



Full wwPDB EM Validation Report ⓘ

Mar 5, 2026 – 05:56 PM UTC

PDB ID : 3FOH / pdb_00003foh
EMDB ID : EMD-1126
Title : Fitting of gp18M crystal structure into 3D cryo-EM reconstruction of bacteriophage T4 extended tail
Authors : Aksyuk, A.A.; Leiman, P.G.; Kurochkina, L.P.; Shneider, M.M.; Kostyuchenko, V.A.; Mesyanzhinov, V.V.; Rossmann, M.G.
Deposited on : 2008-12-30
Resolution : 15.00 Å (reported)
Based on initial model : 3FOA

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:

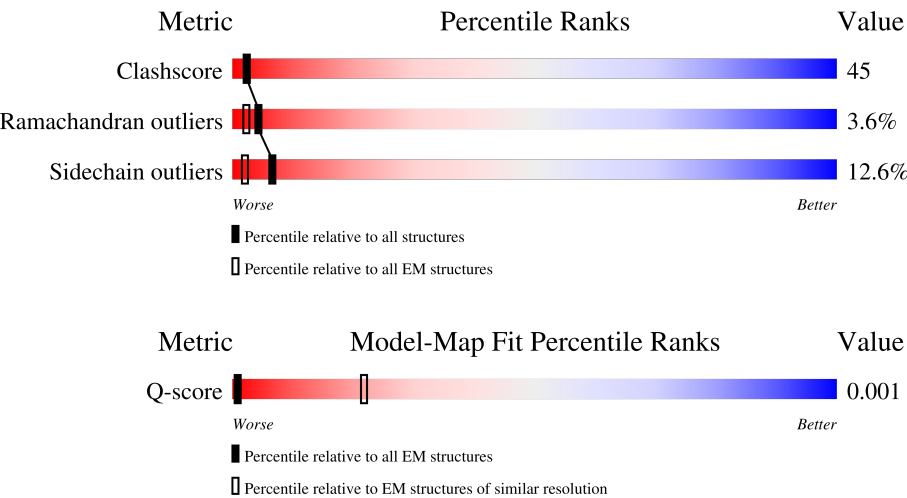
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh &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

1 Overall quality at a glance i

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

The reported resolution of this entry is 15.00 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	37 (14.50 - 15.50)

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

Mol	Chain	Length	Quality of chain
1	A	510	50% 37% 46% 10% 6%
1	B	510	36% 36% 46% 10% 6%
1	C	510	28% 37% 45% 11% 6%
1	D	510	27% 37% 46% 10% 6%

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Mol	Chain	Length	Quality of chain
1	E	510	45% 36% 46% 10% • 6%
1	F	510	57% 36% 46% 10% • 6%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 21678 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 Tail sheath protein Gp18.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	479	Total	C	N	O	S	0	0
			3613	2279	600	726	8		
1	B	479	Total	C	N	O	S	0	0
			3613	2279	600	726	8		
1	C	479	Total	C	N	O	S	0	0
			3613	2279	600	726	8		
1	D	479	Total	C	N	O	S	0	0
			3613	2279	600	726	8		
1	E	479	Total	C	N	O	S	0	0
			3613	2279	600	726	8		
1	F	479	Total	C	N	O	S	0	0
			3613	2279	600	726	8		

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	100	GLU	ASP	SEE REMARK 999	UNP P13332
A	148	ALA	GLY	SEE REMARK 999	UNP P13332
A	150	ILE	ASN	SEE REMARK 999	UNP P13332
A	151	ILE	TYR	SEE REMARK 999	UNP P13332
A	301	GLY	GLU	SEE REMARK 999	UNP P13332
A	399	VAL	ALA	SEE REMARK 999	UNP P13332
A	454	TYR	HIS	SEE REMARK 999	UNP P13332
A	510	PRO	ARG	engineered mutation	UNP P13332
B	100	GLU	ASP	SEE REMARK 999	UNP P13332
B	148	ALA	GLY	SEE REMARK 999	UNP P13332
B	150	ILE	ASN	SEE REMARK 999	UNP P13332
B	151	ILE	TYR	SEE REMARK 999	UNP P13332
B	301	GLY	GLU	SEE REMARK 999	UNP P13332
B	399	VAL	ALA	SEE REMARK 999	UNP P13332
B	454	TYR	HIS	SEE REMARK 999	UNP P13332
B	510	PRO	ARG	engineered mutation	UNP P13332
C	100	GLU	ASP	SEE REMARK 999	UNP P13332
C	148	ALA	GLY	SEE REMARK 999	UNP P13332

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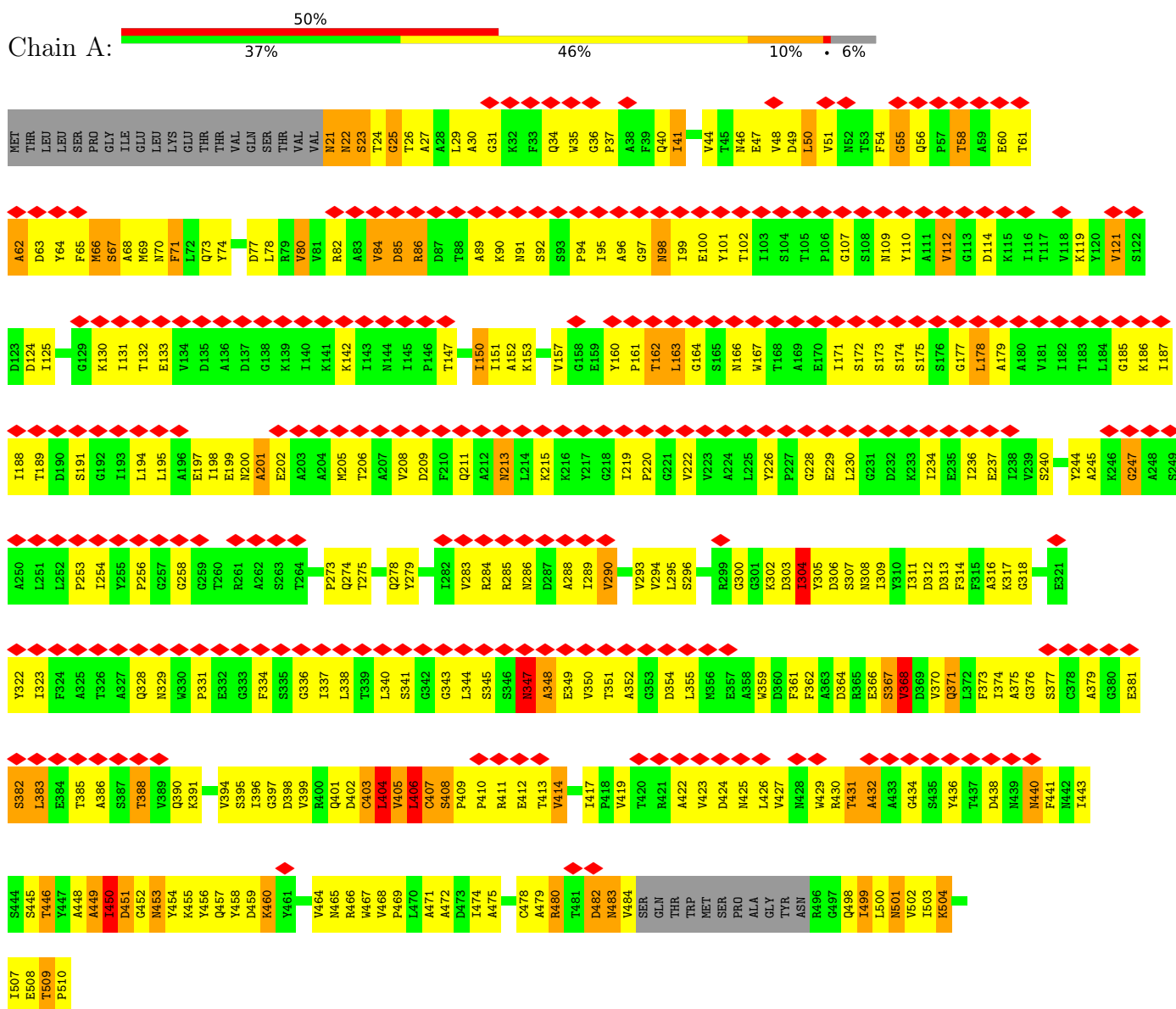
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Chain	Residue	Modelled	Actual	Comment	Reference
C	150	ILE	ASN	SEE REMARK 999	UNP P13332
C	151	ILE	TYR	SEE REMARK 999	UNP P13332
C	301	GLY	GLU	SEE REMARK 999	UNP P13332
C	399	VAL	ALA	SEE REMARK 999	UNP P13332
C	454	TYR	HIS	SEE REMARK 999	UNP P13332
C	510	PRO	ARG	engineered mutation	UNP P13332
D	100	GLU	ASP	SEE REMARK 999	UNP P13332
D	148	ALA	GLY	SEE REMARK 999	UNP P13332
D	150	ILE	ASN	SEE REMARK 999	UNP P13332
D	151	ILE	TYR	SEE REMARK 999	UNP P13332
D	301	GLY	GLU	SEE REMARK 999	UNP P13332
D	399	VAL	ALA	SEE REMARK 999	UNP P13332
D	454	TYR	HIS	SEE REMARK 999	UNP P13332
D	510	PRO	ARG	engineered mutation	UNP P13332
E	100	GLU	ASP	SEE REMARK 999	UNP P13332
E	148	ALA	GLY	SEE REMARK 999	UNP P13332
E	150	ILE	ASN	SEE REMARK 999	UNP P13332
E	151	ILE	TYR	SEE REMARK 999	UNP P13332
E	301	GLY	GLU	SEE REMARK 999	UNP P13332
E	399	VAL	ALA	SEE REMARK 999	UNP P13332
E	454	TYR	HIS	SEE REMARK 999	UNP P13332
E	510	PRO	ARG	engineered mutation	UNP P13332
F	100	GLU	ASP	SEE REMARK 999	UNP P13332
F	148	ALA	GLY	SEE REMARK 999	UNP P13332
F	150	ILE	ASN	SEE REMARK 999	UNP P13332
F	151	ILE	TYR	SEE REMARK 999	UNP P13332
F	301	GLY	GLU	SEE REMARK 999	UNP P13332
F	399	VAL	ALA	SEE REMARK 999	UNP P13332
F	454	TYR	HIS	SEE REMARK 999	UNP P13332
F	510	PRO	ARG	engineered mutation	UNP P13332

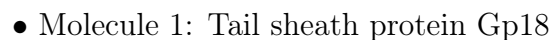
3 Residue-property plots

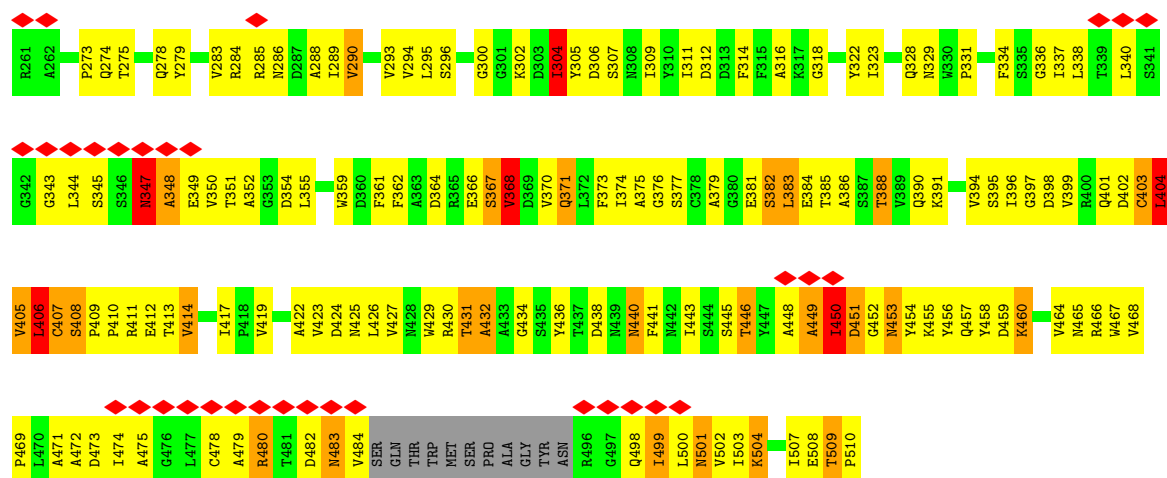
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion 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: Tail sheath protein Gp18

