	824	THR
6	GK	43	CYS
6	GK	62	LYS
6	GK	63	LYS
6	GK	70	VAL
6	GK	101	GLU
6	GK	122	LYS
6	GK	129	GLU
6	GK	156	THR
6	GK	166	THR
6	GK	181	GLU
7	GL	104	LYS
7	GL	110	GLN
7	GL	130	ASP
7	GL	154	LEU
7	GL	155	LEU
7	GL	160	GLN
7	GL	190	MET
7	GL	201	LYS
7	GL	211	LYS
8	GM	69	VAL
8	GM	70	MET

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Mol	Chain	Res	Type
8	GM	79	VAL
8	GM	171	ARG
8	GM	194	LEU
8	GM	196	MET
8	GM	262	GLU
8	GM	301	MET
8	GM	303	GLU
8	GM	310	LYS
9	GN	39	ASP
9	GN	67	THR
9	GN	71	THR
9	GN	76	MET
9	GN	113	SER
9	GN	114	LEU
1	HC	57	MET
1	HC	59	LEU
1	HC	60	LYS
1	HC	61	GLU
1	HC	66	VAL
1	HC	69	GLN
1	HC	123	GLN
1	HC	137	GLN
1	HC	157	LYS
1	HC	182	ARG
1	HC	185	LYS
3	HF	207	ARG
3	HF	229	THR
3	HF	238	ILE
3	HF	243	MET
3	HF	250	LEU
3	HF	263	LYS
3	HF	267	GLU
4	HG	795	GLN
4	HG	796	GLN
4	HG	802	LEU
4	HG	806	THR
4	HG	808	LEU
4	HG	810	GLN
4	HG	815	VAL
4	HG	819	VAL
4	HG	822	GLU
4	HG	824	THR

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Mol	Chain	Res	Type
6	HK	43	CYS
6	HK	62	LYS
6	HK	63	LYS
6	HK	70	VAL
6	HK	101	GLU
6	HK	122	LYS
6	HK	129	GLU
6	HK	156	THR
6	HK	166	THR
6	HK	181	GLU
7	HL	104	LYS
7	HL	110	GLN
7	HL	130	ASP
7	HL	154	LEU
7	HL	155	LEU
7	HL	160	GLN
7	HL	190	MET
7	HL	201	LYS
7	HL	211	LYS
8	HM	69	VAL
8	HM	70	MET
8	HM	79	VAL
8	HM	171	ARG
8	HM	194	LEU
8	HM	196	MET
8	HM	262	GLU
8	HM	301	MET
8	HM	303	GLU
8	HM	310	LYS
9	HN	39	ASP
9	HN	67	THR
9	HN	71	THR
9	HN	76	MET
9	HN	113	SER
9	HN	114	LEU
1	IC	57	MET
1	IC	59	LEU
1	IC	60	LYS
1	IC	61	GLU
1	IC	66	VAL
1	IC	69	GLN
1	IC	97	ASN

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Mol	Chain	Res	Type
1	IC	123	GLN
1	IC	137	GLN
1	IC	157	LYS
1	IC	182	ARG
1	IC	185	LYS
3	IF	207	ARG
3	IF	229	THR
3	IF	238	ILE
3	IF	243	MET
3	IF	250	LEU
3	IF	263	LYS
3	IF	267	GLU
4	IG	795	GLN
4	IG	796	GLN
4	IG	802	LEU
4	IG	806	THR
4	IG	808	LEU
4	IG	810	GLN
4	IG	815	VAL
4	IG	819	VAL
4	IG	822	GLU
4	IG	824	THR
5	IH	225	ASN
6	IK	43	CYS
6	IK	62	LYS
6	IK	63	LYS
6	IK	70	VAL
6	IK	100	ASN
6	IK	101	GLU
6	IK	122	LYS
6	IK	129	GLU
6	IK	156	THR
6	IK	166	THR
6	IK	181	GLU
7	IL	104	LYS
7	IL	110	GLN
7	IL	130	ASP
7	IL	154	LEU
7	IL	155	LEU
7	IL	160	GLN
7	IL	190	MET
7	IL	201	LYS

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Mol	Chain	Res	Type
7	IL	211	LYS
8	IM	69	VAL
8	IM	70	MET
8	IM	79	VAL
8	IM	171	ARG
8	IM	194	LEU
8	IM	196	MET
8	IM	262	GLU
8	IM	301	MET
8	IM	303	GLU
8	IM	310	LYS
9	IN	39	ASP
9	IN	67	THR
9	IN	71	THR
9	IN	76	MET
9	IN	113	SER
9	IN	114	LEU
2	Id	69	ASN
1	JC	57	MET
1	JC	59	LEU
1	JC	60	LYS
1	JC	61	GLU
1	JC	66	VAL
1	JC	69	GLN
1	JC	123	GLN
1	JC	137	GLN
1	JC	157	LYS
1	JC	182	ARG
1	JC	185	LYS
3	JF	207	ARG
3	JF	229	THR
3	JF	238	ILE
3	JF	243	MET
3	JF	250	LEU
3	JF	263	LYS
3	JF	267	GLU
4	JG	795	GLN
4	JG	796	GLN
4	JG	802	LEU
4	JG	806	THR
4	JG	808	LEU
4	JG	810	GLN

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Mol	Chain	Res	Type
4	JG	815	VAL
4	JG	819	VAL
4	JG	822	GLU
4	JG	824	THR
6	JK	43	CYS
6	JK	62	LYS
6	JK	63	LYS
6	JK	70	VAL
6	JK	101	GLU
6	JK	122	LYS
6	JK	129	GLU
6	JK	156	THR
6	JK	166	THR
6	JK	181	GLU
7	JL	104	LYS
7	JL	110	GLN
7	JL	130	ASP
7	JL	154	LEU
7	JL	155	LEU
7	JL	160	GLN
7	JL	190	MET
7	JL	201	LYS
7	JL	211	LYS
8	JM	69	VAL
8	JM	70	MET
8	JM	79	VAL
8	JM	171	ARG
8	JM	194	LEU
8	JM	196	MET
8	JM	262	GLU
8	JM	301	MET
8	JM	303	GLU
8	JM	310	LYS
9	JN	39	ASP
9	JN	67	THR
9	JN	71	THR
9	JN	76	MET
9	JN	113	SER
9	JN	114	LEU
1	KC	57	MET
1	KC	59	LEU
1	KC	60	LYS