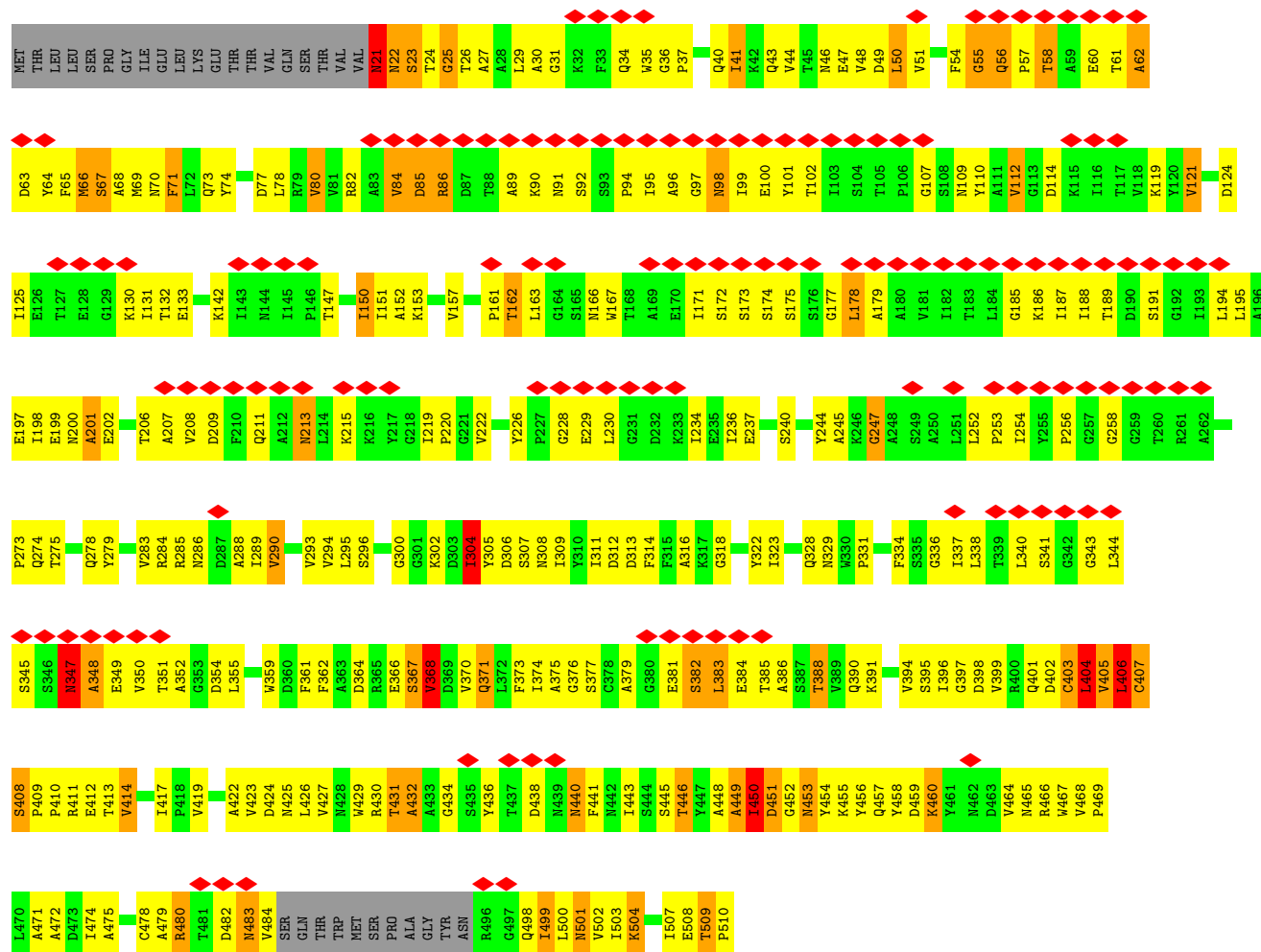


• Molecule 1: Tail sheath protein Gp18

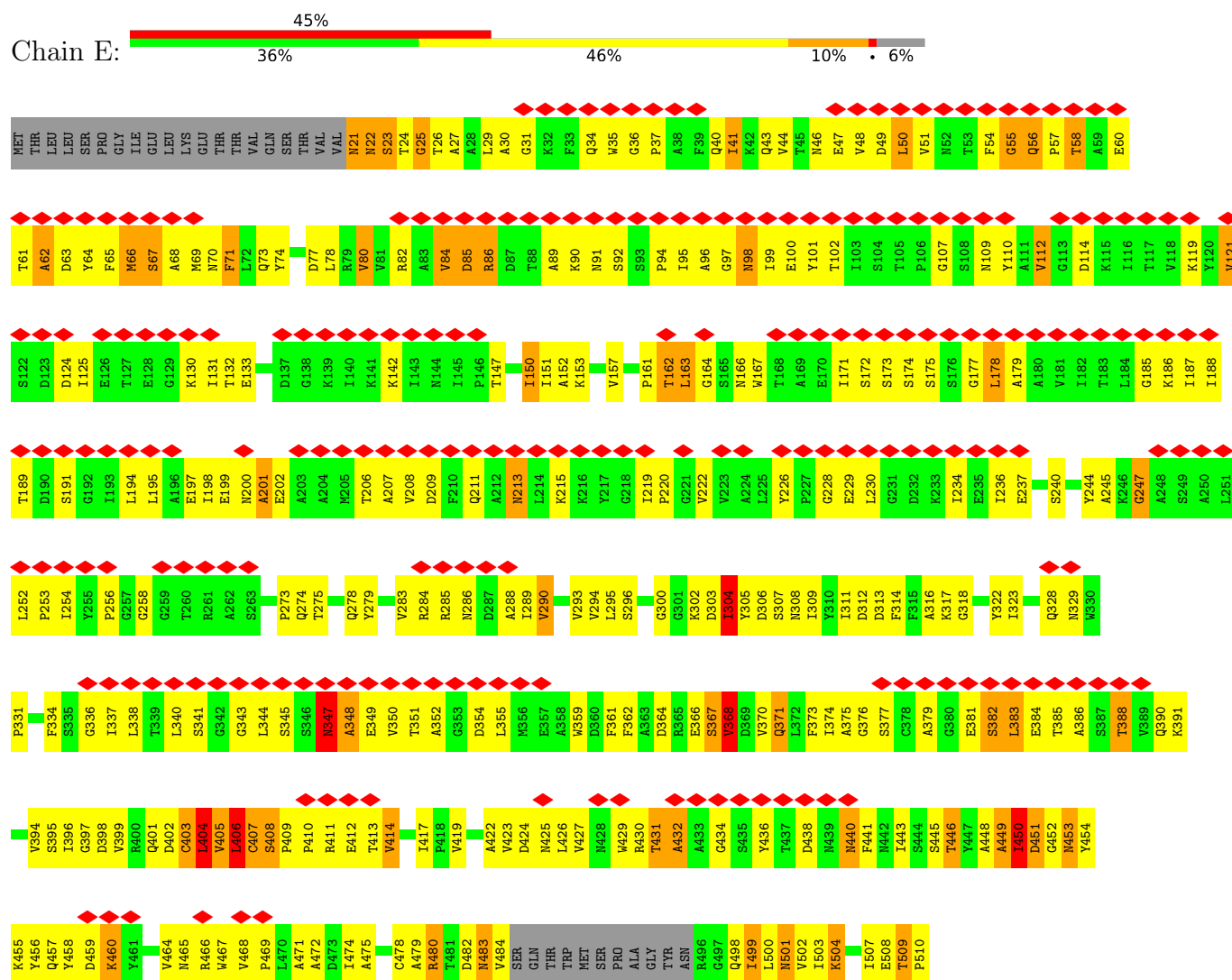




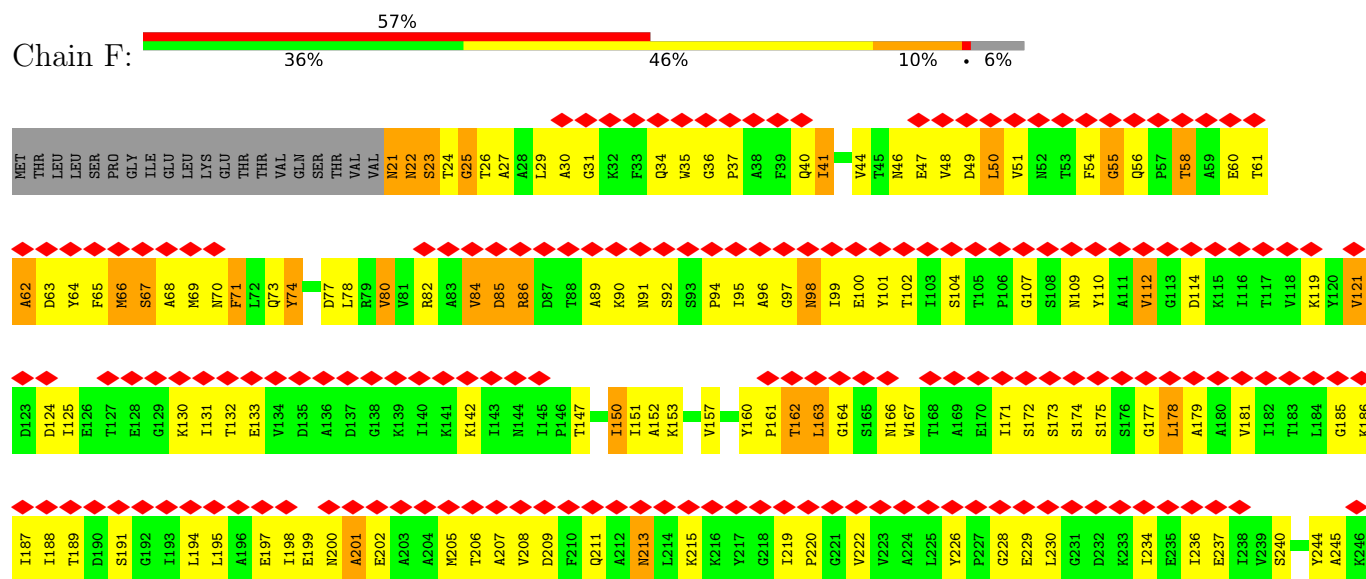
• Molecule 1: Tail sheath protein Gp18

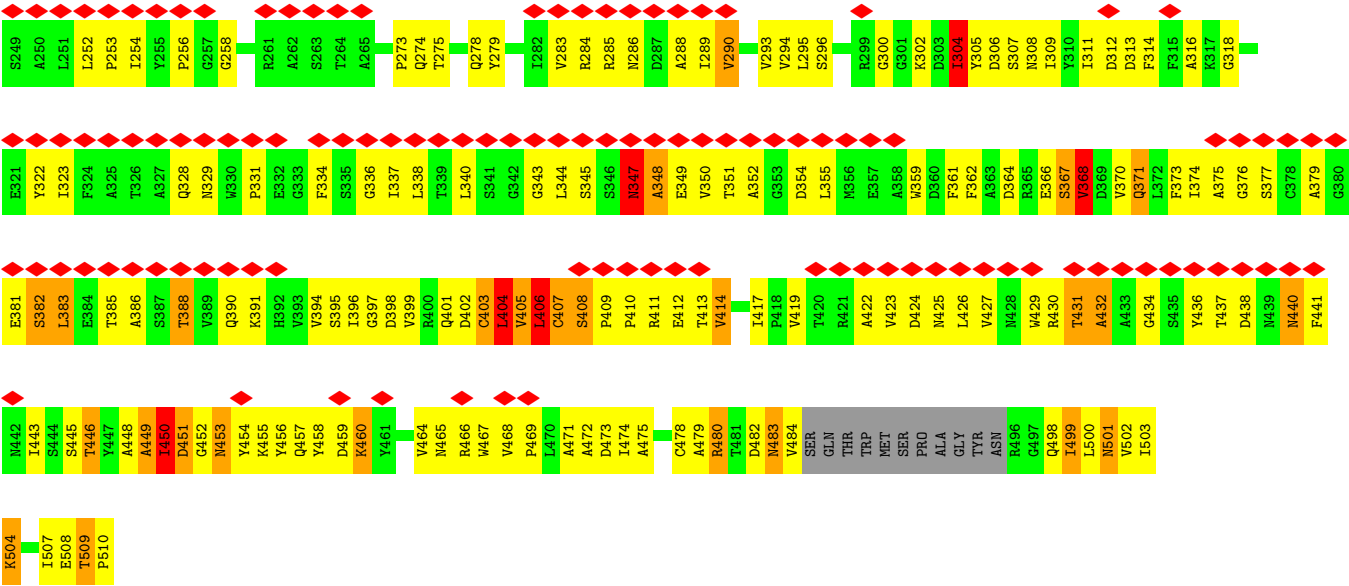


• Molecule 1: Tail sheath protein Gp18



• Molecule 1: Tail sheath protein Gp18





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C6	Depositor
Number of particles used	3029	Depositor
Resolution determination method	Not provided	
CTF correction method	Not provided	
Microscope	FEI/PHILIPS CM300FEG/ST	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	Not provided	
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	4000	Depositor
Magnification	45000	Depositor
Image detector	Not provided	
Maximum map value	7.953	Depositor
Minimum map value	-2.953	Depositor
Average map value	0.059	Depositor
Map value standard deviation	0.533	Depositor
Recommended contour level	1.01	Depositor
Map size ($\text{\AA}$)	714.6, 714.6, 1508.6	wwPDB
Map dimensions	180, 180, 380	wwPDB
Map angles ($^\circ$)	90, 90, 90	wwPDB
Pixel spacing ($\text{\AA}$)	3.97, 3.97, 3.97	Depositor

5 Model quality

5.1 Standard geometry

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.47	4/3677 (0.1%)	1.09	30/5001 (0.6%)
1	B	0.47	4/3677 (0.1%)	1.09	30/5001 (0.6%)
1	C	0.47	4/3677 (0.1%)	1.09	30/5001 (0.6%)
1	D	0.47	4/3677 (0.1%)	1.08	30/5001 (0.6%)
1	E	0.47	4/3677 (0.1%)	1.09	30/5001 (0.6%)
1	F	0.47	4/3677 (0.1%)	1.09	30/5001 (0.6%)
All	All	0.47	24/22062 (0.1%)	1.09	180/30006 (0.6%)

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
1	A	0	4
1	B	0	4
1	C	0	4
1	D	0	4
1	E	0	4
1	F	0	4
All	All	0	24

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	22	ASN	C-N	-7.25	1.23	1.33
1	E	22	ASN	C-N	-7.25	1.23	1.33
1	D	22	ASN	C-N	-7.19	1.23	1.33
1	C	22	ASN	C-N	-7.17	1.23	1.33
1	F	22	ASN	C-N	-7.17	1.23	1.33
1	A	22	ASN	C-N	-7.16	1.23	1.33
1	B	23	SER	C-N	-6.93	1.24	1.33
1	E	23	SER	C-N	-6.93	1.24	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	23	SER	C-N	-6.93	1.24	1.33
1	C	23	SER	C-N	-6.90	1.24	1.33
1	F	23	SER	C-N	-6.90	1.24	1.33
1	A	23	SER	C-N	-6.88	1.24	1.33
1	F	23	SER	C-O	6.85	1.32	1.24
1	B	23	SER	C-O	6.84	1.32	1.24
1	C	23	SER	C-O	6.84	1.32	1.24
1	E	23	SER	C-O	6.84	1.32	1.24
1	A	23	SER	C-O	6.81	1.32	1.24
1	D	23	SER	C-O	6.81	1.32	1.24
1	E	22	ASN	C-O	5.58	1.31	1.24
1	A	22	ASN	C-O	5.54	1.30	1.24
1	D	22	ASN	C-O	5.54	1.30	1.24
1	B	22	ASN	C-O	5.54	1.30	1.24
1	C	22	ASN	C-O	5.48	1.30	1.24
1	F	22	ASN	C-O	5.48	1.30	1.24

All (180) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	55	GLY	N-CA-C	18.97	158.14	113.18
1	F	55	GLY	N-CA-C	18.96	158.12	113.18
1	C	55	GLY	N-CA-C	18.96	158.12	113.18
1	D	55	GLY	N-CA-C	18.96	158.10	113.18
1	E	55	GLY	N-CA-C	18.95	158.10	113.18
1	B	55	GLY	N-CA-C	18.94	158.07	113.18
1	A	112	VAL	N-CA-C	17.42	136.71	112.50
1	B	112	VAL	N-CA-C	17.39	136.68	112.50
1	F	112	VAL	N-CA-C	17.39	136.67	112.50
1	C	112	VAL	N-CA-C	17.38	136.66	112.50
1	E	112	VAL	N-CA-C	17.36	136.63	112.50
1	D	112	VAL	N-CA-C	17.35	136.62	112.50
1	D	405	VAL	N-CA-C	-15.51	85.11	107.75
1	C	405	VAL	N-CA-C	-15.50	85.11	107.75
1	F	405	VAL	N-CA-C	-15.50	85.11	107.75
1	E	405	VAL	N-CA-C	-15.50	85.12	107.75
1	A	405	VAL	N-CA-C	-15.50	85.12	107.75
1	B	405	VAL	N-CA-C	-15.50	85.12	107.75
1	A	404	LEU	N-CA-C	-10.79	90.63	108.32
1	C	404	LEU	N-CA-C	-10.76	90.68	108.32
1	B	404	LEU	N-CA-C	-10.75	90.69	108.32
1	E	404	LEU	N-CA-C	-10.75	90.69	108.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	404	LEU	N-CA-C	-10.75	90.69	108.32
1	D	404	LEU	N-CA-C	-10.74	90.70	108.32
1	A	56	GLN	N-CA-CB	-10.63	90.12	110.10
1	F	56	GLN	N-CA-CB	-10.62	90.14	110.10
1	C	56	GLN	N-CA-CB	-10.61	90.16	110.10
1	D	56	GLN	N-CA-CB	-10.59	90.18	110.10
1	E	56	GLN	N-CA-CB	-10.59	90.19	110.10
1	B	56	GLN	N-CA-CB	-10.58	90.20	110.10
1	A	85	ASP	N-CA-C	-10.33	93.34	109.76
1	C	85	ASP	N-CA-C	-10.31	93.36	109.76
1	F	85	ASP	N-CA-C	-10.31	93.36	109.76
1	B	85	ASP	N-CA-C	-10.31	93.37	109.76
1	D	85	ASP	N-CA-C	-10.31	93.37	109.76
1	E	85	ASP	N-CA-C	-10.31	93.37	109.76
1	A	367	SER	N-CA-CB	-9.40	94.54	110.33
1	D	367	SER	N-CA-CB	-9.40	94.54	110.33
1	B	367	SER	N-CA-CB	-9.39	94.56	110.33
1	C	367	SER	N-CA-CB	-9.39	94.56	110.33
1	F	367	SER	N-CA-CB	-9.39	94.56	110.33
1	E	367	SER	N-CA-CB	-9.37	94.58	110.33
1	C	407	CYS	CB-CA-C	-9.23	91.40	109.68
1	A	407	CYS	CB-CA-C	-9.22	91.42	109.68
1	B	407	CYS	CB-CA-C	-9.22	91.42	109.68
1	E	407	CYS	CB-CA-C	-9.22	91.42	109.68
1	F	407	CYS	CB-CA-C	-9.21	91.44	109.68
1	F	366	GLU	CB-CA-C	8.73	125.28	110.79
1	E	366	GLU	CB-CA-C	8.72	125.27	110.79
1	C	366	GLU	CB-CA-C	8.71	125.26	110.79
1	B	366	GLU	CB-CA-C	8.71	125.24	110.79
1	D	366	GLU	CB-CA-C	8.70	125.24	110.79
1	D	407	CYS	CB-CA-C	-8.69	91.43	109.94
1	A	366	GLU	CB-CA-C	8.68	125.20	110.79
1	E	406	LEU	N-CA-C	-7.91	96.51	109.40
1	C	406	LEU	N-CA-C	-7.91	96.51	109.40
1	B	404	LEU	CB-CA-C	7.90	123.25	109.89
1	D	406	LEU	N-CA-C	-7.90	96.52	109.40
1	F	406	LEU	N-CA-C	-7.90	96.53	109.40
1	C	404	LEU	CB-CA-C	7.90	123.23	109.89
1	B	406	LEU	N-CA-C	-7.89	96.54	109.40
1	F	404	LEU	CB-CA-C	7.89	123.22	109.89
1	E	404	LEU	CB-CA-C	7.88	123.22	109.89
1	D	404	LEU	CB-CA-C	7.88	123.20	109.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	406	LEU	N-CA-C	-7.87	96.57	109.40
1	A	404	LEU	CB-CA-C	7.87	123.19	109.89
1	F	366	GLU	N-CA-C	-7.65	102.94	111.28
1	C	366	GLU	N-CA-C	-7.62	102.97	111.28
1	B	366	GLU	N-CA-C	-7.62	102.98	111.28
1	E	366	GLU	N-CA-C	-7.62	102.98	111.28
1	A	366	GLU	N-CA-C	-7.60	102.99	111.28
1	D	366	GLU	N-CA-C	-7.58	103.01	111.28
1	F	347	ASN	N-CA-C	-7.46	103.98	113.23
1	E	347	ASN	N-CA-C	-7.45	103.99	113.23
1	A	347	ASN	N-CA-C	-7.45	103.99	113.23
1	D	347	ASN	N-CA-C	-7.43	104.01	113.23
1	C	347	ASN	N-CA-C	-7.43	104.02	113.23
1	B	347	ASN	N-CA-C	-7.42	104.03	113.23
1	F	85	ASP	CB-CA-C	7.36	121.84	109.84
1	C	85	ASP	CB-CA-C	7.36	121.83	109.84
1	A	85	ASP	CB-CA-C	7.33	121.80	109.84
1	D	85	ASP	CB-CA-C	7.33	121.79	109.84
1	B	85	ASP	CB-CA-C	7.33	121.79	109.84
1	E	85	ASP	CB-CA-C	7.33	121.78	109.84
1	C	86	ARG	N-CA-C	-6.87	104.37	112.89
1	F	86	ARG	N-CA-C	-6.87	104.37	112.89
1	A	86	ARG	N-CA-C	-6.87	104.38	112.89
1	D	86	ARG	N-CA-C	-6.87	104.38	112.89
1	E	86	ARG	N-CA-C	-6.87	104.38	112.89
1	B	405	VAL	N-CA-CB	6.86	120.55	111.64
1	C	405	VAL	N-CA-CB	6.86	120.55	111.64
1	F	405	VAL	N-CA-CB	6.86	120.55	111.64
1	D	405	VAL	N-CA-CB	6.84	120.53	111.64
1	A	405	VAL	N-CA-CB	6.84	120.53	111.64
1	B	86	ARG	N-CA-C	-6.83	104.42	112.89
1	E	405	VAL	N-CA-CB	6.79	120.47	111.64
1	B	449	ALA	N-CA-C	-6.79	97.35	108.41
1	E	449	ALA	N-CA-C	-6.79	97.35	108.41
1	F	449	ALA	N-CA-C	-6.78	97.37	108.41
1	A	449	ALA	N-CA-C	-6.77	97.37	108.41
1	C	449	ALA	N-CA-C	-6.77	97.37	108.41
1	D	449	ALA	N-CA-C	-6.77	97.38	108.41
1	B	179	ALA	N-CA-C	6.76	125.21	110.80
1	A	179	ALA	N-CA-C	6.76	125.20	110.80
1	C	179	ALA	N-CA-C	6.76	125.20	110.80
1	F	179	ALA	N-CA-C	6.76	125.19	110.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	179	ALA	N-CA-C	6.74	125.15	110.80
1	D	179	ALA	N-CA-C	6.74	125.15	110.80
1	B	21	ASN	CA-C-N	6.73	134.39	121.54
1	B	21	ASN	C-N-CA	6.73	134.39	121.54
1	D	21	ASN	CA-C-N	6.72	134.38	121.54
1	D	21	ASN	C-N-CA	6.72	134.38	121.54
1	A	21	ASN	CA-C-N	6.70	134.33	121.54
1	A	21	ASN	C-N-CA	6.70	134.33	121.54
1	E	21	ASN	CA-C-N	6.70	134.33	121.54
1	E	21	ASN	C-N-CA	6.70	134.33	121.54
1	B	408	SER	N-CA-CB	-6.69	97.18	110.70
1	F	408	SER	N-CA-CB	-6.69	97.19	110.70
1	D	408	SER	N-CA-CB	-6.69	97.19	110.70
1	C	21	ASN	CA-C-N	6.68	134.30	121.54
1	C	21	ASN	C-N-CA	6.68	134.30	121.54
1	C	408	SER	N-CA-CB	-6.68	97.21	110.70
1	A	408	SER	N-CA-CB	-6.67	97.22	110.70
1	F	21	ASN	CA-C-N	6.67	134.28	121.54
1	F	21	ASN	C-N-CA	6.67	134.28	121.54
1	E	408	SER	N-CA-CB	-6.67	97.24	110.70
1	E	406	LEU	CB-CA-C	6.58	121.94	109.37
1	C	406	LEU	CB-CA-C	6.58	121.94	109.37
1	D	406	LEU	CB-CA-C	6.58	121.94	109.37
1	B	406	LEU	CB-CA-C	6.57	121.91	109.37
1	F	406	LEU	CB-CA-C	6.56	121.90	109.37
1	A	406	LEU	CB-CA-C	6.54	121.87	109.37
1	B	163	LEU	CB-CA-C	-6.21	107.70	115.89
1	D	163	LEU	CB-CA-C	-6.20	107.70	115.89
1	C	163	LEU	CB-CA-C	-6.20	107.71	115.89
1	F	163	LEU	CB-CA-C	-6.20	107.71	115.89
1	E	163	LEU	CB-CA-C	-6.19	107.72	115.89
1	B	178	LEU	N-CA-CB	-6.18	100.05	110.49
1	F	178	LEU	N-CA-CB	-6.17	100.06	110.49
1	A	178	LEU	N-CA-CB	-6.17	100.06	110.49
1	D	178	LEU	N-CA-CB	-6.17	100.06	110.49
1	A	163	LEU	CB-CA-C	-6.16	107.76	115.89
1	C	178	LEU	N-CA-CB	-6.16	100.08	110.49
1	E	178	LEU	N-CA-CB	-6.16	100.08	110.49
1	E	449	ALA	CB-CA-C	6.05	120.20	110.16
1	C	449	ALA	CB-CA-C	6.04	120.19	110.16
1	A	449	ALA	CB-CA-C	6.03	120.16	110.16
1	F	449	ALA	CB-CA-C	6.02	120.16	110.16

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	449	ALA	CB-CA-C	6.02	120.15	110.16
1	B	449	ALA	CB-CA-C	6.01	120.14	110.16
1	D	446	THR	N-CA-C	-5.83	106.01	113.01
1	B	446	THR	N-CA-C	-5.81	106.03	113.01
1	E	446	THR	N-CA-C	-5.80	106.05	113.01
1	C	446	THR	N-CA-C	-5.79	106.06	113.01
1	F	446	THR	N-CA-C	-5.76	106.10	113.01
1	A	446	THR	N-CA-C	-5.76	106.10	113.01
1	E	368	VAL	N-CA-C	5.39	115.62	107.75
1	F	368	VAL	N-CA-C	5.38	115.61	107.75
1	B	112	VAL	CB-CA-C	-5.38	96.92	111.34
1	B	368	VAL	N-CA-C	5.38	115.60	107.75
1	A	112	VAL	CB-CA-C	-5.37	96.94	111.34
1	D	112	VAL	CB-CA-C	-5.37	96.95	111.34
1	C	112	VAL	CB-CA-C	-5.37	96.96	111.34
1	F	112	VAL	CB-CA-C	-5.36	96.97	111.34
1	A	368	VAL	N-CA-C	5.36	115.57	107.75
1	E	112	VAL	CB-CA-C	-5.36	96.98	111.34
1	C	368	VAL	N-CA-C	5.34	115.55	107.75
1	D	368	VAL	N-CA-C	5.34	115.54	107.75
1	E	396	ILE	N-CA-C	-5.26	107.70	112.12
1	F	396	ILE	N-CA-C	-5.21	107.74	112.12
1	B	396	ILE	N-CA-C	-5.19	107.76	112.12
1	A	396	ILE	N-CA-C	-5.18	107.77	112.12
1	C	396	ILE	N-CA-C	-5.18	107.77	112.12
1	D	396	ILE	N-CA-C	-5.17	107.77	112.12
1	F	367	SER	N-CA-C	5.16	119.55	112.68
1	B	367	SER	N-CA-C	5.16	119.54	112.68
1	A	367	SER	N-CA-C	5.15	119.53	112.68
1	D	367	SER	N-CA-C	5.15	119.53	112.68
1	E	367	SER	N-CA-C	5.15	119.53	112.68
1	C	367	SER	N-CA-C	5.14	119.52	112.68

There are no chirality outliers.

All (24) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	21	ASN	Peptide
1	A	450	ILE	Peptide
1	A	451	ASP	Peptide
1	A	452	GLY	Peptide
1	B	21	ASN	Peptide

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Mol	Chain	Res	Type	Group
1	B	450	ILE	Peptide
1	B	451	ASP	Peptide
1	B	452	GLY	Peptide
1	C	21	ASN	Peptide
1	C	450	ILE	Peptide
1	C	451	ASP	Peptide
1	C	452	GLY	Peptide
1	D	21	ASN	Peptide
1	D	450	ILE	Peptide
1	D	451	ASP	Peptide
1	D	452	GLY	Peptide
1	E	21	ASN	Peptide
1	E	450	ILE	Peptide
1	E	451	ASP	Peptide
1	E	452	GLY	Peptide
1	F	21	ASN	Peptide
1	F	450	ILE	Peptide
1	F	451	ASP	Peptide
1	F	452	GLY	Peptide

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3613	0	3554	320	0
1	B	3613	0	3554	319	0
1	C	3613	0	3554	318	0
1	D	3613	0	3554	321	0
1	E	3613	0	3554	320	0
1	F	3613	0	3554	319	0
All	All	21678	0	21324	1917	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 45.

All (1917) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:409:PRO:O	1:E:454:TYR:CE1	1.65	1.50
1:F:409:PRO:O	1:F:454:TYR:CE1	1.65	1.49
1:D:409:PRO:O	1:D:454:TYR:CE1	1.65	1.48
1:A:409:PRO:O	1:A:454:TYR:CE1	1.65	1.48
1:C:409:PRO:O	1:C:454:TYR:CE1	1.65	1.47
1:B:409:PRO:O	1:B:454:TYR:CE1	1.65	1.47
1:E:409:PRO:CD	1:E:451:ASP:O	1.75	1.34
1:F:409:PRO:CD	1:F:451:ASP:O	1.75	1.34
1:D:409:PRO:CD	1:D:451:ASP:O	1.75	1.34
1:C:409:PRO:CD	1:C:451:ASP:O	1.75	1.33
1:A:409:PRO:CD	1:A:451:ASP:O	1.75	1.33
1:B:409:PRO:CD	1:B:451:ASP:O	1.75	1.32
1:D:379:ALA:CB	1:D:454:TYR:OH	1.84	1.25
1:C:379:ALA:CB	1:C:454:TYR:OH	1.84	1.25
1:B:379:ALA:CB	1:B:454:TYR:OH	1.84	1.25
1:E:379:ALA:CB	1:E:454:TYR:OH	1.84	1.24
1:F:379:ALA:CB	1:F:454:TYR:OH	1.84	1.24
1:A:379:ALA:CB	1:A:454:TYR:OH	1.84	1.24
1:E:379:ALA:HB2	1:E:454:TYR:OH	1.37	1.23
1:D:379:ALA:HB2	1:D:454:TYR:OH	1.37	1.22
1:C:379:ALA:HB2	1:C:454:TYR:OH	1.37	1.21
1:F:379:ALA:HB2	1:F:454:TYR:OH	1.37	1.20
1:A:379:ALA:HB2	1:A:454:TYR:OH	1.37	1.18
1:B:379:ALA:HB2	1:B:454:TYR:OH	1.37	1.17
1:C:409:PRO:HD3	1:C:451:ASP:O	0.98	1.15
1:B:409:PRO:HD3	1:B:451:ASP:O	0.98	1.15
1:E:409:PRO:HD3	1:E:451:ASP:O	0.98	1.14
1:A:409:PRO:HD3	1:A:451:ASP:O	0.98	1.14
1:F:409:PRO:HD3	1:F:451:ASP:O	0.98	1.14
1:D:409:PRO:HD3	1:D:451:ASP:O	0.98	1.14
1:D:454:TYR:CE2	1:D:469:PRO:HA	1.84	1.13
1:E:454:TYR:CE2	1:E:469:PRO:HA	1.84	1.13
1:A:454:TYR:CE2	1:A:469:PRO:HA	1.84	1.13
1:F:379:ALA:HB1	1:F:454:TYR:CZ	1.83	1.13
1:A:379:ALA:HB1	1:A:454:TYR:CZ	1.83	1.12
1:F:454:TYR:CE2	1:F:469:PRO:HA	1.84	1.13
1:C:454:TYR:CE2	1:C:469:PRO:HA	1.84	1.12
1:D:379:ALA:HB1	1:D:454:TYR:CZ	1.83	1.12
1:B:454:TYR:CE2	1:B:469:PRO:HA	1.84	1.12
1:E:379:ALA:HB1	1:E:454:TYR:CZ	1.83	1.12
1:B:379:ALA:HB1	1:B:454:TYR:CZ	1.83	1.11
1:C:379:ALA:HB1	1:C:454:TYR:CZ	1.83	1.11

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:379:ALA:CB	1:A:454:TYR:CZ	2.35	1.10
1:F:379:ALA:CB	1:F:454:TYR:CZ	2.35	1.10
1:B:379:ALA:CB	1:B:454:TYR:CZ	2.35	1.09
1:D:409:PRO:O	1:D:454:TYR:CZ	2.06	1.09
1:F:409:PRO:O	1:F:454:TYR:CZ	2.06	1.09
1:A:409:PRO:O	1:A:454:TYR:CZ	2.06	1.09
1:B:409:PRO:O	1:B:454:TYR:CZ	2.06	1.09
1:E:379:ALA:CB	1:E:454:TYR:CZ	2.35	1.08
1:D:379:ALA:CB	1:D:454:TYR:CZ	2.35	1.08
1:B:51:VAL:HA	1:B:55:GLY:HA2	1.36	1.08
1:E:409:PRO:O	1:E:454:TYR:CZ	2.06	1.08
1:C:379:ALA:CB	1:C:454:TYR:CZ	2.35	1.08
1:B:450:ILE:HG12	1:B:451:ASP:H	0.94	1.07
1:A:450:ILE:HG12	1:A:451:ASP:H	0.94	1.07
1:A:454:TYR:CD2	1:A:469:PRO:HA	1.89	1.07
1:C:409:PRO:O	1:C:454:TYR:CZ	2.06	1.07
1:A:51:VAL:HA	1:A:55:GLY:HA2	1.36	1.07
1:E:454:TYR:CD2	1:E:469:PRO:HA	1.89	1.07
1:F:454:TYR:CD2	1:F:469:PRO:HA	1.89	1.07
1:B:454:TYR:CD2	1:B:469:PRO:HA	1.89	1.07
1:C:51:VAL:HA	1:C:55:GLY:HA2	1.36	1.06
1:F:51:VAL:HA	1:F:55:GLY:HA2	1.36	1.06
1:C:450:ILE:HG12	1:C:451:ASP:H	0.94	1.06
1:D:454:TYR:CD2	1:D:469:PRO:HA	1.89	1.06
1:C:454:TYR:CD2	1:C:469:PRO:HA	1.89	1.06
1:F:450:ILE:HG12	1:F:451:ASP:H	0.94	1.05
1:E:51:VAL:HA	1:E:55:GLY:HA2	1.36	1.05
1:D:450:ILE:HG12	1:D:451:ASP:H	0.94	1.04
1:E:450:ILE:HG12	1:E:451:ASP:N	1.67	1.04
1:D:51:VAL:HA	1:D:55:GLY:HA2	1.36	1.04
1:E:450:ILE:HG12	1:E:451:ASP:H	0.94	1.03
1:D:450:ILE:HG12	1:D:451:ASP:N	1.67	1.01
1:C:450:ILE:HG12	1:C:451:ASP:N	1.67	0.99
1:A:450:ILE:HG12	1:A:451:ASP:N	1.67	0.98
1:B:450:ILE:HG12	1:B:451:ASP:N	1.67	0.98
1:F:450:ILE:HG12	1:F:451:ASP:N	1.67	0.98
1:D:409:PRO:O	1:D:454:TYR:HE1	1.16	0.97
1:C:228:GLY:HA2	1:C:345:SER:HB3	1.47	0.97
1:D:228:GLY:HA2	1:D:345:SER:HB3	1.47	0.96
1:B:228:GLY:HA2	1:B:345:SER:HB3	1.47	0.96
1:E:112:VAL:CG1	1:E:112:VAL:O	2.14	0.96