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Mol	Chain	Res	Type
1	KC	61	GLU
1	KC	66	VAL
1	KC	69	GLN
1	KC	97	ASN
1	KC	123	GLN
1	KC	137	GLN
1	KC	157	LYS
1	KC	182	ARG
1	KC	185	LYS
3	KF	207	ARG
3	KF	229	THR
3	KF	238	ILE
3	KF	243	MET
3	KF	250	LEU
3	KF	263	LYS
3	KF	267	GLU
4	KG	795	GLN
4	KG	796	GLN
4	KG	802	LEU
4	KG	806	THR
4	KG	808	LEU
4	KG	810	GLN
4	KG	815	VAL
4	KG	819	VAL
4	KG	822	GLU
4	KG	824	THR
6	KK	43	CYS
6	KK	62	LYS
6	KK	63	LYS
6	KK	70	VAL
6	KK	101	GLU
6	KK	122	LYS
6	KK	129	GLU
6	KK	156	THR
6	KK	166	THR
6	KK	181	GLU
7	KL	104	LYS
7	KL	110	GLN
7	KL	130	ASP
7	KL	154	LEU
7	KL	155	LEU
7	KL	160	GLN

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Mol	Chain	Res	Type
7	KL	190	MET
7	KL	201	LYS
7	KL	211	LYS
8	KM	69	VAL
8	KM	70	MET
8	KM	79	VAL
8	KM	171	ARG
8	KM	194	LEU
8	KM	196	MET
8	KM	262	GLU
8	KM	301	MET
8	KM	303	GLU
8	KM	310	LYS
9	KN	39	ASP
9	KN	67	THR
9	KN	71	THR
9	KN	76	MET
9	KN	113	SER
9	KN	114	LEU
1	LC	57	MET
1	LC	59	LEU
1	LC	60	LYS
1	LC	61	GLU
1	LC	66	VAL
1	LC	69	GLN
1	LC	123	GLN
1	LC	137	GLN
1	LC	157	LYS
1	LC	182	ARG
1	LC	185	LYS
3	LF	207	ARG
3	LF	229	THR
3	LF	238	ILE
3	LF	243	MET
3	LF	250	LEU
3	LF	263	LYS
3	LF	267	GLU
4	LG	795	GLN
4	LG	796	GLN
4	LG	802	LEU
4	LG	806	THR
4	LG	808	LEU

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Mol	Chain	Res	Type
4	LG	810	GLN
4	LG	815	VAL
4	LG	819	VAL
4	LG	822	GLU
4	LG	824	THR
6	LK	43	CYS
6	LK	62	LYS
6	LK	63	LYS
6	LK	70	VAL
6	LK	101	GLU
6	LK	122	LYS
6	LK	129	GLU
6	LK	156	THR
6	LK	166	THR
6	LK	181	GLU
7	LL	104	LYS
7	LL	110	GLN
7	LL	130	ASP
7	LL	154	LEU
7	LL	155	LEU
7	LL	160	GLN
7	LL	190	MET
7	LL	201	LYS
7	LL	211	LYS
8	LM	69	VAL
8	LM	70	MET
8	LM	79	VAL
8	LM	171	ARG
8	LM	194	LEU
8	LM	196	MET
8	LM	262	GLU
8	LM	301	MET
8	LM	303	GLU
8	LM	310	LYS
9	LN	39	ASP
9	LN	67	THR
9	LN	71	THR
9	LN	76	MET
9	LN	113	SER
9	LN	114	LEU
1	MC	35	GLN
1	MC	57	MET

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Mol	Chain	Res	Type
1	MC	59	LEU
1	MC	60	LYS
1	MC	61	GLU
1	MC	66	VAL
1	MC	69	GLN
1	MC	123	GLN
1	MC	137	GLN
1	MC	157	LYS
1	MC	182	ARG
1	MC	185	LYS
3	MF	207	ARG
3	MF	229	THR
3	MF	238	ILE
3	MF	243	MET
3	MF	250	LEU
3	MF	263	LYS
3	MF	267	GLU
4	MG	795	GLN
4	MG	796	GLN
4	MG	802	LEU
4	MG	806	THR
4	MG	808	LEU
4	MG	810	GLN
4	MG	815	VAL
4	MG	819	VAL
4	MG	822	GLU
4	MG	824	THR
6	MK	43	CYS
6	MK	62	LYS
6	MK	63	LYS
6	MK	70	VAL
6	MK	101	GLU
6	MK	122	LYS
6	MK	129	GLU
6	MK	156	THR
6	MK	166	THR
6	MK	181	GLU
7	ML	104	LYS
7	ML	110	GLN
7	ML	130	ASP
7	ML	154	LEU
7	ML	155	LEU

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Mol	Chain	Res	Type
7	ML	160	GLN
7	ML	190	MET
7	ML	201	LYS
7	ML	211	LYS
8	MM	69	VAL
8	MM	70	MET
8	MM	79	VAL
8	MM	171	ARG
8	MM	194	LEU
8	MM	196	MET
8	MM	262	GLU
8	MM	301	MET
8	MM	303	GLU
8	MM	310	LYS
9	MN	39	ASP
9	MN	67	THR
9	MN	71	THR
9	MN	76	MET
9	MN	113	SER
9	MN	114	LEU
3	VF	207	ARG
3	VF	229	THR
3	VF	238	ILE
3	VF	243	MET
3	VF	250	LEU
3	VF	263	LYS
3	VF	267	GLU
4	VG	795	GLN
4	VG	796	GLN
4	VG	802	LEU
4	VG	806	THR
4	VG	808	LEU
4	VG	810	GLN
4	VG	815	VAL
4	VG	819	VAL
4	VG	822	GLU
4	VG	824	THR
3	WF	207	ARG
3	WF	229	THR
3	WF	238	ILE
3	WF	243	MET
3	WF	250	LEU

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Mol	Chain	Res	Type
3	WF	263	LYS
3	WF	267	GLU
4	WG	795	GLN
4	WG	796	GLN
4	WG	802	LEU
4	WG	806	THR
4	WG	808	LEU
4	WG	810	GLN
4	WG	815	VAL
4	WG	819	VAL
4	WG	822	GLU
4	WG	824	THR
3	XF	207	ARG
3	XF	229	THR
3	XF	238	ILE
3	XF	243	MET
3	XF	250	LEU
3	XF	263	LYS
3	XF	267	GLU
4	XG	795	GLN
4	XG	796	GLN
4	XG	802	LEU
4	XG	806	THR
4	XG	808	LEU
4	XG	810	GLN
4	XG	815	VAL
4	XG	819	VAL
4	XG	822	GLU
4	XG	824	THR
3	YF	207	ARG
3	YF	229	THR
3	YF	238	ILE
3	YF	243	MET
3	YF	250	LEU
3	YF	263	LYS
3	YF	267	GLU
4	YG	795	GLN
4	YG	796	GLN
4	YG	802	LEU
4	YG	806	THR
4	YG	808	LEU
4	YG	810	GLN