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:228:GLY:HA2	1:E:345:SER:HB3	1.47	0.95
1:D:112:VAL:O	1:D:112:VAL:CG1	2.14	0.95
1:B:112:VAL:O	1:B:112:VAL:CG1	2.14	0.95
1:C:112:VAL:O	1:C:112:VAL:CG1	2.14	0.95
1:A:112:VAL:O	1:A:112:VAL:CG1	2.14	0.94
1:F:112:VAL:O	1:F:112:VAL:CG1	2.14	0.94
1:A:228:GLY:HA2	1:A:345:SER:HB3	1.47	0.93
1:F:228:GLY:HA2	1:F:345:SER:HB3	1.47	0.93
1:E:409:PRO:O	1:E:454:TYR:HE1	1.16	0.93
1:B:409:PRO:O	1:B:454:TYR:HE1	1.16	0.92
1:F:408:SER:CA	1:F:451:ASP:HB3	2.00	0.92
1:F:409:PRO:O	1:F:454:TYR:HE1	1.15	0.92
1:A:408:SER:CA	1:A:451:ASP:HB3	2.00	0.92
1:D:408:SER:CA	1:D:451:ASP:HB3	2.00	0.92
1:B:408:SER:CA	1:B:451:ASP:HB3	2.00	0.92
1:E:408:SER:CA	1:E:451:ASP:HB3	2.00	0.91
1:C:408:SER:CA	1:C:451:ASP:HB3	2.00	0.91
1:C:23:SER:OG	1:C:483:ASN:CB	2.19	0.91
1:D:23:SER:OG	1:D:483:ASN:CB	2.19	0.91
1:C:409:PRO:O	1:C:454:TYR:HE1	1.15	0.91
1:B:23:SER:OG	1:B:483:ASN:CB	2.19	0.90
1:A:23:SER:OG	1:A:483:ASN:CB	2.19	0.90
1:E:23:SER:OG	1:E:483:ASN:CB	2.19	0.90
1:F:23:SER:OG	1:F:483:ASN:CB	2.19	0.90
1:A:409:PRO:O	1:A:454:TYR:HE1	1.16	0.89
1:E:408:SER:HA	1:E:451:ASP:HB3	1.54	0.89
1:C:508:GLU:HG2	1:C:509:THR:HG22	1.55	0.89
1:E:508:GLU:HG2	1:E:509:THR:HG22	1.55	0.89
1:B:508:GLU:HG2	1:B:509:THR:HG22	1.55	0.88
1:F:508:GLU:HG2	1:F:509:THR:HG22	1.55	0.88
1:C:408:SER:HA	1:C:451:ASP:HB3	1.54	0.88
1:D:391:LYS:HZ3	1:D:440:ASN:HD21	1.20	0.88
1:D:508:GLU:HG2	1:D:509:THR:HG22	1.55	0.88
1:A:508:GLU:HG2	1:A:509:THR:HG22	1.55	0.88
1:B:408:SER:HA	1:B:451:ASP:HB3	1.54	0.88
1:D:408:SER:HA	1:D:451:ASP:HB3	1.54	0.88
1:F:408:SER:HA	1:F:451:ASP:HB3	1.54	0.88
1:F:23:SER:CB	1:F:483:ASN:HB3	2.04	0.87
1:A:41:ILE:HD11	1:A:361:PHE:HB3	1.57	0.87
1:A:23:SER:CB	1:A:483:ASN:HB3	2.04	0.87
1:C:41:ILE:HD11	1:C:361:PHE:HB3	1.57	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:408:SER:HA	1:A:451:ASP:HB3	1.54	0.87
1:D:391:LYS:NZ	1:D:440:ASN:ND2	2.23	0.87
1:E:23:SER:CB	1:E:483:ASN:HB3	2.04	0.87
1:A:509:THR:OG1	1:A:510:PRO:OXT	1.93	0.86
1:B:41:ILE:HD11	1:B:361:PHE:HB3	1.57	0.86
1:C:23:SER:CB	1:C:483:ASN:HB3	2.04	0.86
1:B:23:SER:CB	1:B:483:ASN:HB3	2.04	0.86
1:C:391:LYS:NZ	1:C:440:ASN:ND2	2.23	0.86
1:C:23:SER:OG	1:C:483:ASN:HB3	1.76	0.86
1:D:23:SER:CB	1:D:483:ASN:HB3	2.04	0.86
1:D:41:ILE:HD11	1:D:361:PHE:HB3	1.57	0.86
1:F:509:THR:OG1	1:F:510:PRO:OXT	1.93	0.86
1:D:391:LYS:HE3	1:D:440:ASN:O	1.75	0.86
1:E:391:LYS:NZ	1:E:440:ASN:ND2	2.23	0.86
1:F:41:ILE:HD11	1:F:361:PHE:HB3	1.57	0.86
1:C:391:LYS:HE3	1:C:440:ASN:O	1.76	0.86
1:D:23:SER:OG	1:D:483:ASN:HB3	1.76	0.85
1:F:391:LYS:HE3	1:F:440:ASN:O	1.75	0.85
1:E:23:SER:OG	1:E:483:ASN:HB3	1.76	0.85
1:A:391:LYS:NZ	1:A:440:ASN:ND2	2.23	0.85
1:B:391:LYS:NZ	1:B:440:ASN:ND2	2.23	0.85
1:D:509:THR:OG1	1:D:510:PRO:OXT	1.93	0.85
1:F:391:LYS:NZ	1:F:440:ASN:ND2	2.23	0.85
1:C:509:THR:OG1	1:C:510:PRO:OXT	1.93	0.85
1:F:23:SER:OG	1:F:483:ASN:HB3	1.76	0.85
1:F:23:SER:CB	1:F:483:ASN:CB	2.55	0.85
1:A:391:LYS:HE3	1:A:440:ASN:O	1.76	0.84
1:B:391:LYS:HE3	1:B:440:ASN:O	1.76	0.84
1:B:509:THR:OG1	1:B:510:PRO:OXT	1.93	0.84
1:A:23:SER:CB	1:A:483:ASN:CB	2.55	0.84
1:B:23:SER:OG	1:B:483:ASN:HB3	1.76	0.84
1:E:41:ILE:HD11	1:E:361:PHE:HB3	1.57	0.84
1:E:391:LYS:HE3	1:E:440:ASN:O	1.76	0.84
1:D:23:SER:CB	1:D:483:ASN:CB	2.55	0.84
1:E:509:THR:OG1	1:E:510:PRO:OXT	1.93	0.84
1:E:23:SER:CB	1:E:483:ASN:CB	2.55	0.84
1:C:23:SER:CB	1:C:483:ASN:CB	2.55	0.84
1:C:454:TYR:HE2	1:C:469:PRO:HA	1.43	0.84
1:D:331:PRO:HB2	1:D:334:PHE:HB2	1.60	0.83
1:B:23:SER:CB	1:B:483:ASN:CB	2.55	0.83
1:D:454:TYR:HE2	1:D:469:PRO:HA	1.43	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:331:PRO:HB2	1:C:334:PHE:HB2	1.60	0.83
1:F:331:PRO:HB2	1:F:334:PHE:HB2	1.60	0.83
1:B:454:TYR:HE2	1:B:469:PRO:HA	1.43	0.83
1:C:94:PRO:HB2	1:C:219:ILE:HD12	1.61	0.83
1:A:94:PRO:HB2	1:A:219:ILE:HD12	1.61	0.83
1:E:331:PRO:HB2	1:E:334:PHE:HB2	1.60	0.83
1:F:94:PRO:HB2	1:F:219:ILE:HD12	1.61	0.83
1:A:454:TYR:HE2	1:A:469:PRO:HA	1.43	0.83
1:A:331:PRO:HB2	1:A:334:PHE:HB2	1.60	0.83
1:B:331:PRO:HB2	1:B:334:PHE:HB2	1.60	0.83
1:A:23:SER:OG	1:A:483:ASN:HB3	1.76	0.82
1:C:391:LYS:HZ1	1:C:440:ASN:HD21	1.25	0.82
1:D:94:PRO:HB2	1:D:219:ILE:HD12	1.61	0.82
1:E:379:ALA:HB1	1:E:454:TYR:OH	1.73	0.81
1:B:94:PRO:HB2	1:B:219:ILE:HD12	1.61	0.81
1:E:110:TYR:CE1	1:E:178:LEU:O	2.33	0.81
1:F:110:TYR:CE1	1:F:178:LEU:O	2.33	0.81
1:E:94:PRO:HB2	1:E:219:ILE:HD12	1.61	0.81
1:D:110:TYR:CE1	1:D:178:LEU:O	2.33	0.81
1:A:110:TYR:CE1	1:A:178:LEU:O	2.33	0.81
1:B:110:TYR:CE1	1:B:178:LEU:O	2.33	0.81
1:C:110:TYR:CE1	1:C:178:LEU:O	2.33	0.80
1:D:112:VAL:O	1:D:112:VAL:HG12	1.82	0.80
1:A:23:SER:HB3	1:A:483:ASN:HB3	1.64	0.80
1:B:23:SER:HB3	1:B:483:ASN:HB3	1.64	0.80
1:C:379:ALA:HB1	1:C:454:TYR:OH	1.73	0.80
1:E:23:SER:HB3	1:E:483:ASN:HB3	1.64	0.80
1:F:454:TYR:HE2	1:F:469:PRO:HA	1.43	0.80
1:A:112:VAL:O	1:A:112:VAL:HG12	1.82	0.80
1:C:112:VAL:O	1:C:112:VAL:HG12	1.82	0.80
1:E:112:VAL:O	1:E:112:VAL:HG12	1.82	0.79
1:F:112:VAL:O	1:F:112:VAL:HG12	1.82	0.79
1:F:178:LEU:HD23	1:F:178:LEU:H	1.48	0.79
1:D:23:SER:HB3	1:D:483:ASN:HB3	1.64	0.79
1:B:178:LEU:HD23	1:B:178:LEU:H	1.48	0.79
1:C:23:SER:HB3	1:C:483:ASN:HB3	1.64	0.79
1:C:379:ALA:HB2	1:C:454:TYR:CZ	2.11	0.79
1:E:178:LEU:HD23	1:E:178:LEU:H	1.48	0.79
1:F:479:ALA:HA	1:F:484:VAL:HG11	1.65	0.78
1:A:479:ALA:HA	1:A:484:VAL:HG11	1.65	0.78
1:C:509:THR:OG1	1:C:510:PRO:CA	2.31	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:509:THR:OG1	1:D:510:PRO:CA	2.31	0.78
1:E:479:ALA:HA	1:E:484:VAL:HG11	1.65	0.78
1:F:23:SER:HB3	1:F:483:ASN:HB3	1.64	0.78
1:B:509:THR:OG1	1:B:510:PRO:CA	2.31	0.78
1:D:391:LYS:HZ3	1:D:440:ASN:ND2	1.79	0.78
1:B:112:VAL:O	1:B:112:VAL:HG12	1.82	0.78
1:D:479:ALA:HA	1:D:484:VAL:HG11	1.65	0.78
1:D:178:LEU:HD23	1:D:178:LEU:H	1.48	0.78
1:E:509:THR:OG1	1:E:510:PRO:CA	2.32	0.78
1:A:509:THR:OG1	1:A:510:PRO:CA	2.32	0.77
1:E:407:CYS:O	1:E:451:ASP:HB2	1.85	0.77
1:A:407:CYS:O	1:A:451:ASP:CB	2.32	0.77
1:B:479:ALA:HA	1:B:484:VAL:HG11	1.65	0.77
1:C:84:VAL:HG13	1:C:85:ASP:O	1.85	0.77
1:C:178:LEU:HD23	1:C:178:LEU:H	1.48	0.77
1:F:509:THR:OG1	1:F:510:PRO:CA	2.32	0.77
1:D:407:CYS:O	1:D:451:ASP:HB2	1.85	0.77
1:D:172:SER:C	1:D:174:SER:HA	2.09	0.77
1:F:172:SER:C	1:F:174:SER:HA	2.09	0.77
1:B:84:VAL:HG13	1:B:85:ASP:O	1.85	0.77
1:C:172:SER:C	1:C:174:SER:HA	2.09	0.77
1:C:407:CYS:O	1:C:451:ASP:CB	2.33	0.77
1:E:454:TYR:HE2	1:E:469:PRO:HA	1.43	0.77
1:F:84:VAL:HG13	1:F:85:ASP:O	1.85	0.77
1:F:407:CYS:O	1:F:451:ASP:HB2	1.85	0.77
1:B:172:SER:C	1:B:174:SER:HA	2.09	0.77
1:F:178:LEU:H	1:F:178:LEU:CD2	1.98	0.77
1:A:84:VAL:HG13	1:A:85:ASP:O	1.85	0.77
1:A:172:SER:C	1:A:174:SER:HA	2.09	0.77
1:B:407:CYS:O	1:B:451:ASP:CB	2.32	0.77
1:C:479:ALA:HA	1:C:484:VAL:HG11	1.65	0.77
1:E:172:SER:C	1:E:174:SER:HA	2.09	0.77
1:B:178:LEU:H	1:B:178:LEU:CD2	1.98	0.77
1:C:178:LEU:H	1:C:178:LEU:CD2	1.98	0.77
1:F:407:CYS:O	1:F:451:ASP:CB	2.32	0.77
1:F:171:ILE:HG22	1:F:172:SER:H	1.50	0.76
1:A:171:ILE:HG22	1:A:172:SER:H	1.50	0.76
1:D:407:CYS:O	1:D:451:ASP:CB	2.32	0.76
1:A:178:LEU:H	1:A:178:LEU:CD2	1.98	0.76
1:C:451:ASP:OD1	1:C:474:ILE:HD13	1.86	0.76
1:D:451:ASP:OD1	1:D:474:ILE:HD13	1.86	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:171:ILE:HG22	1:E:172:SER:H	1.50	0.76
1:A:178:LEU:H	1:A:178:LEU:HD23	1.48	0.76
1:C:407:CYS:O	1:C:451:ASP:HB2	1.85	0.76
1:D:84:VAL:HG13	1:D:85:ASP:O	1.85	0.76
1:E:178:LEU:H	1:E:178:LEU:CD2	1.98	0.76
1:E:407:CYS:O	1:E:451:ASP:CB	2.32	0.76
1:E:451:ASP:OD1	1:E:474:ILE:HD13	1.86	0.76
1:F:451:ASP:OD1	1:F:474:ILE:HD13	1.86	0.76
1:D:178:LEU:H	1:D:178:LEU:CD2	1.98	0.76
1:B:454:TYR:CE2	1:B:469:PRO:CA	2.67	0.76
1:C:454:TYR:CE2	1:C:469:PRO:CA	2.67	0.76
1:B:112:VAL:O	1:B:131:ILE:O	2.04	0.76
1:B:407:CYS:O	1:B:451:ASP:HB2	1.84	0.76
1:D:379:ALA:HB1	1:D:454:TYR:OH	1.73	0.76
1:C:112:VAL:O	1:C:131:ILE:O	2.04	0.76
1:D:112:VAL:O	1:D:131:ILE:O	2.04	0.76
1:E:110:TYR:HE1	1:E:178:LEU:O	1.69	0.76
1:A:112:VAL:O	1:A:131:ILE:O	2.04	0.75
1:A:407:CYS:O	1:A:451:ASP:HB2	1.85	0.75
1:B:171:ILE:HG22	1:B:172:SER:H	1.50	0.75
1:A:379:ALA:HB2	1:A:454:TYR:CZ	2.11	0.75
1:A:451:ASP:OD1	1:A:474:ILE:HD13	1.86	0.75
1:A:454:TYR:CE2	1:A:469:PRO:CA	2.67	0.75
1:E:24:THR:O	1:E:371:GLN:OE1	2.04	0.75
1:F:24:THR:O	1:F:371:GLN:OE1	2.04	0.75
1:B:451:ASP:OD1	1:B:474:ILE:HD13	1.86	0.75
1:D:24:THR:O	1:D:371:GLN:OE1	2.04	0.75
1:D:454:TYR:CE2	1:D:469:PRO:CA	2.67	0.75
1:E:112:VAL:O	1:E:131:ILE:O	2.04	0.75
1:F:509:THR:OG1	1:F:510:PRO:HA	1.87	0.75
1:D:110:TYR:HE1	1:D:178:LEU:O	1.69	0.75
1:E:84:VAL:HG13	1:E:85:ASP:O	1.85	0.75
1:A:22:ASN:O	1:A:23:SER:OG	2.04	0.75
1:B:24:THR:O	1:B:371:GLN:OE1	2.04	0.75
1:C:509:THR:OG1	1:C:510:PRO:HA	1.87	0.75
1:A:509:THR:OG1	1:A:510:PRO:HA	1.87	0.75
1:B:509:THR:OG1	1:B:510:PRO:HA	1.87	0.75
1:C:171:ILE:HG22	1:C:172:SER:H	1.50	0.75
1:D:379:ALA:HB2	1:D:454:TYR:CZ	2.11	0.75
1:C:24:THR:O	1:C:371:GLN:OE1	2.04	0.74
1:D:171:ILE:HG22	1:D:172:SER:H	1.50	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:112:VAL:O	1:F:131:ILE:O	2.04	0.74
1:A:26:THR:HA	1:A:77:ASP:OD1	1.87	0.74
1:A:509:THR:H	1:A:510:PRO:HA	1.52	0.74
1:E:26:THR:HA	1:E:77:ASP:OD1	1.87	0.74
1:E:454:TYR:CE2	1:E:469:PRO:CA	2.67	0.74
1:F:110:TYR:HE1	1:F:178:LEU:O	1.69	0.74
1:F:454:TYR:CE2	1:F:469:PRO:CA	2.67	0.74
1:A:24:THR:O	1:A:371:GLN:OE1	2.04	0.74
1:E:509:THR:OG1	1:E:510:PRO:HA	1.87	0.74
1:B:26:THR:HA	1:B:77:ASP:OD1	1.87	0.74
1:A:110:TYR:HE1	1:A:178:LEU:O	1.69	0.74
1:D:26:THR:HA	1:D:77:ASP:OD1	1.87	0.74
1:D:509:THR:OG1	1:D:510:PRO:HA	1.87	0.74
1:D:509:THR:H	1:D:510:PRO:HA	1.52	0.74
1:F:26:THR:HA	1:F:77:ASP:OD1	1.87	0.74
1:F:379:ALA:HB2	1:F:454:TYR:CZ	2.11	0.74
1:F:409:PRO:O	1:F:454:TYR:OH	2.06	0.74
1:E:409:PRO:O	1:E:454:TYR:OH	2.06	0.74
1:F:391:LYS:HZ1	1:F:440:ASN:HD21	1.34	0.74
1:B:450:ILE:CG1	1:B:451:ASP:H	1.89	0.74
1:E:112:VAL:O	1:E:112:VAL:HG13	1.88	0.74
1:A:409:PRO:O	1:A:454:TYR:OH	2.06	0.73
1:B:110:TYR:HE1	1:B:178:LEU:O	1.69	0.73
1:B:391:LYS:HZ1	1:B:440:ASN:HD21	1.33	0.73
1:E:509:THR:H	1:E:510:PRO:HA	1.52	0.73
1:F:22:ASN:O	1:F:23:SER:OG	2.04	0.73
1:C:26:THR:HA	1:C:77:ASP:OD1	1.87	0.73
1:A:509:THR:OG1	1:A:510:PRO:C	2.32	0.73
1:C:509:THR:OG1	1:C:510:PRO:C	2.32	0.73
1:B:509:THR:H	1:B:510:PRO:HA	1.53	0.73
1:C:509:THR:H	1:C:510:PRO:HA	1.52	0.73
1:D:112:VAL:O	1:D:112:VAL:HG13	1.88	0.73
1:B:409:PRO:O	1:B:454:TYR:OH	2.06	0.73
1:F:454:TYR:HD2	1:F:469:PRO:HA	1.53	0.73
1:F:509:THR:H	1:F:510:PRO:HA	1.53	0.73
1:A:112:VAL:O	1:A:112:VAL:HG13	1.88	0.73
1:D:409:PRO:HD2	1:D:451:ASP:O	1.88	0.73
1:E:379:ALA:HB2	1:E:454:TYR:CZ	2.11	0.73
1:F:51:VAL:HG13	1:F:55:GLY:O	1.89	0.73
1:C:110:TYR:HE1	1:C:178:LEU:O	1.69	0.72
1:B:112:VAL:O	1:B:112:VAL:HG13	1.88	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:410:PRO:O	1:B:413:THR:HG22	1.90	0.72
1:D:22:ASN:O	1:D:23:SER:OG	2.04	0.72
1:D:509:THR:OG1	1:D:510:PRO:C	2.32	0.72
1:E:51:VAL:HG13	1:E:55:GLY:O	1.89	0.72
1:B:382:SER:HB2	1:B:385:THR:HG22	1.72	0.72
1:D:409:PRO:O	1:D:454:TYR:OH	2.06	0.72
1:D:410:PRO:O	1:D:413:THR:HG22	1.89	0.72
1:E:391:LYS:HZ3	1:E:440:ASN:HD21	1.36	0.72
1:A:51:VAL:HG13	1:A:55:GLY:O	1.89	0.72
1:C:410:PRO:O	1:C:413:THR:HG22	1.89	0.72
1:D:450:ILE:CG1	1:D:451:ASP:H	1.89	0.72
1:F:112:VAL:O	1:F:112:VAL:HG13	1.88	0.72
1:F:509:THR:OG1	1:F:510:PRO:C	2.32	0.72
1:A:410:PRO:O	1:A:413:THR:HG22	1.89	0.72
1:C:450:ILE:CG1	1:C:451:ASP:H	1.89	0.72
1:E:450:ILE:CG1	1:E:451:ASP:H	1.89	0.72
1:E:509:THR:OG1	1:E:510:PRO:C	2.32	0.72
1:C:409:PRO:O	1:C:454:TYR:OH	2.06	0.72
1:A:382:SER:HB2	1:A:385:THR:HG22	1.72	0.72
1:A:409:PRO:C	1:A:454:TYR:HE1	1.98	0.72
1:B:413:THR:HG23	1:B:414:VAL:HG23	1.72	0.72
1:C:409:PRO:HD2	1:C:451:ASP:O	1.88	0.72
1:A:391:LYS:HZ1	1:A:440:ASN:ND2	1.87	0.72
1:E:409:PRO:C	1:E:454:TYR:HE1	1.98	0.72
1:E:410:PRO:O	1:E:413:THR:HG22	1.89	0.72
1:E:499:ILE:HD13	1:E:499:ILE:H	1.55	0.72
1:A:499:ILE:H	1:A:499:ILE:HD13	1.55	0.71
1:C:112:VAL:O	1:C:112:VAL:HG13	1.88	0.71
1:D:36:GLY:HA2	1:D:82:ARG:HD2	1.72	0.71
1:D:499:ILE:H	1:D:499:ILE:HD13	1.55	0.71
1:F:499:ILE:H	1:F:499:ILE:HD13	1.55	0.71
1:A:36:GLY:HA2	1:A:82:ARG:HD2	1.72	0.71
1:B:454:TYR:HE2	1:B:469:PRO:CA	2.03	0.71
1:B:509:THR:OG1	1:B:510:PRO:C	2.32	0.71
1:F:31:GLY:HA3	1:F:64:TYR:CD2	2.25	0.71
1:F:409:PRO:C	1:F:454:TYR:HE1	1.98	0.71
1:B:31:GLY:HA3	1:B:64:TYR:CD2	2.26	0.71
1:C:454:TYR:HE2	1:C:469:PRO:CA	2.03	0.71
1:D:51:VAL:HG13	1:D:55:GLY:O	1.89	0.71
1:D:394:VAL:HG11	1:D:443:ILE:HD12	1.73	0.71
1:E:23:SER:CB	1:E:483:ASN:HB2	2.21	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:409:PRO:HD2	1:B:451:ASP:O	1.88	0.71
1:F:410:PRO:O	1:F:413:THR:HG22	1.89	0.71
1:A:31:GLY:HA3	1:A:64:TYR:CD2	2.26	0.71
1:C:51:VAL:HG13	1:C:55:GLY:O	1.89	0.71
1:C:382:SER:HB2	1:C:385:THR:HG22	1.71	0.71
1:C:391:LYS:HZ1	1:C:440:ASN:ND2	1.84	0.71
1:F:413:THR:HG23	1:F:414:VAL:HG23	1.72	0.71
1:A:394:VAL:HG11	1:A:443:ILE:HD12	1.73	0.71
1:B:394:VAL:HG11	1:B:443:ILE:HD12	1.73	0.71
1:E:22:ASN:O	1:E:23:SER:OG	2.04	0.71
1:E:454:TYR:HD2	1:E:469:PRO:HA	1.53	0.71
1:B:23:SER:CB	1:B:483:ASN:HB2	2.21	0.71
1:B:51:VAL:HG13	1:B:55:GLY:O	1.89	0.71
1:C:394:VAL:HG11	1:C:443:ILE:HD12	1.73	0.71
1:E:394:VAL:HG11	1:E:443:ILE:HD12	1.73	0.71
1:F:394:VAL:HG11	1:F:443:ILE:HD12	1.73	0.71
1:B:114:ASP:OD2	1:B:175:SER:HB2	1.91	0.71
1:E:31:GLY:HA3	1:E:64:TYR:CD2	2.26	0.71
1:E:36:GLY:HA2	1:E:82:ARG:HD2	1.72	0.71
1:E:406:LEU:HD11	1:E:475:ALA:HB2	1.73	0.71
1:A:454:TYR:HE2	1:A:469:PRO:CA	2.03	0.70
1:B:379:ALA:HB2	1:B:454:TYR:CZ	2.11	0.70
1:D:413:THR:HG23	1:D:414:VAL:HG23	1.72	0.70
1:F:51:VAL:CA	1:F:55:GLY:HA2	2.20	0.70
1:F:406:LEU:HD11	1:F:475:ALA:HB2	1.73	0.70
1:B:36:GLY:HA2	1:B:82:ARG:HD2	1.72	0.70
1:C:36:GLY:HA2	1:C:82:ARG:HD2	1.72	0.70
1:D:454:TYR:HE2	1:D:469:PRO:CA	2.03	0.70
1:C:31:GLY:HA3	1:C:64:TYR:CD2	2.25	0.70
1:E:413:THR:HG23	1:E:414:VAL:HG23	1.72	0.70
1:F:23:SER:CB	1:F:483:ASN:HB2	2.21	0.70
1:F:450:ILE:CG1	1:F:451:ASP:H	1.89	0.70
1:A:23:SER:CB	1:A:483:ASN:HB2	2.21	0.70
1:A:413:THR:HG23	1:A:414:VAL:HG23	1.72	0.70
1:B:499:ILE:HD13	1:B:499:ILE:H	1.55	0.70
1:C:406:LEU:HD11	1:C:475:ALA:HB2	1.73	0.70
1:D:406:LEU:HD11	1:D:475:ALA:HB2	1.73	0.70
1:C:413:THR:HG23	1:C:414:VAL:HG23	1.72	0.70
1:C:499:ILE:HD13	1:C:499:ILE:H	1.55	0.70
1:E:382:SER:HB2	1:E:385:THR:HG22	1.71	0.70
1:F:114:ASP:OD2	1:F:175:SER:HB2	1.91	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:406:LEU:HD11	1:A:475:ALA:HB2	1.73	0.70
1:D:114:ASP:OD2	1:D:175:SER:HB2	1.91	0.70
1:F:36:GLY:HA2	1:F:82:ARG:HD2	1.72	0.70
1:F:409:PRO:HD2	1:F:451:ASP:O	1.88	0.70
1:C:47:GLU:HG3	1:C:69:MET:HG3	1.74	0.70
1:D:31:GLY:HA3	1:D:64:TYR:CD2	2.26	0.70
1:D:382:SER:HB2	1:D:385:THR:HG22	1.71	0.70
1:E:114:ASP:OD2	1:E:175:SER:HB2	1.91	0.70
1:B:406:LEU:HD11	1:B:475:ALA:HB2	1.73	0.70
1:C:22:ASN:O	1:C:23:SER:OG	2.04	0.70
1:A:114:ASP:OD2	1:A:175:SER:HB2	1.91	0.69
1:C:23:SER:CB	1:C:483:ASN:HB2	2.21	0.69
1:B:47:GLU:HG3	1:B:69:MET:HG3	1.74	0.69
1:D:23:SER:CB	1:D:483:ASN:HB2	2.21	0.69
1:F:382:SER:HB2	1:F:385:THR:HG22	1.71	0.69
1:B:22:ASN:O	1:B:23:SER:OG	2.04	0.69
1:A:391:LYS:HZ3	1:A:440:ASN:HD21	1.39	0.69
1:B:379:ALA:HB1	1:B:454:TYR:OH	1.73	0.69
1:C:114:ASP:OD2	1:C:175:SER:HB2	1.91	0.69
1:D:409:PRO:C	1:D:454:TYR:HE1	1.98	0.69
1:E:289:ILE:HD12	1:E:289:ILE:H	1.58	0.69
1:B:51:VAL:CA	1:B:55:GLY:HA2	2.20	0.69
1:B:55:GLY:O	1:B:65:PHE:CE1	2.46	0.69
1:B:215:LYS:HE3	1:B:329:ASN:HD21	1.58	0.69
1:C:454:TYR:HD2	1:C:469:PRO:HA	1.53	0.69
1:D:55:GLY:O	1:D:65:PHE:CE1	2.46	0.69
1:D:289:ILE:H	1:D:289:ILE:HD12	1.57	0.69
1:D:454:TYR:HD2	1:D:469:PRO:HA	1.53	0.69
1:E:51:VAL:CA	1:E:55:GLY:HA2	2.20	0.69
1:E:215:LYS:HE3	1:E:329:ASN:HD21	1.58	0.69
1:E:454:TYR:HE2	1:E:469:PRO:CA	2.03	0.69
1:F:289:ILE:H	1:F:289:ILE:HD12	1.58	0.69
1:A:47:GLU:HG3	1:A:69:MET:HG3	1.74	0.69
1:C:215:LYS:HE3	1:C:329:ASN:HD21	1.58	0.69
1:D:47:GLU:HG3	1:D:69:MET:HG3	1.74	0.69
1:F:55:GLY:O	1:F:65:PHE:CE1	2.46	0.69
1:D:215:LYS:HE3	1:D:329:ASN:HD21	1.58	0.69
1:B:289:ILE:HD12	1:B:289:ILE:H	1.58	0.68
1:B:454:TYR:HD2	1:B:469:PRO:HA	1.53	0.68
1:F:454:TYR:HE2	1:F:469:PRO:CA	2.03	0.68
1:C:289:ILE:H	1:C:289:ILE:HD12	1.57	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:47:GLU:HG3	1:E:69:MET:HG3	1.74	0.68
1:E:455:LYS:HG3	1:E:502:VAL:HG22	1.75	0.68
1:B:455:LYS:HG3	1:B:502:VAL:HG22	1.75	0.68
1:C:455:LYS:HG3	1:C:502:VAL:HG22	1.75	0.68
1:F:391:LYS:HZ2	1:F:440:ASN:ND2	1.92	0.68
1:A:55:GLY:O	1:A:65:PHE:CE1	2.46	0.68
1:A:455:LYS:HG3	1:A:502:VAL:HG22	1.76	0.68
1:D:455:LYS:HG3	1:D:502:VAL:HG22	1.75	0.68
1:F:47:GLU:HG3	1:F:69:MET:HG3	1.74	0.68
1:F:455:LYS:HG3	1:F:502:VAL:HG22	1.76	0.68
1:C:55:GLY:O	1:C:65:PHE:CE1	2.46	0.68
1:A:289:ILE:H	1:A:289:ILE:HD12	1.58	0.68
1:E:23:SER:HB3	1:E:483:ASN:CB	2.22	0.68
1:F:215:LYS:HE3	1:F:329:ASN:HD21	1.58	0.68
1:A:23:SER:HB3	1:A:483:ASN:CB	2.22	0.68
1:A:215:LYS:HE3	1:A:329:ASN:HD21	1.58	0.68
1:E:391:LYS:HZ1	1:E:440:ASN:ND2	1.90	0.68
1:A:427:VAL:O	1:A:431:THR:HG22	1.94	0.68
1:E:55:GLY:O	1:E:65:PHE:CE1	2.46	0.68
1:C:23:SER:HB3	1:C:483:ASN:CB	2.22	0.67
1:E:228:GLY:CA	1:E:345:SER:HB3	2.23	0.67
1:B:427:VAL:O	1:B:431:THR:HG22	1.95	0.67
1:B:450:ILE:CG1	1:B:451:ASP:N	2.53	0.67
1:C:500:LEU:HB2	1:C:501:ASN:OD1	1.95	0.67
1:F:500:LEU:HB2	1:F:501:ASN:OD1	1.95	0.67
1:D:23:SER:HB3	1:D:483:ASN:CB	2.22	0.67
1:D:171:ILE:HG22	1:D:172:SER:N	2.09	0.67
1:F:408:SER:HB2	1:F:471:ALA:HB2	1.77	0.67
1:A:171:ILE:HG22	1:A:172:SER:N	2.09	0.67
1:B:23:SER:HB3	1:B:483:ASN:CB	2.22	0.67
1:B:171:ILE:HG22	1:B:172:SER:N	2.09	0.67
1:B:304:ILE:HG13	1:B:305:TYR:CE2	2.30	0.67
1:D:500:LEU:HB2	1:D:501:ASN:OD1	1.95	0.67
1:E:171:ILE:HG22	1:E:172:SER:N	2.09	0.67
1:A:55:GLY:O	1:A:65:PHE:HE1	1.78	0.67
1:C:408:SER:HB2	1:C:471:ALA:HB2	1.77	0.67
1:C:427:VAL:O	1:C:431:THR:HG22	1.94	0.67
1:E:408:SER:HB2	1:E:471:ALA:HB2	1.77	0.67
1:E:500:LEU:HB2	1:E:501:ASN:OD1	1.95	0.67
1:F:55:GLY:O	1:F:65:PHE:HE1	1.78	0.67
1:A:304:ILE:HG13	1:A:305:TYR:CE2	2.30	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:304:ILE:HG13	1:C:305:TYR:CE2	2.30	0.67
1:A:408:SER:HB2	1:A:471:ALA:HB2	1.77	0.67
1:D:51:VAL:CA	1:D:55:GLY:HA2	2.20	0.67
1:D:408:SER:HB2	1:D:471:ALA:HB2	1.77	0.67
1:F:427:VAL:O	1:F:431:THR:HG22	1.94	0.67
1:F:23:SER:HB3	1:F:483:ASN:CB	2.22	0.67
1:B:96:ALA:HB2	1:B:191:SER:HA	1.77	0.67
1:D:304:ILE:HG13	1:D:305:TYR:CE2	2.30	0.66
1:A:96:ALA:HB2	1:A:191:SER:HA	1.77	0.66
1:C:171:ILE:HG22	1:C:172:SER:N	2.09	0.66
1:D:55:GLY:O	1:D:65:PHE:HE1	1.78	0.66
1:D:379:ALA:CB	1:D:454:TYR:CE2	2.79	0.66
1:A:500:LEU:HB2	1:A:501:ASN:OD1	1.95	0.66
1:E:96:ALA:HB2	1:E:191:SER:HA	1.77	0.66
1:E:304:ILE:HG13	1:E:305:TYR:CE2	2.30	0.66
1:C:379:ALA:CB	1:C:454:TYR:CE2	2.79	0.66
1:D:351:THR:HG23	1:D:354:ASP:H	1.60	0.66
1:E:55:GLY:O	1:E:65:PHE:HE1	1.78	0.66
1:E:351:THR:HG23	1:E:354:ASP:H	1.60	0.66
1:F:228:GLY:CA	1:F:345:SER:HB3	2.23	0.66
1:F:304:ILE:HG13	1:F:305:TYR:CE2	2.30	0.66
1:F:379:ALA:CB	1:F:454:TYR:CE2	2.79	0.66
1:A:228:GLY:CA	1:A:345:SER:HB3	2.24	0.66
1:B:55:GLY:O	1:B:65:PHE:HE1	1.78	0.66
1:B:408:SER:HB2	1:B:471:ALA:HB2	1.77	0.66
1:F:171:ILE:HG22	1:F:172:SER:N	2.09	0.66
1:A:409:PRO:HD2	1:A:451:ASP:O	1.88	0.66
1:B:500:LEU:HB2	1:B:501:ASN:OD1	1.95	0.66
1:C:409:PRO:C	1:C:454:TYR:HE1	1.98	0.66
1:D:427:VAL:O	1:D:431:THR:HG22	1.94	0.66
1:A:450:ILE:CG1	1:A:451:ASP:H	1.89	0.66
1:B:379:ALA:CB	1:B:454:TYR:CE2	2.79	0.66
1:E:427:VAL:O	1:E:431:THR:HG22	1.94	0.66
1:A:51:VAL:CA	1:A:55:GLY:HA2	2.20	0.66
1:A:121:VAL:HG22	1:A:166:ASN:HB3	1.78	0.66
1:A:351:THR:HG23	1:A:354:ASP:H	1.60	0.66
1:B:228:GLY:CA	1:B:345:SER:HB3	2.23	0.66
1:D:96:ALA:HB2	1:D:191:SER:HA	1.77	0.66
1:E:379:ALA:CB	1:E:454:TYR:CE2	2.79	0.66
1:F:121:VAL:HG22	1:F:166:ASN:HB3	1.78	0.66
1:B:351:THR:HG23	1:B:354:ASP:H	1.60	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:409:PRO:C	1:B:454:TYR:HE1	1.98	0.65
1:C:96:ALA:HB2	1:C:191:SER:HA	1.77	0.65
1:C:351:THR:HG23	1:C:354:ASP:H	1.60	0.65
1:F:96:ALA:HB2	1:F:191:SER:HA	1.77	0.65
1:E:121:VAL:HG22	1:E:166:ASN:HB3	1.78	0.65
1:A:379:ALA:CB	1:A:454:TYR:CE2	2.78	0.65
1:B:391:LYS:NZ	1:B:440:ASN:HD21	1.92	0.65
1:C:51:VAL:CA	1:C:55:GLY:HA2	2.20	0.65
1:F:351:THR:HG23	1:F:354:ASP:H	1.60	0.65
1:B:121:VAL:HG22	1:B:166:ASN:HB3	1.78	0.65
1:F:391:LYS:NZ	1:F:440:ASN:HD21	1.92	0.65
1:A:391:LYS:NZ	1:A:440:ASN:HD21	1.92	0.64
1:C:55:GLY:O	1:C:65:PHE:HE1	1.78	0.64
1:C:66:MET:HG2	1:C:468:VAL:HG11	1.80	0.64
1:C:228:GLY:CA	1:C:345:SER:HB3	2.23	0.64
1:A:454:TYR:HD2	1:A:469:PRO:HA	1.53	0.64
1:A:100:GLU:HG3	1:A:186:LYS:H	1.63	0.64
1:B:66:MET:HG2	1:B:468:VAL:HG11	1.80	0.64
1:B:100:GLU:HG3	1:B:186:LYS:H	1.63	0.64
1:C:100:GLU:HG3	1:C:186:LYS:H	1.63	0.64
1:B:391:LYS:HZ2	1:B:440:ASN:ND2	1.93	0.64
1:C:121:VAL:HG22	1:C:166:ASN:HB3	1.78	0.64
1:E:100:GLU:HG3	1:E:186:LYS:H	1.63	0.64
1:D:121:VAL:HG22	1:D:166:ASN:HB3	1.78	0.64
1:D:215:LYS:HE3	1:D:329:ASN:ND2	2.13	0.64
1:D:66:MET:HG2	1:D:468:VAL:HG11	1.80	0.64
1:F:100:GLU:HG3	1:F:186:LYS:H	1.63	0.63