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Mol	Chain	Res	Type
4	YG	815	VAL
4	YG	819	VAL
4	YG	822	GLU
4	YG	824	THR
3	ZF	207	ARG
3	ZF	229	THR
3	ZF	238	ILE
3	ZF	243	MET
3	ZF	250	LEU
3	ZF	263	LYS
3	ZF	267	GLU
4	ZG	795	GLN
4	ZG	796	GLN
4	ZG	802	LEU
4	ZG	806	THR
4	ZG	808	LEU
4	ZG	810	GLN
4	ZG	815	VAL
4	ZG	819	VAL
4	ZG	822	GLU
4	ZG	824	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (181) such sidechains are listed below:

Mol	Chain	Res	Type
1	AC	35	GLN
1	AC	97	ASN
2	AD	32	ASN
5	AH	132	GLN
5	AH	245	GLN
8	AM	40	GLN
8	AM	80	GLN
8	AM	88	ASN
8	AM	164	GLN
9	AN	61	ASN
2	Ad	152	GLN
1	BC	69	GLN
2	BD	32	ASN
5	BH	156	GLN
5	BH	222	ASN
5	BH	234	ASN
5	BH	335	HIS

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Mol	Chain	Res	Type
5	BH	350	HIS
6	BK	147	GLN
7	BL	49	ASN
8	BM	40	GLN
8	BM	80	GLN
8	BM	88	ASN
8	BM	164	GLN
8	BM	193	GLN
9	BN	61	ASN
9	BN	115	GLN
2	Bd	152	GLN
1	CC	35	GLN
3	CF	266	GLN
4	CG	818	GLN
5	CH	156	GLN
5	CH	203	ASN
5	CH	205	GLN
5	CH	222	ASN
5	CH	335	HIS
8	CM	40	GLN
8	CM	80	GLN
8	CM	88	ASN
8	CM	164	GLN
9	CN	61	ASN
1	DC	35	GLN
1	DC	137	GLN
4	DG	818	GLN
5	DH	205	GLN
5	DH	350	HIS
6	DK	72	GLN
6	DK	100	ASN
8	DM	40	GLN
8	DM	80	GLN
8	DM	88	ASN
8	DM	164	GLN
9	DN	61	ASN
9	DN	115	GLN
3	Df	213	GLN
2	ED	32	ASN
2	ED	78	ASN
2	ED	152	GLN
5	EH	308	ASN

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Mol	Chain	Res	Type
5	EH	335	HIS
8	EM	40	GLN
8	EM	80	GLN
8	EM	164	GLN
9	EN	61	ASN
9	EN	115	GLN
1	FC	97	ASN
2	FD	32	ASN
2	FD	138	GLN
3	FF	266	GLN
4	FG	818	GLN
5	FH	245	GLN
8	FM	40	GLN
8	FM	80	GLN
8	FM	88	ASN
8	FM	164	GLN
8	FM	193	GLN
9	FN	61	ASN
1	GC	35	GLN
1	GC	97	ASN
1	GC	137	GLN
1	GC	241	HIS
2	GD	32	ASN
5	GH	222	ASN
5	GH	335	HIS
6	GK	147	GLN
8	GM	40	GLN
8	GM	80	GLN
8	GM	88	ASN
8	GM	164	GLN
8	GM	193	GLN
9	GN	61	ASN
1	HC	35	GLN
1	HC	137	GLN
2	HD	32	ASN
2	HD	138	GLN
3	HF	266	GLN
5	HH	216	GLN
5	HH	222	ASN
5	HH	257	GLN
8	HM	40	GLN
8	HM	80	GLN

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Mol	Chain	Res	Type
8	HM	88	ASN
8	HM	164	GLN
9	HN	61	ASN
2	Hd	146	HIS
2	Hd	152	GLN
3	Hf	213	GLN
1	IC	97	ASN
3	IF	266	GLN
5	IH	335	HIS
5	IH	350	HIS
6	IK	100	ASN
8	IM	40	GLN
8	IM	80	GLN
8	IM	88	ASN
8	IM	164	GLN
9	IN	61	ASN
1	JC	35	GLN
1	JC	123	GLN
2	JD	32	ASN
3	JF	266	GLN
5	JH	118	ASN
5	JH	173	GLN
5	JH	245	GLN
7	JL	232	GLN
8	JM	40	GLN
8	JM	80	GLN
8	JM	88	ASN
8	JM	164	GLN
9	JN	61	ASN
3	Jf	266	GLN
1	KC	97	ASN
2	KD	32	ASN
2	KD	80	GLN
3	KF	266	GLN
5	KH	132	GLN
5	KH	156	GLN
5	KH	234	ASN
5	KH	308	ASN
6	KK	147	GLN
7	KL	232	GLN
8	KM	40	GLN
8	KM	80	GLN

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Mol	Chain	Res	Type
8	KM	88	ASN
8	KM	164	GLN
9	KN	61	ASN
9	KN	73	HIS
2	LD	32	ASN
3	LF	266	GLN
5	LH	216	GLN
6	LK	147	GLN
8	LM	40	GLN
8	LM	80	GLN
8	LM	88	ASN
8	LM	164	GLN
9	LN	61	ASN
1	MC	241	HIS
2	MD	80	GLN
5	MH	234	ASN
8	MM	40	GLN
8	MM	80	GLN
8	MM	88	ASN
8	MM	164	GLN
9	MN	61	ASN
2	Md	80	GLN
5	VH	156	GLN
5	VH	222	ASN
5	VH	225	ASN
5	VH	335	HIS
3	WF	266	GLN
5	WH	123	ASN
5	XH	222	ASN
3	YF	251	GLN
3	YF	266	GLN
5	YH	234	ASN
5	YH	282	HIS
5	YH	335	HIS
4	ZG	818	GLN
5	ZH	203	ASN
5	ZH	245	GLN
5	ZH	308	ASN

5.3.3 RNA

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates

There are no oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

The following chains have linkage breaks:

Mol	Chain	Number of breaks
11	AX	1
11	BX	1
11	CX	1
11	DX	1
11	EX	1
11	FX	1
11	GX	1
11	HX	1
11	IX	1
11	JX	1
11	KX	1
11	LX	1
11	MX	1
11	VX	1
11	WX	1
11	XX	1
11	YX	1
11	ZX	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	AX	38:UNK	C	70:UNK	N	24.20
1	BX	38:UNK	C	70:UNK	N	24.20
1	CX	38:UNK	C	70:UNK	N	24.20
1	DX	38:UNK	C	70:UNK	N	24.20
1	EX	38:UNK	C	70:UNK	N	24.20
1	FX	38:UNK	C	70:UNK	N	24.20
1	GX	38:UNK	C	70:UNK	N	24.20
1	HX	38:UNK	C	70:UNK	N	24.20
1	IX	38:UNK	C	70:UNK	N	24.20
1	JX	38:UNK	C	70:UNK	N	24.20
1	KX	38:UNK	C	70:UNK	N	24.20
1	LX	38:UNK	C	70:UNK	N	24.20
1	MX	38:UNK	C	70:UNK	N	24.20
1	VX	38:UNK	C	70:UNK	N	24.20
1	WX	38:UNK	C	70:UNK	N	24.20
1	XX	38:UNK	C	70:UNK	N	24.20
1	YX	38:UNK	C	70:UNK	N	24.20
1	ZX	38:UNK	C	70:UNK	N	24.20

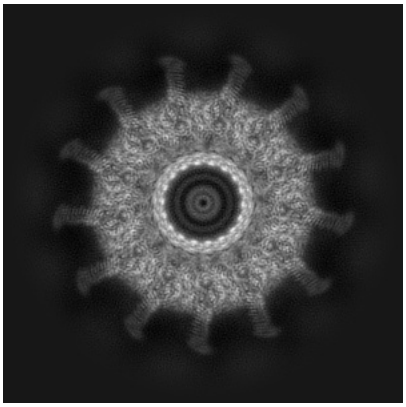
6 Map visualisation

This section contains visualisations of the EMDB entry EMD-24004. These allow visual inspection of the internal detail of the map and identification of artifacts.