1:C:215:LYS:HE3	1:C:329:ASN:ND2	2.13	0.63
1:E:215:LYS:HE3	1:E:329:ASN:ND2	2.13	0.63
1:A:499:ILE:HG13	1:A:502:VAL:HG21	1.81	0.63
1:B:499:ILE:HG13	1:B:502:VAL:HG21	1.81	0.63
1:A:66:MET:HG2	1:A:468:VAL:HG11	1.80	0.63
1:D:100:GLU:HG3	1:D:186:LYS:H	1.63	0.63
1:D:228:GLY:CA	1:D:345:SER:HB3	2.23	0.63
1:B:215:LYS:HE3	1:B:329:ASN:ND2	2.13	0.63
1:A:23:SER:HG	1:A:483:ASN:HB3	1.64	0.62
1:E:66:MET:HG2	1:E:468:VAL:HG11	1.80	0.62
1:C:107:GLY:HA3	1:C:110:TYR:CE1	2.35	0.62
1:F:66:MET:HG2	1:F:468:VAL:HG11	1.80	0.62
1:F:499:ILE:HG13	1:F:502:VAL:HG21	1.81	0.62
1:B:509:THR:H	1:B:510:PRO:CA	2.12	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:499:ILE:HG13	1:C:502:VAL:HG21	1.81	0.62
1:F:509:THR:H	1:F:510:PRO:CA	2.12	0.62
1:F:215:LYS:HE3	1:F:329:ASN:ND2	2.13	0.62
1:A:509:THR:H	1:A:510:PRO:CA	2.12	0.62
1:B:213:ASN:HD22	1:B:213:ASN:N	1.98	0.62
1:D:509:THR:H	1:D:510:PRO:CA	2.12	0.62
1:A:213:ASN:HD22	1:A:213:ASN:N	1.98	0.62
1:A:407:CYS:O	1:A:450:ILE:HA	2.00	0.62
1:D:407:CYS:O	1:D:450:ILE:HA	2.00	0.62
1:B:107:GLY:HA3	1:B:110:TYR:CE1	2.34	0.62
1:C:213:ASN:HD22	1:C:213:ASN:N	1.98	0.62
1:B:407:CYS:O	1:B:450:ILE:HA	2.00	0.62
1:C:509:THR:H	1:C:510:PRO:CA	2.12	0.62
1:F:107:GLY:HA3	1:F:110:TYR:CE1	2.35	0.62
1:D:213:ASN:HD22	1:D:213:ASN:N	1.98	0.62
1:D:499:ILE:HG13	1:D:502:VAL:HG21	1.81	0.62
1:E:107:GLY:HA3	1:E:110:TYR:HE1	1.65	0.62
1:E:107:GLY:HA3	1:E:110:TYR:CE1	2.34	0.61
1:B:278:GLN:HG2	1:B:296:SER:HB2	1.82	0.61
1:C:278:GLN:HG2	1:C:296:SER:HB2	1.82	0.61
1:D:107:GLY:HA3	1:D:110:TYR:CE1	2.34	0.61
1:F:213:ASN:HD22	1:F:213:ASN:N	1.98	0.61
1:A:107:GLY:HA3	1:A:110:TYR:HE1	1.65	0.61
1:C:455:LYS:HE2	1:C:502:VAL:HG22	1.82	0.61
1:D:107:GLY:HA3	1:D:110:TYR:HE1	1.65	0.61
1:D:455:LYS:HE2	1:D:502:VAL:HG22	1.82	0.61
1:E:509:THR:H	1:E:510:PRO:CA	2.12	0.61
1:A:107:GLY:HA3	1:A:110:TYR:CE1	2.34	0.61
1:A:215:LYS:HE3	1:A:329:ASN:ND2	2.13	0.61
1:F:407:CYS:O	1:F:450:ILE:HA	2.00	0.61
1:E:213:ASN:HD22	1:E:213:ASN:N	1.98	0.61
1:E:409:PRO:HD2	1:E:451:ASP:O	1.88	0.61
1:B:50:LEU:C	1:B:50:LEU:HD12	2.26	0.61
1:A:379:ALA:HB1	1:A:454:TYR:OH	1.73	0.61
1:B:107:GLY:HA3	1:B:110:TYR:HE1	1.65	0.61
1:C:107:GLY:HA3	1:C:110:TYR:HE1	1.65	0.61
1:D:278:GLN:HG2	1:D:296:SER:HB2	1.83	0.61
1:E:499:ILE:HG13	1:E:502:VAL:HG21	1.81	0.61
1:C:407:CYS:O	1:C:450:ILE:HA	2.00	0.61
1:B:455:LYS:HE2	1:B:502:VAL:HG22	1.82	0.61
1:E:391:LYS:CE	1:E:440:ASN:O	2.49	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:407:CYS:O	1:E:450:ILE:HA	2.00	0.61
1:F:107:GLY:HA3	1:F:110:TYR:HE1	1.65	0.61
1:D:50:LEU:C	1:D:50:LEU:HD12	2.26	0.60
1:A:278:GLN:HG2	1:A:296:SER:HB2	1.83	0.60
1:D:391:LYS:CE	1:D:440:ASN:O	2.49	0.60
1:C:509:THR:N	1:C:510:PRO:HA	2.14	0.60
1:E:278:GLN:HG2	1:E:296:SER:HB2	1.83	0.60
1:F:161:PRO:HB3	1:F:187:ILE:HB	1.83	0.60
1:C:161:PRO:HB3	1:C:187:ILE:HB	1.83	0.60
1:E:97:GLY:HA3	1:E:256:PRO:HG2	1.84	0.60
1:E:455:LYS:HE2	1:E:502:VAL:HG22	1.82	0.60
1:F:455:LYS:HE2	1:F:502:VAL:HG22	1.82	0.60
1:A:50:LEU:C	1:A:50:LEU:HD12	2.26	0.60
1:A:161:PRO:HB3	1:A:187:ILE:HB	1.83	0.60
1:C:50:LEU:C	1:C:50:LEU:HD12	2.26	0.60
1:E:50:LEU:C	1:E:50:LEU:HD12	2.26	0.60
1:F:97:GLY:HA3	1:F:256:PRO:HG2	1.84	0.60
1:F:509:THR:N	1:F:510:PRO:HA	2.14	0.60
1:E:509:THR:N	1:E:510:PRO:HA	2.14	0.60
1:F:50:LEU:C	1:F:50:LEU:HD12	2.26	0.60
1:B:347:ASN:O	1:B:348:ALA:C	2.45	0.59
1:F:278:GLN:HG2	1:F:296:SER:HB2	1.83	0.59
1:A:347:ASN:O	1:A:348:ALA:C	2.45	0.59
1:C:347:ASN:O	1:C:348:ALA:C	2.45	0.59
1:B:66:MET:HA	1:B:66:MET:HE3	1.84	0.59
1:E:419:VAL:HA	1:E:422:ALA:HB3	1.84	0.59
1:A:455:LYS:HE2	1:A:502:VAL:HG22	1.82	0.59
1:B:97:GLY:HA3	1:B:256:PRO:HG2	1.84	0.59
1:E:161:PRO:HB3	1:E:187:ILE:HB	1.83	0.59
1:E:391:LYS:NZ	1:E:440:ASN:HD21	1.92	0.59
1:B:161:PRO:HB3	1:B:187:ILE:HB	1.83	0.59
1:B:391:LYS:CE	1:B:440:ASN:O	2.49	0.59
1:C:382:SER:O	1:C:383:LEU:C	2.45	0.59
1:D:66:MET:HA	1:D:66:MET:HE3	1.84	0.59
1:D:178:LEU:CD2	1:D:178:LEU:N	2.66	0.59
1:F:419:VAL:HA	1:F:422:ALA:HB3	1.84	0.59
1:C:391:LYS:CE	1:C:440:ASN:O	2.49	0.59
1:A:97:GLY:HA3	1:A:256:PRO:HG2	1.84	0.59
1:C:97:GLY:HA3	1:C:256:PRO:HG2	1.84	0.59
1:D:60:GLU:HG2	1:D:347:ASN:HB2	1.85	0.59
1:D:161:PRO:HB3	1:D:187:ILE:HB	1.83	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:382:SER:O	1:F:383:LEU:C	2.45	0.59
1:D:347:ASN:O	1:D:348:ALA:C	2.45	0.58
1:D:419:VAL:HA	1:D:422:ALA:HB3	1.84	0.58
1:F:501:ASN:OD1	1:F:501:ASN:N	2.36	0.58
1:C:60:GLU:HG2	1:C:347:ASN:HB2	1.85	0.58
1:D:509:THR:N	1:D:510:PRO:HA	2.14	0.58
1:E:382:SER:O	1:E:383:LEU:C	2.45	0.58
1:A:391:LYS:CE	1:A:440:ASN:O	2.49	0.58
1:B:178:LEU:CD2	1:B:178:LEU:N	2.66	0.58
1:F:347:ASN:O	1:F:348:ALA:C	2.45	0.58
1:F:373:PHE:HB2	1:F:405:VAL:HA	1.86	0.58
1:C:66:MET:HA	1:C:66:MET:HE3	1.85	0.58
1:D:97:GLY:HA3	1:D:256:PRO:HG2	1.84	0.58
1:D:429:TRP:CZ2	1:D:441:PHE:HB2	2.39	0.58
1:E:347:ASN:O	1:E:348:ALA:C	2.45	0.58
1:E:429:TRP:CZ2	1:E:441:PHE:HB2	2.39	0.58
1:F:456:TYR:HB3	1:F:504:LYS:HB3	1.86	0.58
1:A:373:PHE:HB2	1:A:405:VAL:HA	1.86	0.58
1:B:398:ASP:O	1:B:401:GLN:HG3	2.04	0.58
1:C:398:ASP:O	1:C:401:GLN:HG3	2.04	0.58
1:F:398:ASP:O	1:F:401:GLN:HG3	2.04	0.58
1:B:456:TYR:HB3	1:B:504:LYS:HB3	1.85	0.58
1:E:60:GLU:HG2	1:E:347:ASN:HB2	1.85	0.58
1:E:373:PHE:HB2	1:E:405:VAL:HA	1.86	0.58
1:A:133:GLU:HB3	1:A:142:LYS:HB3	1.86	0.58
1:A:419:VAL:HA	1:A:422:ALA:HB3	1.84	0.58
1:E:455:LYS:HG2	1:E:456:TYR:N	2.19	0.58
1:F:133:GLU:HB3	1:F:142:LYS:HB3	1.86	0.58
1:A:429:TRP:CZ2	1:A:441:PHE:HB2	2.39	0.58
1:B:419:VAL:HA	1:B:422:ALA:HB3	1.84	0.58
1:C:373:PHE:HB2	1:C:404:LEU:O	2.04	0.58
1:D:382:SER:O	1:D:383:LEU:C	2.45	0.58
1:E:66:MET:HA	1:E:66:MET:HE3	1.85	0.58
1:E:398:ASP:O	1:E:401:GLN:HG3	2.04	0.58
1:E:456:TYR:O	1:E:503:ILE:HG12	2.04	0.58
1:F:373:PHE:HB2	1:F:404:LEU:O	2.04	0.58
1:F:456:TYR:O	1:F:503:ILE:HG12	2.04	0.58
1:A:382:SER:O	1:A:383:LEU:C	2.45	0.58
1:A:501:ASN:OD1	1:A:501:ASN:N	2.36	0.58
1:C:429:TRP:CZ2	1:C:441:PHE:HB2	2.39	0.58
1:D:456:TYR:O	1:D:503:ILE:HG12	2.04	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:456:TYR:O	1:A:503:ILE:HG12	2.04	0.57
1:B:501:ASN:OD1	1:B:501:ASN:N	2.36	0.57
1:E:133:GLU:HB3	1:E:142:LYS:HB3	1.86	0.57
1:E:178:LEU:CD2	1:E:178:LEU:N	2.66	0.57
1:F:66:MET:HA	1:F:66:MET:HE3	1.85	0.57
1:A:398:ASP:O	1:A:401:GLN:HG3	2.04	0.57
1:B:456:TYR:O	1:B:503:ILE:HG12	2.04	0.57
1:E:23:SER:OG	1:E:483:ASN:HB2	2.04	0.57
1:F:455:LYS:HG2	1:F:456:TYR:N	2.19	0.57
1:C:114:ASP:CG	1:C:175:SER:HB2	2.29	0.57
1:C:456:TYR:O	1:C:503:ILE:HG12	2.04	0.57
1:D:398:ASP:O	1:D:401:GLN:HG3	2.04	0.57
1:E:114:ASP:CG	1:E:175:SER:HB2	2.29	0.57
1:E:373:PHE:HB2	1:E:404:LEU:O	2.04	0.57
1:F:60:GLU:HG2	1:F:347:ASN:HB2	1.85	0.57
1:F:114:ASP:CG	1:F:175:SER:HB2	2.29	0.57
1:F:429:TRP:CZ2	1:F:441:PHE:HB2	2.39	0.57
1:A:114:ASP:CG	1:A:175:SER:HB2	2.29	0.57
1:B:60:GLU:HG2	1:B:347:ASN:HB2	1.85	0.57
1:B:382:SER:O	1:B:383:LEU:C	2.45	0.57
1:B:429:TRP:CZ2	1:B:441:PHE:HB2	2.39	0.57
1:D:373:PHE:HB2	1:D:405:VAL:HA	1.86	0.57
1:A:455:LYS:HG2	1:A:456:TYR:N	2.19	0.57
1:A:456:TYR:HB3	1:A:504:LYS:HB3	1.86	0.57
1:C:499:ILE:HG13	1:C:502:VAL:CG2	2.34	0.57
1:D:114:ASP:CG	1:D:175:SER:HB2	2.29	0.57
1:D:373:PHE:HB2	1:D:404:LEU:O	2.04	0.57
1:D:501:ASN:OD1	1:D:501:ASN:N	2.36	0.57
1:B:373:PHE:HB2	1:B:404:LEU:O	2.04	0.57
1:C:419:VAL:HA	1:C:422:ALA:HB3	1.84	0.57
1:C:501:ASN:OD1	1:C:501:ASN:N	2.36	0.57
1:D:456:TYR:HB3	1:D:504:LYS:HB3	1.85	0.57
1:E:501:ASN:OD1	1:E:501:ASN:N	2.36	0.57
1:A:66:MET:HE3	1:A:66:MET:HA	1.85	0.57
1:B:133:GLU:HB3	1:B:142:LYS:HB3	1.86	0.57
1:C:455:LYS:HG2	1:C:456:TYR:N	2.19	0.57
1:F:379:ALA:HB1	1:F:454:TYR:OH	1.73	0.57
1:E:456:TYR:HB3	1:E:504:LYS:HB3	1.86	0.57
1:E:499:ILE:HG13	1:E:502:VAL:CG2	2.35	0.57
1:A:60:GLU:HG2	1:A:347:ASN:HB2	1.85	0.57
1:A:499:ILE:HG13	1:A:502:VAL:CG2	2.35	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:114:ASP:CG	1:B:175:SER:HB2	2.29	0.57
1:D:499:ILE:HG13	1:D:502:VAL:CG2	2.35	0.57
1:F:391:LYS:CE	1:F:440:ASN:O	2.49	0.57
1:B:373:PHE:HB2	1:B:405:VAL:HA	1.86	0.57
1:A:373:PHE:HB2	1:A:404:LEU:O	2.04	0.56
1:C:373:PHE:HB2	1:C:405:VAL:HA	1.86	0.56
1:D:455:LYS:HG2	1:D:456:TYR:N	2.19	0.56
1:C:456:TYR:HB3	1:C:504:LYS:HB3	1.86	0.56
1:A:431:THR:O	1:A:432:ALA:HB2	2.06	0.56
1:B:84:VAL:CG1	1:B:85:ASP:O	2.54	0.56
1:B:454:TYR:HE2	1:B:469:PRO:CB	2.19	0.56
1:B:499:ILE:HG13	1:B:502:VAL:CG2	2.35	0.56
1:C:133:GLU:HB3	1:C:142:LYS:HB3	1.86	0.56
1:C:178:LEU:CD2	1:C:178:LEU:N	2.66	0.56
1:D:133:GLU:HB3	1:D:142:LYS:HB3	1.86	0.56
1:D:431:THR:O	1:D:432:ALA:HB2	2.06	0.56
1:B:455:LYS:HG2	1:B:456:TYR:N	2.19	0.56
1:F:499:ILE:HG13	1:F:502:VAL:CG2	2.35	0.56
1:F:431:THR:O	1:F:432:ALA:HB2	2.06	0.56
1:A:84:VAL:CG1	1:A:85:ASP:O	2.54	0.56
1:C:431:THR:O	1:C:432:ALA:HB2	2.06	0.56
1:C:503:ILE:O	1:C:504:LYS:HB2	2.06	0.56
1:F:23:SER:HG	1:F:483:ASN:HB3	1.69	0.56
1:B:23:SER:OG	1:B:483:ASN:HB2	2.04	0.56
1:B:509:THR:N	1:B:510:PRO:HA	2.14	0.56
1:C:409:PRO:C	1:C:454:TYR:CE1	2.71	0.56
1:D:89:ALA:HB3	1:D:194:LEU:HD11	1.88	0.56
1:F:236:ILE:HD12	1:F:236:ILE:H	1.71	0.56
1:F:503:ILE:O	1:F:504:LYS:HB2	2.06	0.56
1:C:236:ILE:HD12	1:C:236:ILE:H	1.71	0.55
1:D:454:TYR:HE2	1:D:469:PRO:CB	2.19	0.55
1:E:46:ASN:HB2	1:E:49:ASP:HB2	1.89	0.55
1:E:283:VAL:HG21	1:E:323:ILE:HD13	1.88	0.55
1:C:46:ASN:HB2	1:C:49:ASP:HB2	1.89	0.55
1:C:84:VAL:CG1	1:C:85:ASP:O	2.54	0.55
1:D:236:ILE:H	1:D:236:ILE:HD12	1.72	0.55
1:E:454:TYR:HE2	1:E:469:PRO:CB	2.19	0.55
1:F:283:VAL:HG21	1:F:323:ILE:HD13	1.88	0.55
1:B:236:ILE:H	1:B:236:ILE:HD12	1.71	0.55
1:C:234:ILE:HB	1:C:340:LEU:HD12	1.89	0.55
1:F:381:GLU:HB3	1:F:385:THR:HG23	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:23:SER:OG	1:A:483:ASN:HB2	2.04	0.55
1:A:503:ILE:O	1:A:504:LYS:HB2	2.06	0.55
1:C:89:ALA:HB3	1:C:194:LEU:HD11	1.88	0.55
1:C:150:ILE:HG22	1:C:167:TRP:CZ3	2.42	0.55
1:D:283:VAL:HG21	1:D:323:ILE:HD13	1.88	0.55
1:D:381:GLU:HB3	1:D:385:THR:HG23	1.89	0.55
1:E:150:ILE:HG22	1:E:167:TRP:CZ3	2.42	0.55
1:E:381:GLU:HB3	1:E:385:THR:HG23	1.89	0.55
1:F:46:ASN:HB2	1:F:49:ASP:HB2	1.88	0.55
1:F:234:ILE:HB	1:F:340:LEU:HD12	1.89	0.55
1:A:150:ILE:HG22	1:A:167:TRP:CZ3	2.42	0.55
1:A:454:TYR:HE2	1:A:469:PRO:CB	2.19	0.55
1:B:46:ASN:HB2	1:B:49:ASP:HB2	1.89	0.55
1:D:84:VAL:CG1	1:D:85:ASP:O	2.54	0.55
1:D:503:ILE:O	1:D:504:LYS:HB2	2.06	0.55
1:F:84:VAL:CG1	1:F:85:ASP:O	2.54	0.55
1:A:71:PHE:C	1:A:71:PHE:CD2	2.85	0.55
1:B:431:THR:O	1:B:432:ALA:HB2	2.06	0.55
1:D:455:LYS:HE2	1:D:502:VAL:CG2	2.37	0.55
1:E:89:ALA:HB3	1:E:194:LEU:HD11	1.88	0.55
1:E:455:LYS:HE2	1:E:502:VAL:CG2	2.37	0.55