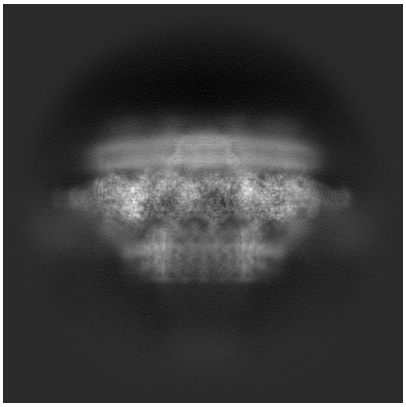
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections

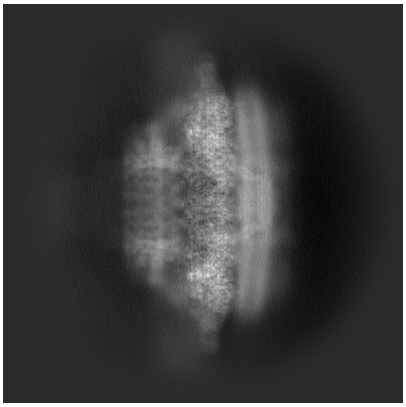
6.1.1 Primary map



X



Y

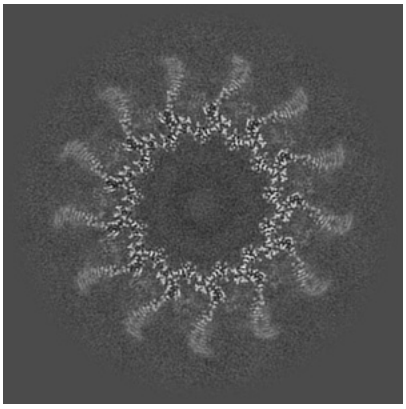


Z

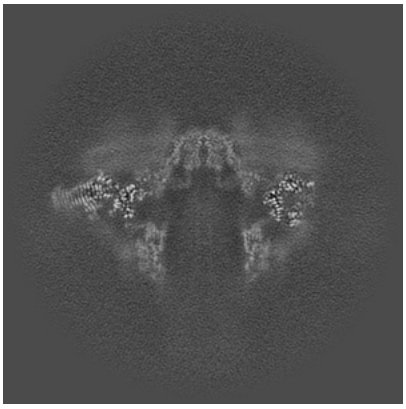
The images above show the map projected in three orthogonal directions.

6.2 Central slices

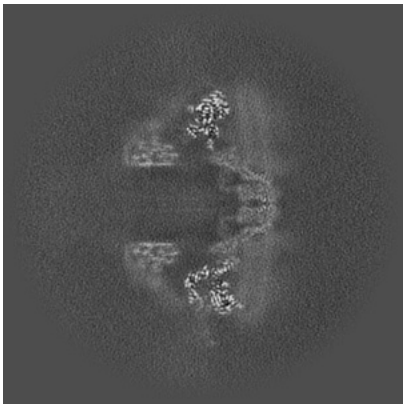
6.2.1 Primary map



X Index: 255



Y Index: 255

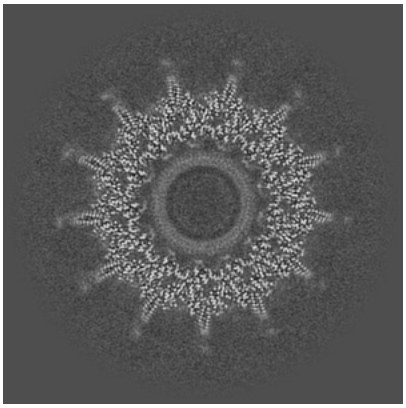


Z Index: 255

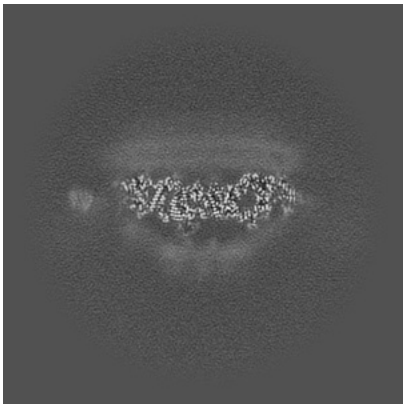
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices (i)

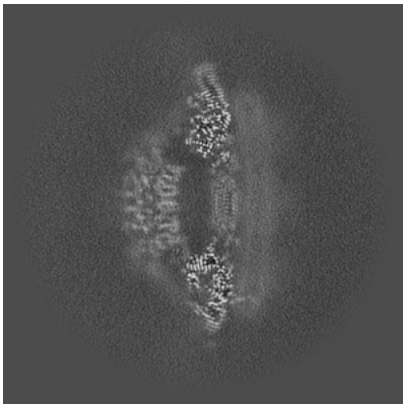
6.3.1 Primary map



X Index: 274



Y Index: 164



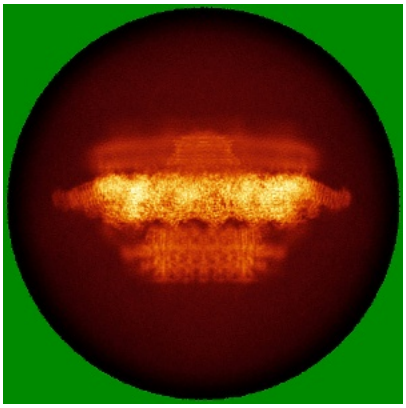
Z Index: 309

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

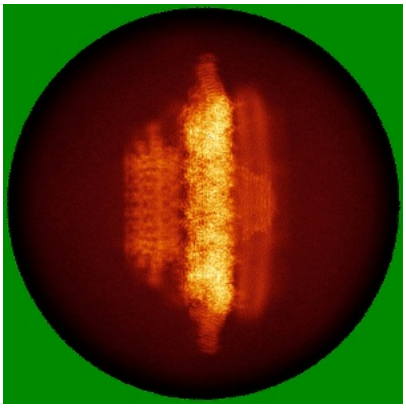
6.4 Orthogonal standard-deviation projections (False-color) (i)

6.4.1 Primary map

X



Y

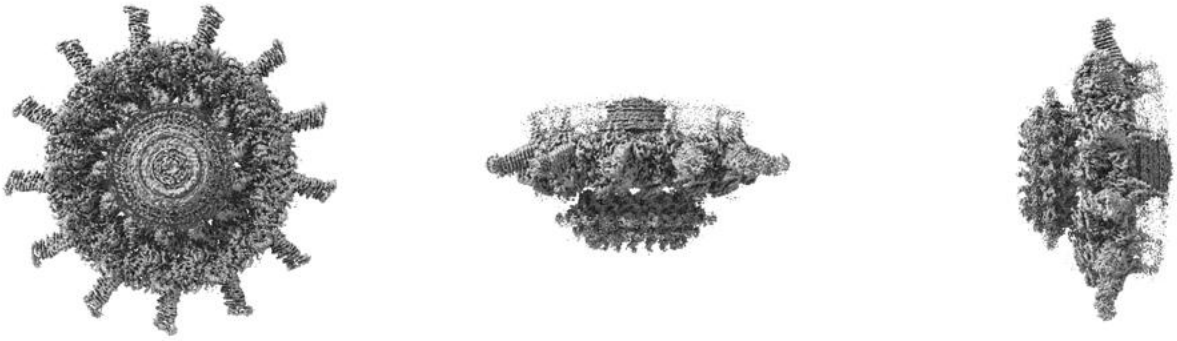


Z

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

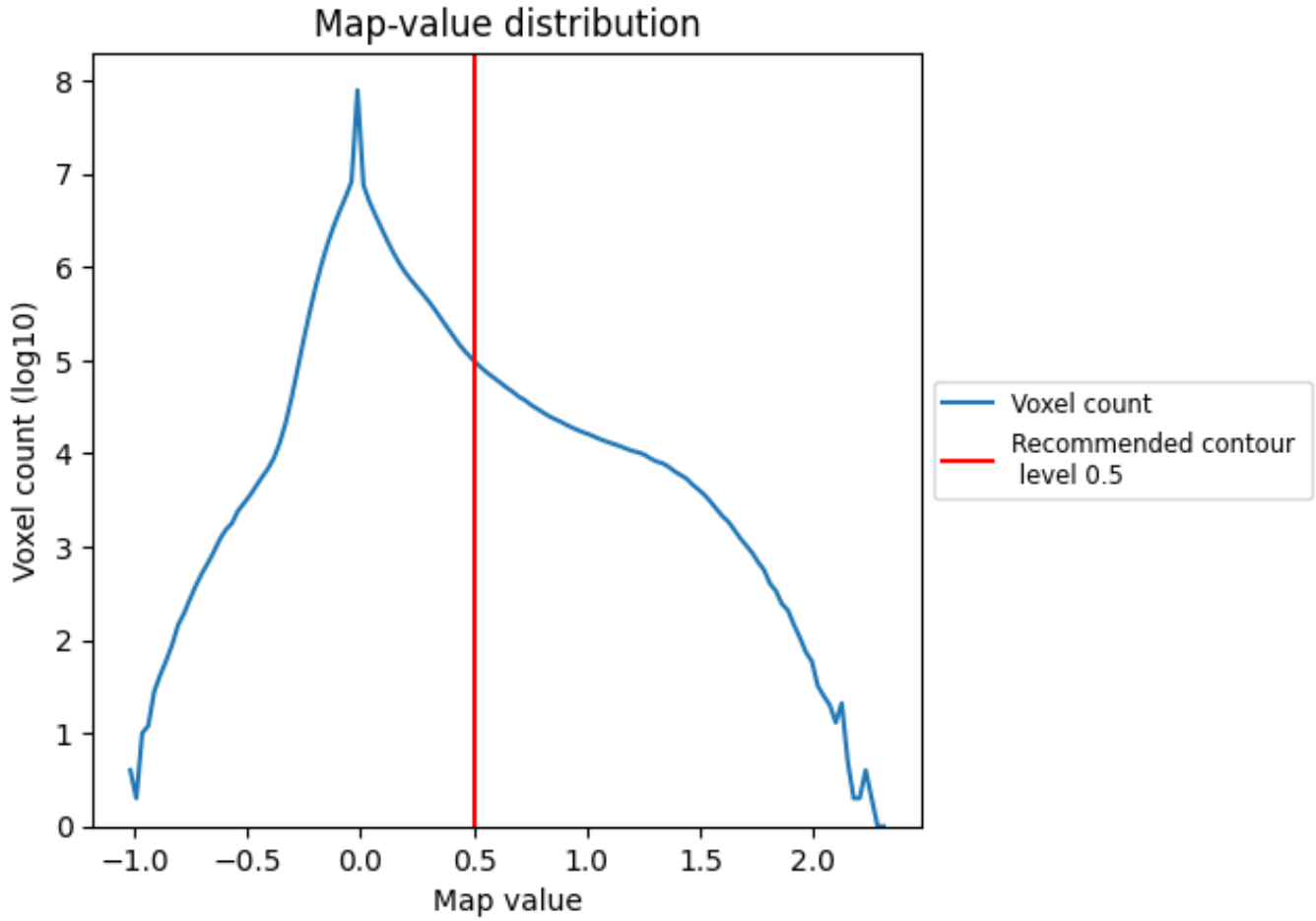
6.6 Mask visualisation (i)