1:A:498:GLN:OE1	1:A:498:GLN:HA	2.07	0.55
1:B:409:PRO:C	1:B:454:TYR:CE1	2.71	0.55
1:C:454:TYR:HE2	1:C:469:PRO:CB	2.19	0.55
1:C:455:LYS:HE2	1:C:502:VAL:CG2	2.37	0.55
1:E:431:THR:O	1:E:432:ALA:HB2	2.06	0.55
1:A:150:ILE:HG22	1:A:167:TRP:HZ3	1.72	0.55
1:A:234:ILE:HB	1:A:340:LEU:HD12	1.89	0.55
1:B:150:ILE:HG22	1:B:167:TRP:CZ3	2.42	0.55
1:B:274:GLN:H	1:B:278:GLN:NE2	2.05	0.55
1:D:46:ASN:HB2	1:D:49:ASP:HB2	1.89	0.55
1:D:150:ILE:HG22	1:D:167:TRP:CZ3	2.42	0.55
1:E:274:GLN:H	1:E:278:GLN:NE2	2.05	0.55
1:F:454:TYR:HE2	1:F:469:PRO:CB	2.19	0.55
1:A:236:ILE:H	1:A:236:ILE:HD12	1.71	0.55
1:B:71:PHE:C	1:B:71:PHE:CD2	2.85	0.55
1:C:283:VAL:HG21	1:C:323:ILE:HD13	1.88	0.55
1:C:498:GLN:OE1	1:C:498:GLN:HA	2.07	0.55
1:D:302:LYS:HD3	1:D:306:ASP:HA	1.88	0.55
1:E:302:LYS:HD3	1:E:306:ASP:HA	1.89	0.55
1:B:302:LYS:HD3	1:B:306:ASP:HA	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:498:GLN:OE1	1:B:498:GLN:HA	2.07	0.55
1:A:46:ASN:HB2	1:A:49:ASP:HB2	1.89	0.54
1:B:283:VAL:HG21	1:B:323:ILE:HD13	1.88	0.54
1:C:206:THR:O	1:C:206:THR:HG22	2.07	0.54
1:D:498:GLN:HA	1:D:498:GLN:OE1	2.07	0.54
1:B:234:ILE:HB	1:B:340:LEU:HD12	1.89	0.54
1:C:381:GLU:HB3	1:C:385:THR:HG23	1.89	0.54
1:D:234:ILE:HB	1:D:340:LEU:HD12	1.89	0.54
1:E:503:ILE:O	1:E:504:LYS:HB2	2.06	0.54
1:F:71:PHE:C	1:F:71:PHE:CD2	2.85	0.54
1:A:283:VAL:HG21	1:A:323:ILE:HD13	1.88	0.54
1:A:381:GLU:HB3	1:A:385:THR:HG23	1.89	0.54
1:B:503:ILE:O	1:B:504:LYS:HB2	2.06	0.54
1:D:51:VAL:HG22	1:D:65:PHE:HZ	1.73	0.54
1:E:236:ILE:H	1:E:236:ILE:HD12	1.72	0.54
1:F:150:ILE:HG22	1:F:167:TRP:HZ3	1.72	0.54
1:B:206:THR:O	1:B:206:THR:HG22	2.07	0.54
1:F:498:GLN:OE1	1:F:498:GLN:HA	2.07	0.54
1:B:150:ILE:HG22	1:B:167:TRP:HZ3	1.72	0.54
1:B:455:LYS:HE2	1:B:502:VAL:CG2	2.37	0.54
1:D:206:THR:HG22	1:D:206:THR:O	2.07	0.54
1:D:274:GLN:H	1:D:278:GLN:NE2	2.05	0.54
1:B:89:ALA:HB3	1:B:194:LEU:HD11	1.88	0.54
1:C:71:PHE:C	1:C:71:PHE:CD2	2.85	0.54
1:E:84:VAL:CG1	1:E:85:ASP:O	2.54	0.54
1:F:89:ALA:HB3	1:F:194:LEU:HD11	1.88	0.54
1:F:302:LYS:HD3	1:F:306:ASP:HA	1.88	0.54
1:C:274:GLN:H	1:C:278:GLN:NE2	2.05	0.54
1:E:498:GLN:OE1	1:E:498:GLN:HA	2.07	0.54
1:F:150:ILE:HG22	1:F:167:TRP:CZ3	2.42	0.54
1:F:206:THR:O	1:F:206:THR:HG22	2.07	0.54
1:A:206:THR:HG22	1:A:206:THR:O	2.07	0.54
1:B:381:GLU:HB3	1:B:385:THR:HG23	1.89	0.54
1:D:150:ILE:HG22	1:D:167:TRP:HZ3	1.72	0.54
1:E:71:PHE:C	1:E:71:PHE:CD2	2.85	0.54
1:A:51:VAL:HG22	1:A:65:PHE:HZ	1.73	0.54
1:A:455:LYS:HE2	1:A:502:VAL:CG2	2.37	0.54
1:D:71:PHE:C	1:D:71:PHE:CD2	2.85	0.54
1:E:150:ILE:HG22	1:E:167:TRP:HZ3	1.72	0.54
1:E:244:TYR:CD2	1:E:273:PRO:HD2	2.43	0.54
1:F:244:TYR:CD2	1:F:273:PRO:HD2	2.43	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:302:LYS:HD3	1:A:306:ASP:HA	1.88	0.54
1:B:244:TYR:CD2	1:B:273:PRO:HD2	2.43	0.54
1:C:302:LYS:HD3	1:C:306:ASP:HA	1.88	0.54
1:D:509:THR:HG1	1:D:510:PRO:HA	1.71	0.54
1:E:90:LYS:HB2	1:E:344:LEU:HB3	1.90	0.54
1:E:234:ILE:HB	1:E:340:LEU:HD12	1.89	0.54
1:F:90:LYS:HB2	1:F:344:LEU:HB3	1.90	0.54
1:A:89:ALA:HB3	1:A:194:LEU:HD11	1.88	0.53
1:B:51:VAL:HG22	1:B:65:PHE:HZ	1.73	0.53
1:E:51:VAL:HG22	1:E:65:PHE:HZ	1.73	0.53
1:E:206:THR:O	1:E:206:THR:HG22	2.07	0.53
1:F:173:SER:N	1:F:174:SER:HA	2.22	0.53
1:F:455:LYS:HE2	1:F:502:VAL:CG2	2.37	0.53
1:A:90:LYS:HB2	1:A:344:LEU:HB3	1.90	0.53
1:B:90:LYS:HB2	1:B:344:LEU:HB3	1.90	0.53
1:C:408:SER:CB	1:C:451:ASP:HB3	2.39	0.53
1:F:178:LEU:CD2	1:F:178:LEU:N	2.66	0.53
1:F:409:PRO:C	1:F:454:TYR:CE1	2.71	0.53
1:A:244:TYR:CD2	1:A:273:PRO:HD2	2.43	0.53
1:A:274:GLN:H	1:A:278:GLN:NE2	2.05	0.53
1:A:371:GLN:HB3	1:A:484:VAL:HG23	1.90	0.53
1:D:90:LYS:HB2	1:D:344:LEU:HB3	1.90	0.53
1:D:408:SER:CB	1:D:451:ASP:HB3	2.39	0.53
1:A:245:ALA:C	1:A:247:GLY:H	2.16	0.53
1:B:371:GLN:HB3	1:B:484:VAL:HG23	1.90	0.53
1:C:51:VAL:HG22	1:C:65:PHE:HZ	1.73	0.53
1:D:244:TYR:CD2	1:D:273:PRO:HD2	2.43	0.53
1:E:245:ALA:C	1:E:247:GLY:H	2.16	0.53
1:A:409:PRO:C	1:A:454:TYR:CE1	2.71	0.53
1:C:90:LYS:HB2	1:C:344:LEU:HB3	1.90	0.53
1:C:172:SER:O	1:C:174:SER:HA	2.08	0.53
1:C:244:TYR:CD2	1:C:273:PRO:HD2	2.43	0.53
1:F:274:GLN:H	1:F:278:GLN:NE2	2.05	0.53
1:A:173:SER:N	1:A:174:SER:HA	2.22	0.53
1:B:451:ASP:OD2	1:B:471:ALA:N	2.42	0.53
1:D:409:PRO:C	1:D:454:TYR:CE1	2.71	0.53
1:E:23:SER:HG	1:E:483:ASN:HB3	1.70	0.53
1:A:289:ILE:HD12	1:A:289:ILE:N	2.24	0.53
1:A:397:GLY:HA2	1:A:403:CYS:SG	2.49	0.53
1:C:150:ILE:HG22	1:C:167:TRP:HZ3	1.72	0.53
1:C:440:ASN:C	1:C:440:ASN:HD22	2.17	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:451:ASP:OD2	1:C:471:ALA:N	2.42	0.53
1:F:245:ALA:C	1:F:247:GLY:H	2.15	0.53
1:F:408:SER:CB	1:F:451:ASP:HB3	2.38	0.53
1:C:371:GLN:HB3	1:C:484:VAL:HG23	1.90	0.53
1:D:371:GLN:HB3	1:D:484:VAL:HG23	1.90	0.53
1:D:440:ASN:HD22	1:D:440:ASN:C	2.17	0.53
1:D:451:ASP:OD2	1:D:471:ALA:N	2.42	0.53
1:B:228:GLY:HA2	1:B:345:SER:CB	2.32	0.53
1:B:397:GLY:HA2	1:B:403:CYS:SG	2.49	0.53
1:C:228:GLY:HA2	1:C:345:SER:H	1.74	0.53
1:B:172:SER:O	1:B:174:SER:HA	2.08	0.53
1:D:172:SER:O	1:D:174:SER:HA	2.08	0.53
1:E:371:GLN:HB3	1:E:484:VAL:HG23	1.90	0.53
1:E:408:SER:CB	1:E:451:ASP:HB3	2.39	0.53
1:F:371:GLN:HB3	1:F:484:VAL:HG23	1.90	0.53
1:A:228:GLY:HA2	1:A:345:SER:H	1.74	0.52
1:A:451:ASP:OD2	1:A:471:ALA:N	2.42	0.52
1:A:509:THR:N	1:A:510:PRO:HA	2.14	0.52
1:B:245:ALA:C	1:B:247:GLY:H	2.15	0.52
1:E:374:ILE:HG23	1:E:472:ALA:HA	1.91	0.52
1:B:228:GLY:HA2	1:B:345:SER:H	1.74	0.52
1:C:275:THR:OG1	1:C:278:GLN:HG3	2.09	0.52
1:A:408:SER:CB	1:A:451:ASP:HB3	2.39	0.52
1:D:502:VAL:HG12	1:D:504:LYS:H	1.75	0.52
1:E:399:VAL:O	1:E:399:VAL:HG12	2.09	0.52
1:F:23:SER:OG	1:F:483:ASN:HB2	2.04	0.52
1:F:502:VAL:HG12	1:F:504:LYS:H	1.75	0.52
1:B:173:SER:N	1:B:174:SER:HA	2.22	0.52
1:B:440:ASN:C	1:B:440:ASN:HD22	2.17	0.52
1:D:245:ALA:C	1:D:247:GLY:H	2.15	0.52
1:E:440:ASN:C	1:E:440:ASN:HD22	2.17	0.52
1:E:451:ASP:OD2	1:E:471:ALA:N	2.42	0.52
1:C:198:ILE:HG23	1:C:201:ALA:HB2	1.92	0.52
1:D:198:ILE:HG23	1:D:201:ALA:HB2	1.92	0.52
1:D:374:ILE:HG23	1:D:472:ALA:HA	1.91	0.52
1:E:397:GLY:HA2	1:E:403:CYS:SG	2.49	0.52
1:F:198:ILE:HG23	1:F:201:ALA:HB2	1.92	0.52
1:F:228:GLY:HA2	1:F:345:SER:H	1.74	0.52
1:F:275:THR:OG1	1:F:278:GLN:HG3	2.10	0.52
1:A:51:VAL:HG22	1:A:65:PHE:CZ	2.45	0.52
1:A:172:SER:O	1:A:174:SER:HA	2.08	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:275:THR:OG1	1:A:278:GLN:HG3	2.09	0.52
1:B:29:LEU:O	1:B:80:VAL:HA	2.10	0.52
1:C:245:ALA:C	1:C:247:GLY:H	2.16	0.52
1:D:397:GLY:HA2	1:D:403:CYS:SG	2.49	0.52
1:E:172:SER:O	1:E:174:SER:HA	2.08	0.52
1:E:173:SER:N	1:E:174:SER:HA	2.22	0.52
1:A:198:ILE:HG23	1:A:201:ALA:HB2	1.92	0.52
1:E:228:GLY:HA2	1:E:345:SER:N	2.25	0.52
1:E:275:THR:OG1	1:E:278:GLN:HG3	2.09	0.52
1:E:391:LYS:HE3	1:E:441:PHE:HA	1.92	0.52
1:F:51:VAL:HG22	1:F:65:PHE:HZ	1.73	0.52
1:F:289:ILE:HD12	1:F:289:ILE:N	2.24	0.52
1:F:391:LYS:HE3	1:F:441:PHE:HA	1.92	0.52
1:A:440:ASN:HD22	1:A:440:ASN:C	2.17	0.52
1:C:391:LYS:HZ2	1:C:440:ASN:ND2	2.03	0.52
1:D:275:THR:OG1	1:D:278:GLN:HG3	2.09	0.52
1:A:502:VAL:HG12	1:A:504:LYS:H	1.75	0.52
1:B:51:VAL:HG22	1:B:65:PHE:CZ	2.45	0.52
1:B:399:VAL:HG12	1:B:399:VAL:O	2.09	0.52
1:B:408:SER:CB	1:B:451:ASP:HB3	2.39	0.52
1:C:173:SER:N	1:C:174:SER:HA	2.22	0.52
1:C:391:LYS:HE3	1:C:441:PHE:HA	1.92	0.52
1:C:397:GLY:HA2	1:C:403:CYS:SG	2.49	0.52
1:D:228:GLY:HA2	1:D:345:SER:N	2.25	0.52
1:D:391:LYS:HE3	1:D:441:PHE:HA	1.92	0.52
1:F:51:VAL:HG22	1:F:65:PHE:CZ	2.45	0.52
1:F:172:SER:O	1:F:174:SER:HA	2.08	0.52
1:F:228:GLY:HA2	1:F:345:SER:N	2.25	0.52
1:F:374:ILE:HG23	1:F:472:ALA:HA	1.91	0.52
1:A:374:ILE:HG23	1:A:472:ALA:HA	1.91	0.52
1:B:84:VAL:HG13	1:B:89:ALA:HB2	1.92	0.52
1:B:352:ALA:O	1:B:355:LEU:HB2	2.10	0.52
1:B:374:ILE:HG23	1:B:472:ALA:HA	1.91	0.52
1:E:228:GLY:HA2	1:E:345:SER:H	1.74	0.52
1:F:397:GLY:HA2	1:F:403:CYS:SG	2.49	0.52
1:B:198:ILE:HG23	1:B:201:ALA:HB2	1.92	0.51
1:B:454:TYR:HE2	1:B:469:PRO:HB3	1.75	0.51
1:D:29:LEU:O	1:D:80:VAL:HA	2.10	0.51
1:E:198:ILE:HG23	1:E:201:ALA:HB2	1.92	0.51
1:D:411:ARG:O	1:D:412:GLU:C	2.54	0.51
1:D:454:TYR:HE2	1:D:469:PRO:HB3	1.75	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:84:VAL:HG13	1:E:89:ALA:HB2	1.92	0.51
1:A:237:GLU:HG3	1:A:337:ILE:CD1	2.41	0.51
1:A:411:ARG:O	1:A:412:GLU:C	2.54	0.51
1:B:275:THR:OG1	1:B:278:GLN:HG3	2.10	0.51
1:C:29:LEU:O	1:C:80:VAL:HA	2.10	0.51
1:C:237:GLU:HG3	1:C:337:ILE:CD1	2.41	0.51
1:C:399:VAL:O	1:C:399:VAL:HG12	2.09	0.51
1:D:173:SER:N	1:D:174:SER:HA	2.22	0.51
1:E:51:VAL:HG22	1:E:65:PHE:CZ	2.45	0.51
1:F:451:ASP:OD2	1:F:471:ALA:N	2.42	0.51
1:C:453:ASN:HD22	1:C:453:ASN:H	1.58	0.51
1:F:440:ASN:HD22	1:F:440:ASN:C	2.17	0.51
1:B:502:VAL:HG12	1:B:504:LYS:H	1.75	0.51
1:C:25:GLY:HA2	1:C:484:VAL:HG21	1.93	0.51
1:C:51:VAL:HG22	1:C:65:PHE:CZ	2.45	0.51
1:D:51:VAL:HG22	1:D:65:PHE:CZ	2.45	0.51
1:D:228:GLY:HA2	1:D:345:SER:H	1.74	0.51
1:E:352:ALA:O	1:E:355:LEU:HB2	2.10	0.51
1:F:352:ALA:O	1:F:355:LEU:HB2	2.10	0.51
1:A:391:LYS:HE3	1:A:441:PHE:HA	1.92	0.51
1:B:391:LYS:HE3	1:B:441:PHE:HA	1.92	0.51
1:B:453:ASN:H	1:B:453:ASN:HD22	1.58	0.51
1:C:228:GLY:HA2	1:C:345:SER:N	2.25	0.51
1:C:352:ALA:O	1:C:355:LEU:HB2	2.11	0.51
1:C:454:TYR:HE2	1:C:469:PRO:HB3	1.75	0.51
1:A:30:ALA:HB3	1:A:359:TRP:CD2	2.46	0.51
1:B:25:GLY:HA2	1:B:484:VAL:HG21	1.93	0.51
1:B:237:GLU:HG3	1:B:337:ILE:CD1	2.41	0.51
1:B:411:ARG:O	1:B:412:GLU:C	2.54	0.51
1:E:289:ILE:HD12	1:E:289:ILE:N	2.24	0.51
1:F:71:PHE:C	1:F:73:GLN:H	2.19	0.51
1:F:399:VAL:O	1:F:399:VAL:HG12	2.09	0.51
1:A:228:GLY:HA2	1:A:345:SER:N	2.25	0.51
1:C:84:VAL:HG13	1:C:89:ALA:HB2	1.92	0.51
1:C:502:VAL:HG12	1:C:504:LYS:H	1.75	0.51
1:C:508:GLU:CG	1:C:509:THR:HG22	2.36	0.51
1:D:71:PHE:C	1:D:73:GLN:H	2.19	0.51
1:D:161:PRO:O	1:D:186:LYS:HB3	2.11	0.51
1:D:399:VAL:O	1:D:399:VAL:HG12	2.09	0.51
1:E:71:PHE:C	1:E:73:GLN:H	2.19	0.51
1:E:411:ARG:O	1:E:412:GLU:C	2.54	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:194:LEU:HD23	1:F:195:LEU:N	2.26	0.51
1:A:29:LEU:O	1:A:80:VAL:HA	2.10	0.51
1:A:84:VAL:HG13	1:A:89:ALA:HB2	1.92	0.51
1:A:194:LEU:HD23	1:A:195:LEU:N	2.26	0.51
1:D:352:ALA:O	1:D:355:LEU:HB2	2.10	0.51
1:E:71:PHE:C	1:E:73:GLN:N	2.69	0.51
1:E:454:TYR:HE2	1:E:469:PRO:HB3	1.75	0.51
1:E:502:VAL:HG12	1:E:504:LYS:H	1.75	0.51
1:B:71:PHE:C	1:B:73:GLN:H	2.19	0.51
1:B:228:GLY:HA2	1:B:345:SER:N	2.25	0.51
1:D:194:LEU:HD23	1:D:195:LEU:N	2.26	0.51
1:D:300:GLY:O	1:D:302:LYS:HG3	2.11	0.51
1:D:453:ASN:HD22	1:D:453:ASN:H	1.58	0.51
1:E:161:PRO:O	1:E:186:LYS:HB3	2.11	0.51