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

7 Map analysis

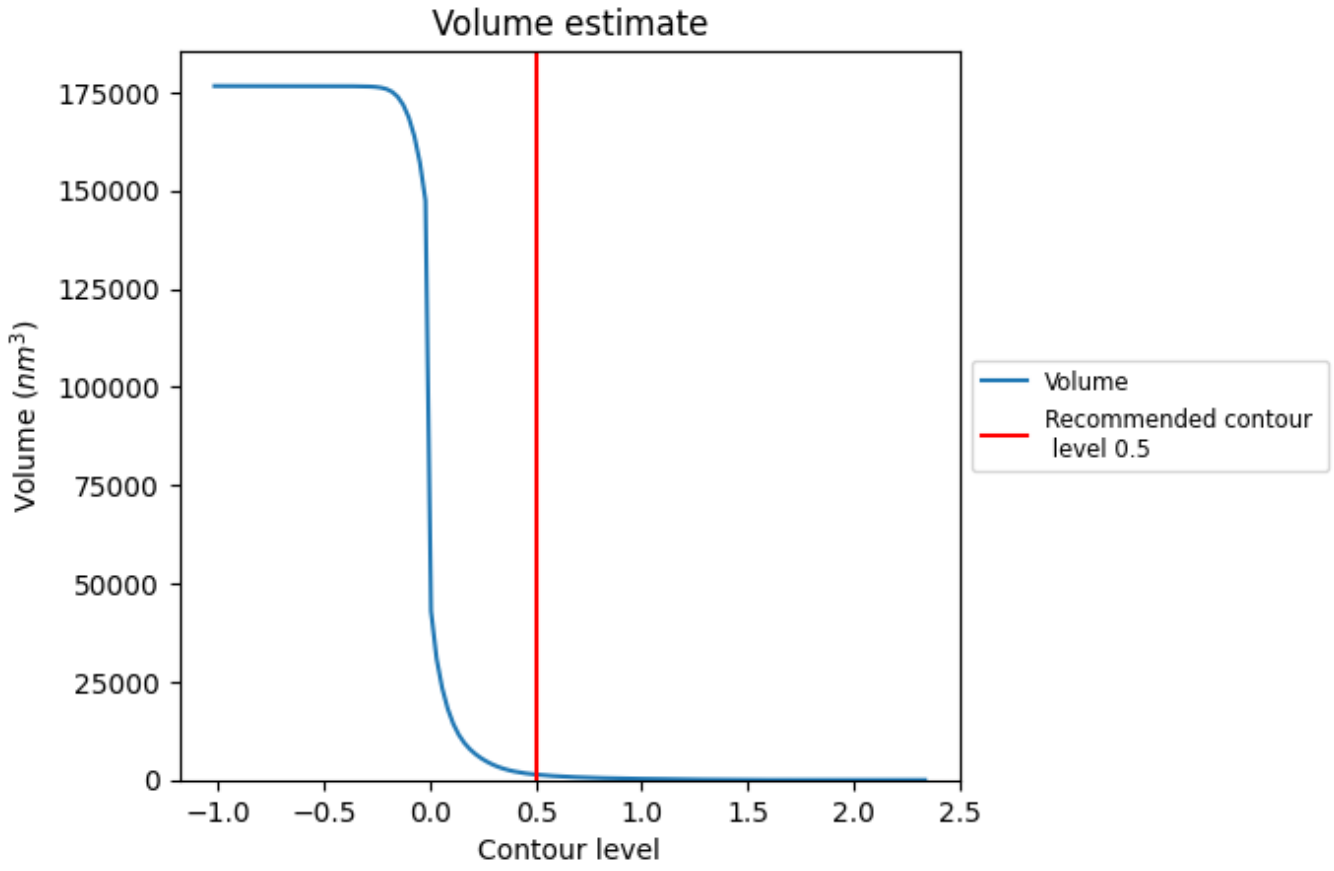
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution



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.

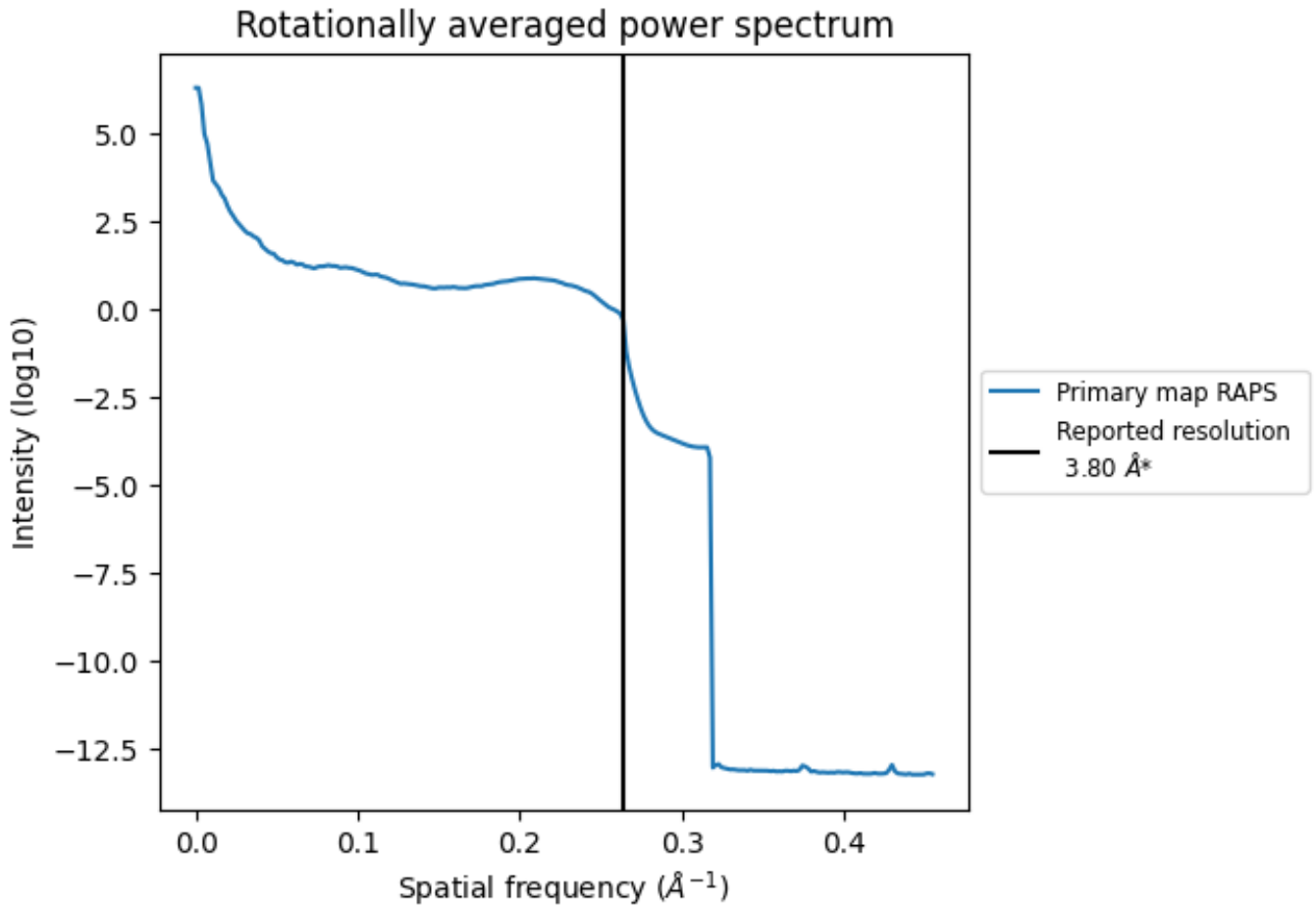
7.2 Volume estimate



The volume at the recommended contour level is 1376 nm^3; this corresponds to an approximate mass of 1243 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 0.263 $Å^{-1}$

8 Fourier-Shell correlation (i)

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

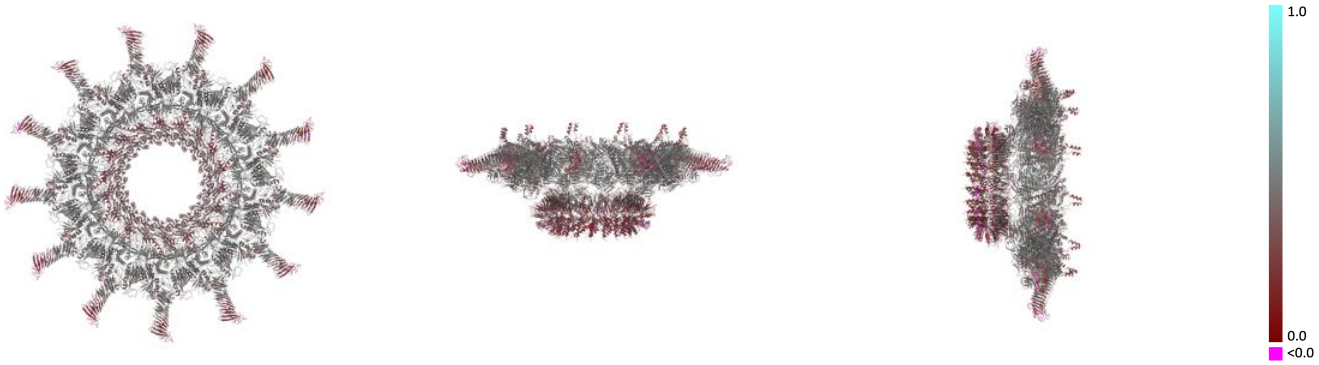
9 Map-model fit

This section contains information regarding the fit between EMDB map EMD-24004 and PDB model 7MUC. Per-residue inclusion information can be found in section 3 on page 22.

9.1 Map-model overlay

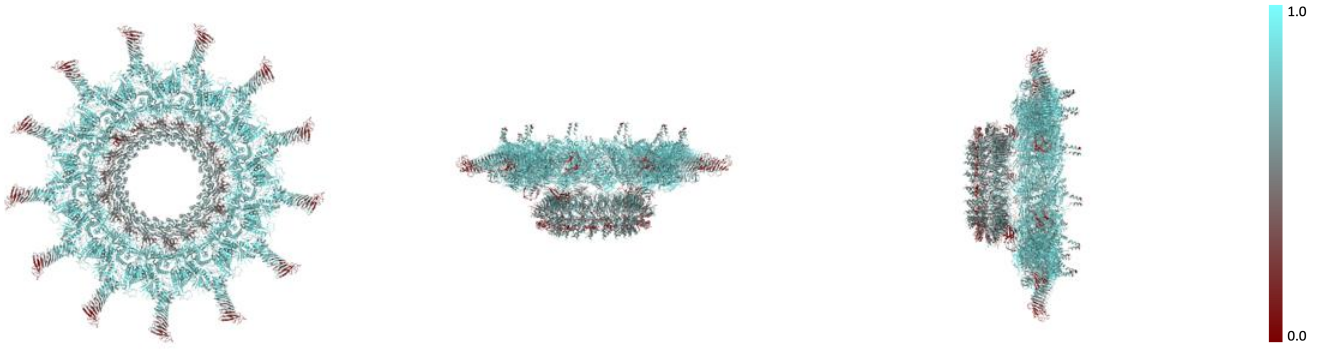
This section was not generated.

9.2 Q-score mapped to coordinate model



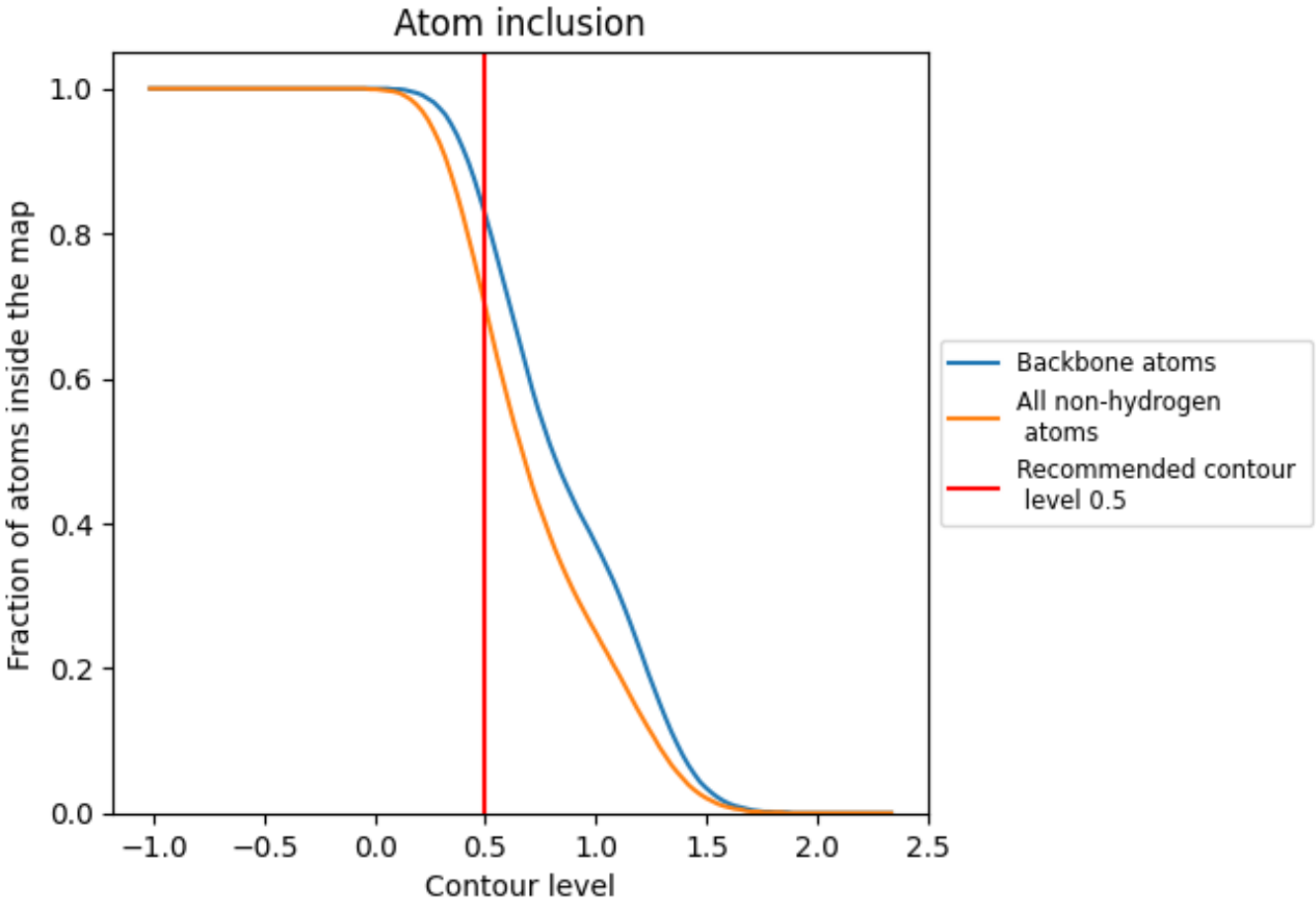
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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



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 (0.5).

9.4 Atom inclusion (i)



At the recommended contour level, 83% of all backbone atoms, 70% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary

The table lists the average atom inclusion at the recommended contour level (0.5) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.6980	0.3930
AC	0.8190	0.4550
AD	0.8240	0.4650
AF	0.4500	0.1820
AG	0.6250	0.2810
AH	0.7010	0.3920
AK	0.8420	0.4630
AL	0.8060	0.4360
AM	0.5890	0.3380
AN	0.7550	0.4370
AU	0.9780	0.5100
AX	0.2000	0.1820
Ad	0.8210	0.4560
Af	0.6810	0.3970
BC	0.8200	0.4520
BD	0.8260	0.4700
BF	0.4500	0.2010
BG	0.5960	0.2570
BH	0.7030	0.3960
BK	0.8310	0.4620
BL	0.8020	0.4370
BM	0.5840	0.3440
BN	0.7740	0.4500
BU	0.9780	0.4820
BX	0.2460	0.1950
Bd	0.8180	0.4450
Bf	0.6950	0.4160
CC	0.8130	0.4560
CD	0.8070	0.4570
CF	0.4630	0.2440
CG	0.6140	0.2620
CH	0.7140	0.3900
CK	0.8270	0.4540
CL	0.7960	0.4250
CM	0.5760	0.3460

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Chain	Atom inclusion	Q-score
CN	0.7760	0.4520
CU	0.9560	0.4740
CX	0.2290	0.2120
Cd	0.8120	0.4340
Cf	0.6990	0.4220
DC	0.8190	0.4540
DD	0.8100	0.4530
DF	0.4500	0.2320
DG	0.6210	0.2840
DH	0.6940	0.3890
DK	0.8180	0.4550
DL	0.7780	0.4060
DM	0.5820	0.3490
DN	0.7660	0.4460
DU	0.9780	0.4920
DX	0.1880	0.1930
Dd	0.7940	0.4430
Df	0.6400	0.3980
EC	0.8050	0.4460
ED	0.8070	0.4580
EF	0.4250	0.2200
EG	0.6360	0.2780
EH	0.6890	0.3750
EK	0.8260	0.4550
EL	0.7920	0.4160
EM	0.5730	0.3520
EN	0.7640	0.4370
EU	0.9330	0.4800
EX	0.2420	0.2010
Ed	0.8130	0.4420
Ef	0.6700	0.4050
FC	0.8140	0.4510
FD	0.8170	0.4560
FF	0.4220	0.2120
FG	0.6210	0.2620
FH	0.6860	0.3820
FK	0.8230	0.4560
FL	0.7870	0.4180
FM	0.5760	0.3380
FN	0.7520	0.4420
FU	0.9780	0.4960
FX	0.1750	0.1330

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Chain	Atom inclusion	Q-score
Fd	0.8020	0.4400
Ff	0.6670	0.4050
GC	0.8220	0.4500
GD	0.8070	0.4550
GF	0.4290	0.2140
GG	0.6030	0.2800
GH	0.6920	0.3900
GK	0.8220	0.4580
GL	0.8030	0.4330
GM	0.5740	0.3360
GN	0.7660	0.4420
GU	1.0000	0.4930
GX	0.1380	0.1290
Gd	0.8120	0.4500
Gf	0.6720	0.4280
HC	0.8160	0.4510
HD	0.8140	0.4590
HF	0.4460	0.2430
HG	0.6030	0.2610
HH	0.7010	0.3910
HK	0.8270	0.4580
HL	0.8200	0.4290
HM	0.5780	0.3400
HN	0.7730	0.4470
HU	0.9780	0.4860
HX	0.2460	0.1990
Hd	0.7990	0.4530
Hf	0.6540	0.4050
IC	0.8060	0.4400
ID	0.8010	0.4520
IF	0.4690	0.2400
IG	0.5880	0.2820
IH	0.6970	0.3920
IK	0.8150	0.4440
IL	0.8020	0.4310
IM	0.5610	0.3410
IN	0.7600	0.4540
IU	0.9330	0.4920
IX	0.2080	0.1450
Id	0.7900	0.4370
If	0.6900	0.4050
JC	0.8020	0.4300

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Chain	Atom inclusion	Q-score
JD	0.8080	0.4530
JF	0.4610	0.2370
JG	0.5440	0.2460
JH	0.7090	0.3910
JK	0.8150	0.4500
JL	0.8010	0.4270
JM	0.5790	0.3450
JN	0.7660	0.4470
JU	0.9780	0.5140
JX	0.1960	0.2150
Jd	0.8140	0.4430
Jf	0.6720	0.4180
KC	0.8100	0.4390
KD	0.8160	0.4540
KF	0.5010	0.2500
KG	0.5810	0.2350
KH	0.6990	0.3820
KK	0.8210	0.4490
KL	0.8070	0.4250
KM	0.5820	0.3400
KN	0.7710	0.4510
KU	0.9330	0.5000
KX	0.2710	0.2290
Kd	0.8030	0.4480
Kf	0.6930	0.4170
LC	0.8160	0.4490
LD	0.8170	0.4600
LF	0.4780	0.2500
LG	0.5920	0.2450
LH	0.6940	0.3830
LK	0.8320	0.4530
LL	0.8080	0.4230
LM	0.6000	0.3570
LN	0.7810	0.4580
LU	0.8890	0.4570
LX	0.2120	0.2290
Ld	0.8210	0.4480
Lf	0.6970	0.4160
MC	0.8250	0.4510
MD	0.8140	0.4600
MF	0.4670	0.2240
MG	0.6030	0.2910

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Chain	Atom inclusion	Q-score
MH	0.7230	0.4020
MK	0.8300	0.4590
ML	0.8120	0.4330
MM	0.6080	0.3410
MN	0.7730	0.4500
MU	0.9560	0.4940
MX	0.2250	0.2070
Md	0.8130	0.4450
Mf	0.6810	0.4110
VF	0.4710	0.2340
VG	0.5920	0.2670
VH	0.5600	0.3600
VX	0.2630	0.2330
WF	0.4120	0.1920
WG	0.6100	0.2230
WH	0.4280	0.3360
WX	0.1750	0.1610
XF	0.4520	0.2420
XG	0.6100	0.2740
XH	0.5640	0.3700
XX	0.1960	0.1580
YF	0.4800	0.2430
YG	0.6180	0.2900
YH	0.5060	0.3500
YX	0.2290	0.2210
ZF	0.4310	0.2010
ZG	0.5880	0.2450
ZH	0.4770	0.3490
ZX	0.2170	0.2010