1:B:30:ALA:HB3	1:B:359:TRP:CD2	2.46	0.50
1:C:23:SER:OG	1:C:483:ASN:HB2	2.04	0.50
1:C:30:ALA:HB3	1:C:359:TRP:CD2	2.46	0.50
1:D:30:ALA:HB3	1:D:359:TRP:CD2	2.46	0.50
1:E:25:GLY:HA2	1:E:484:VAL:HG21	1.93	0.50
1:F:450:ILE:CG1	1:F:451:ASP:N	2.53	0.50
1:A:352:ALA:O	1:A:355:LEU:HB2	2.11	0.50
1:A:399:VAL:O	1:A:399:VAL:HG12	2.09	0.50
1:A:453:ASN:HD22	1:A:453:ASN:H	1.58	0.50
1:B:220:PRO:HD2	1:B:338:LEU:HD11	1.94	0.50
1:C:374:ILE:HG23	1:C:472:ALA:HA	1.91	0.50
1:C:411:ARG:O	1:C:412:GLU:C	2.54	0.50
1:F:29:LEU:O	1:F:80:VAL:HA	2.10	0.50
1:F:84:VAL:HG13	1:F:89:ALA:HB2	1.92	0.50
1:F:453:ASN:HD22	1:F:453:ASN:H	1.58	0.50
1:A:220:PRO:HD2	1:A:338:LEU:HD11	1.94	0.50
1:D:84:VAL:HG13	1:D:89:ALA:HB2	1.92	0.50
1:E:237:GLU:HG3	1:E:337:ILE:CD1	2.41	0.50
1:F:71:PHE:C	1:F:73:GLN:N	2.69	0.50
1:F:220:PRO:HD2	1:F:338:LEU:HD11	1.94	0.50
1:D:237:GLU:HG3	1:D:337:ILE:CD1	2.41	0.50
1:D:289:ILE:HD12	1:D:289:ILE:N	2.24	0.50
1:D:407:CYS:N	1:D:449:ALA:O	2.45	0.50
1:E:29:LEU:O	1:E:80:VAL:HA	2.10	0.50
1:E:300:GLY:O	1:E:302:LYS:HG3	2.11	0.50
1:F:30:ALA:HB3	1:F:359:TRP:CD2	2.46	0.50
1:F:411:ARG:O	1:F:412:GLU:C	2.54	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:161:PRO:O	1:C:186:LYS:HB3	2.11	0.50
1:C:194:LEU:HD23	1:C:195:LEU:N	2.26	0.50
1:C:450:ILE:CG1	1:C:451:ASP:N	2.53	0.50
1:E:30:ALA:HB3	1:E:359:TRP:CD2	2.46	0.50
1:E:194:LEU:HD23	1:E:195:LEU:N	2.26	0.50
1:F:454:TYR:HE2	1:F:469:PRO:HB3	1.75	0.50
1:B:161:PRO:O	1:B:186:LYS:HB3	2.11	0.50
1:B:194:LEU:HD23	1:B:195:LEU:N	2.26	0.50
1:C:71:PHE:C	1:C:73:GLN:H	2.19	0.50
1:C:228:GLY:HA2	1:C:345:SER:CB	2.32	0.50
1:F:70:ASN:HB3	1:F:457:GLN:HE22	1.76	0.50
1:A:454:TYR:HE2	1:A:469:PRO:HB3	1.75	0.50
1:C:220:PRO:HD2	1:C:338:LEU:HD11	1.94	0.50
1:D:70:ASN:HB3	1:D:457:GLN:HE22	1.76	0.50
1:E:453:ASN:H	1:E:453:ASN:HD22	1.58	0.50
1:F:237:GLU:HG3	1:F:337:ILE:CD1	2.41	0.50
1:F:508:GLU:CG	1:F:509:THR:HG22	2.36	0.50
1:C:300:GLY:O	1:C:302:LYS:HG3	2.11	0.50
1:C:407:CYS:N	1:C:449:ALA:O	2.44	0.50
1:D:25:GLY:HA2	1:D:484:VAL:HG21	1.93	0.50
1:D:220:PRO:HD2	1:D:338:LEU:HD11	1.94	0.50
1:D:362:PHE:HA	1:D:368:VAL:HG21	1.94	0.50
1:A:71:PHE:C	1:A:73:GLN:H	2.19	0.49
1:B:254:ILE:HG12	1:B:337:ILE:HB	1.93	0.49
1:B:300:GLY:O	1:B:302:LYS:HG3	2.11	0.49
1:F:25:GLY:HA2	1:F:484:VAL:HG21	1.93	0.49
1:F:236:ILE:HD12	1:F:236:ILE:N	2.27	0.49
1:A:300:GLY:O	1:A:302:LYS:HG3	2.11	0.49
1:B:70:ASN:HB3	1:B:457:GLN:HE22	1.76	0.49
1:C:70:ASN:HB3	1:C:457:GLN:HE22	1.76	0.49
1:E:236:ILE:HD12	1:E:236:ILE:N	2.27	0.49
1:F:161:PRO:O	1:F:186:LYS:HB3	2.11	0.49
1:F:362:PHE:HA	1:F:368:VAL:HG21	1.94	0.49
1:A:25:GLY:HA2	1:A:484:VAL:HG21	1.93	0.49
1:A:161:PRO:O	1:A:186:LYS:HB3	2.11	0.49
1:A:236:ILE:HD12	1:A:236:ILE:N	2.27	0.49
1:A:109:ASN:O	1:A:177:GLY:HA3	2.13	0.49
1:A:254:ILE:HG12	1:A:337:ILE:HB	1.93	0.49
1:A:362:PHE:HA	1:A:368:VAL:HG21	1.94	0.49
1:C:362:PHE:HA	1:C:368:VAL:HG21	1.94	0.49
1:D:71:PHE:C	1:D:73:GLN:N	2.69	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:220:PRO:HD2	1:E:338:LEU:HD11	1.94	0.49
1:B:407:CYS:N	1:B:449:ALA:O	2.44	0.49
1:C:71:PHE:C	1:C:73:GLN:N	2.69	0.49
1:D:383:LEU:O	1:D:386:ALA:HB3	2.13	0.49
1:E:382:SER:CB	1:E:385:THR:HG22	2.42	0.49
1:E:383:LEU:O	1:E:386:ALA:HB3	2.13	0.49
1:F:383:LEU:O	1:F:386:ALA:HB3	2.13	0.49
1:A:70:ASN:HB3	1:A:457:GLN:HE22	1.76	0.49
1:C:289:ILE:HD12	1:C:289:ILE:N	2.24	0.49
1:C:304:ILE:HG13	1:C:305:TYR:CD2	2.47	0.49
1:D:254:ILE:HG12	1:D:337:ILE:HB	1.93	0.49
1:E:70:ASN:HB3	1:E:457:GLN:HE22	1.76	0.49
1:E:228:GLY:HA2	1:E:345:SER:CB	2.32	0.49
1:F:109:ASN:O	1:F:177:GLY:HA3	2.13	0.49
1:F:254:ILE:HG12	1:F:337:ILE:HB	1.93	0.49
1:A:71:PHE:C	1:A:73:GLN:N	2.69	0.49
1:A:178:LEU:CD2	1:A:178:LEU:N	2.66	0.49
1:F:300:GLY:O	1:F:302:LYS:HG3	2.11	0.49
1:A:304:ILE:HG13	1:A:305:TYR:CD2	2.47	0.49
1:E:376:GLY:HA2	1:E:390:GLN:NE2	2.28	0.49
1:E:456:TYR:HB2	1:E:467:TRP:CZ3	2.48	0.49
1:B:109:ASN:O	1:B:177:GLY:HA3	2.13	0.49
1:B:236:ILE:HD12	1:B:236:ILE:N	2.27	0.49
1:B:304:ILE:HG13	1:B:305:TYR:CD2	2.47	0.49
1:D:456:TYR:HB2	1:D:467:TRP:CZ3	2.48	0.49
1:F:376:GLY:HA2	1:F:390:GLN:NE2	2.28	0.49
1:B:383:LEU:O	1:B:386:ALA:HB3	2.13	0.49
1:D:109:ASN:O	1:D:177:GLY:HA3	2.13	0.49
1:E:109:ASN:O	1:E:177:GLY:HA3	2.13	0.49
1:E:445:SER:HB3	1:E:448:ALA:HB2	1.95	0.49
1:F:456:TYR:HB2	1:F:467:TRP:CZ3	2.48	0.49
1:A:456:TYR:HB2	1:A:467:TRP:CZ3	2.48	0.48
1:B:215:LYS:CE	1:B:329:ASN:HD21	2.25	0.48
1:D:236:ILE:HD12	1:D:236:ILE:N	2.27	0.48
1:D:304:ILE:HG13	1:D:305:TYR:CD2	2.47	0.48
1:E:407:CYS:N	1:E:449:ALA:O	2.44	0.48
1:B:289:ILE:HD12	1:B:289:ILE:N	2.24	0.48
1:B:453:ASN:H	1:B:453:ASN:ND2	2.12	0.48
1:A:408:SER:HA	1:A:451:ASP:O	2.14	0.48
1:C:236:ILE:HD12	1:C:236:ILE:N	2.27	0.48
1:C:254:ILE:HG12	1:C:337:ILE:HB	1.93	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:508:GLU:CG	1:E:509:THR:HG22	2.36	0.48
1:A:508:GLU:CG	1:A:509:THR:HG22	2.36	0.48
1:B:23:SER:HG	1:B:483:ASN:HB3	1.77	0.48
1:C:453:ASN:H	1:C:453:ASN:ND2	2.11	0.48
1:E:254:ILE:HG12	1:E:337:ILE:HB	1.93	0.48
1:E:304:ILE:HG13	1:E:305:TYR:CD2	2.47	0.48
1:E:453:ASN:H	1:E:453:ASN:ND2	2.11	0.48
1:F:445:SER:HB3	1:F:448:ALA:HB2	1.95	0.48
1:F:453:ASN:H	1:F:453:ASN:ND2	2.11	0.48
1:C:376:GLY:HA2	1:C:390:GLN:NE2	2.28	0.48
1:C:383:LEU:O	1:C:386:ALA:HB3	2.13	0.48
1:C:456:TYR:HB2	1:C:467:TRP:CZ3	2.48	0.48
1:E:63:ASP:O	1:E:67:SER:HB2	2.14	0.48
1:A:453:ASN:H	1:A:453:ASN:ND2	2.11	0.48
1:A:509:THR:N	1:A:510:PRO:CA	2.73	0.48
1:B:376:GLY:HA2	1:B:390:GLN:NE2	2.28	0.48
1:B:413:THR:OG1	1:B:425:ASN:HB3	2.14	0.48
1:B:456:TYR:HB2	1:B:467:TRP:CZ3	2.48	0.48
1:B:508:GLU:HA	1:B:509:THR:HA	1.59	0.48
1:E:408:SER:HA	1:E:451:ASP:O	2.14	0.48
1:F:304:ILE:HG13	1:F:305:TYR:CD2	2.47	0.48
1:F:407:CYS:N	1:F:449:ALA:O	2.45	0.48
1:B:71:PHE:C	1:B:73:GLN:N	2.69	0.48
1:C:408:SER:HA	1:C:451:ASP:O	2.14	0.48
1:D:408:SER:HA	1:D:451:ASP:O	2.14	0.48
1:D:456:TYR:CZ	1:D:465:ASN:HB3	2.49	0.48
1:A:284:ARG:HA	1:A:288:ALA:O	2.14	0.48
1:D:284:ARG:HA	1:D:288:ALA:O	2.14	0.48
1:E:362:PHE:HA	1:E:368:VAL:HG21	1.94	0.48
1:A:376:GLY:HA2	1:A:390:GLN:NE2	2.28	0.48
1:A:407:CYS:N	1:A:449:ALA:O	2.44	0.48
1:A:456:TYR:CZ	1:A:465:ASN:HB3	2.49	0.48
1:B:125:ILE:HD12	1:B:153:LYS:HG2	1.96	0.48
1:B:284:ARG:HA	1:B:288:ALA:O	2.14	0.48
1:B:456:TYR:CZ	1:B:465:ASN:HB3	2.49	0.48
1:B:509:THR:N	1:B:510:PRO:CA	2.73	0.48
1:C:109:ASN:O	1:C:177:GLY:HA3	2.13	0.48
1:D:63:ASP:O	1:D:67:SER:HB2	2.14	0.48
1:D:453:ASN:H	1:D:453:ASN:ND2	2.11	0.48
1:B:362:PHE:HA	1:B:368:VAL:HG21	1.94	0.48
1:C:285:ARG:O	1:C:286:ASN:HB2	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:413:THR:OG1	1:C:425:ASN:HB3	2.14	0.48
1:C:456:TYR:CZ	1:C:465:ASN:HB3	2.49	0.48
1:D:502:VAL:HG12	1:D:503:ILE:N	2.29	0.48
1:E:58:THR:H	1:E:61:THR:HB	1.79	0.48
1:E:125:ILE:HD12	1:E:153:LYS:HG2	1.96	0.48
1:E:284:ARG:HA	1:E:288:ALA:O	2.14	0.48
1:E:456:TYR:CZ	1:E:465:ASN:HB3	2.49	0.48
1:A:58:THR:H	1:A:61:THR:HB	1.79	0.47
1:A:63:ASP:O	1:A:67:SER:HB2	2.14	0.47
1:B:30:ALA:HB3	1:B:359:TRP:CE2	2.49	0.47
1:C:30:ALA:HB3	1:C:359:TRP:CE2	2.49	0.47
1:D:30:ALA:HB3	1:D:359:TRP:CE2	2.49	0.47
1:D:228:GLY:HA2	1:D:345:SER:CB	2.32	0.47
1:D:413:THR:OG1	1:D:425:ASN:HB3	2.14	0.47
1:E:450:ILE:CG1	1:E:451:ASP:N	2.53	0.47
1:F:63:ASP:O	1:F:67:SER:HB2	2.14	0.47
1:F:382:SER:CB	1:F:385:THR:HG22	2.42	0.47
1:F:408:SER:HA	1:F:451:ASP:O	2.14	0.47
1:A:383:LEU:O	1:A:386:ALA:HB3	2.13	0.47
1:A:450:ILE:CG1	1:A:451:ASP:N	2.53	0.47
1:C:125:ILE:HD12	1:C:153:LYS:HG2	1.96	0.47
1:D:58:THR:H	1:D:61:THR:HB	1.79	0.47
1:D:285:ARG:O	1:D:286:ASN:HB2	2.14	0.47
1:E:30:ALA:HB3	1:E:359:TRP:CE2	2.49	0.47
1:E:409:PRO:C	1:E:454:TYR:CE1	2.71	0.47
1:F:58:THR:H	1:F:61:THR:HB	1.79	0.47
1:F:413:THR:OG1	1:F:425:ASN:HB3	2.14	0.47
1:A:445:SER:HB3	1:A:448:ALA:HB2	1.95	0.47
1:D:125:ILE:HD12	1:D:153:LYS:HG2	1.96	0.47
1:D:376:GLY:HA2	1:D:390:GLN:NE2	2.28	0.47
1:D:391:LYS:HZ1	1:D:440:ASN:ND2	2.09	0.47
1:D:445:SER:HB3	1:D:448:ALA:HB2	1.95	0.47
1:F:30:ALA:HB3	1:F:359:TRP:CE2	2.49	0.47
1:F:125:ILE:HD12	1:F:153:LYS:HG2	1.96	0.47
1:A:30:ALA:HB3	1:A:359:TRP:CE2	2.49	0.47
1:A:407:CYS:O	1:A:451:ASP:HB3	2.15	0.47
1:B:408:SER:HA	1:B:451:ASP:O	2.14	0.47
1:C:215:LYS:CE	1:C:329:ASN:HD21	2.25	0.47
1:A:413:THR:OG1	1:A:425:ASN:HB3	2.14	0.47
1:A:453:ASN:ND2	1:A:453:ASN:N	2.62	0.47
1:B:63:ASP:O	1:B:67:SER:HB2	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:499:ILE:HD13	1:B:499:ILE:N	2.28	0.47
1:C:502:VAL:HG12	1:C:503:ILE:N	2.29	0.47
1:A:125:ILE:HD12	1:A:153:LYS:HG2	1.96	0.47
1:B:285:ARG:O	1:B:286:ASN:HB2	2.14	0.47
1:C:284:ARG:HA	1:C:288:ALA:O	2.14	0.47
1:D:68:ALA:O	1:D:69:MET:C	2.58	0.47
1:D:508:GLU:CG	1:D:509:THR:HG22	2.36	0.47
1:F:456:TYR:CZ	1:F:465:ASN:HB3	2.49	0.47
1:A:379:ALA:HB2	1:A:454:TYR:CE2	2.46	0.47
1:A:382:SER:CB	1:A:385:THR:HG22	2.42	0.47
1:B:58:THR:H	1:B:61:THR:HB	1.79	0.47
1:B:171:ILE:CG2	1:B:172:SER:H	2.25	0.47
1:B:350:VAL:HG13	1:B:354:ASP:HB2	1.97	0.47
1:B:382:SER:CB	1:B:385:THR:HG22	2.42	0.47
1:B:407:CYS:O	1:B:451:ASP:HB3	2.14	0.47
1:C:27:ALA:HB2	1:C:71:PHE:CZ	2.50	0.47
1:C:379:ALA:HB2	1:C:454:TYR:CE2	2.46	0.47
1:C:445:SER:HB3	1:C:448:ALA:HB2	1.95	0.47
1:E:413:THR:OG1	1:E:425:ASN:HB3	2.14	0.47
1:E:453:ASN:ND2	1:E:453:ASN:N	2.62	0.47
1:F:285:ARG:O	1:F:286:ASN:HB2	2.14	0.47
1:F:407:CYS:O	1:F:451:ASP:N	2.48	0.47
1:A:285:ARG:O	1:A:286:ASN:HB2	2.14	0.47
1:B:502:VAL:HG12	1:B:503:ILE:N	2.29	0.47
1:C:63:ASP:O	1:C:67:SER:HB2	2.14	0.47
1:D:197:GLU:H	1:D:197:GLU:CD	2.23	0.47
1:D:215:LYS:CE	1:D:329:ASN:HD21	2.25	0.47
1:E:407:CYS:O	1:E:451:ASP:N	2.48	0.47
1:A:350:VAL:HG13	1:A:354:ASP:HB2	1.97	0.47
1:C:407:CYS:O	1:C:451:ASP:N	2.48	0.47
1:D:27:ALA:HB2	1:D:71:PHE:CZ	2.50	0.47
1:D:350:VAL:HG13	1:D:354:ASP:HB2	1.97	0.47
1:A:100:GLU:HG2	1:A:186:LYS:O	2.15	0.47
1:A:293:VAL:HG22	1:A:294:VAL:N	2.30	0.47
1:B:379:ALA:HB2	1:B:454:TYR:CE2	2.46	0.47
1:B:445:SER:HB3	1:B:448:ALA:HB2	1.95	0.47
1:C:23:SER:HG	1:C:483:ASN:HB3	1.77	0.47
1:C:290:VAL:HG11	1:C:322:TYR:CD1	2.50	0.47
1:E:502:VAL:HG12	1:E:503:ILE:N	2.29	0.47
1:F:27:ALA:HB2	1:F:71:PHE:CZ	2.50	0.47
1:F:197:GLU:H	1:F:197:GLU:CD	2.23	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:407:CYS:O	1:A:451:ASP:N	2.48	0.46
1:B:27:ALA:HB2	1:B:71:PHE:CZ	2.50	0.46
1:B:290:VAL:HG11	1:B:322:TYR:CD1	2.50	0.46
1:B:508:GLU:CG	1:B:509:THR:HG22	2.36	0.46
1:C:350:VAL:HG13	1:C:354:ASP:HB2	1.97	0.46
1:C:453:ASN:ND2	1:C:453:ASN:N	2.62	0.46
1:E:163:LEU:HB3	1:E:164:GLY:H	1.63	0.46
1:F:100:GLU:HG2	1:F:186:LYS:O	2.16	0.46
1:F:502:VAL:HG12	1:F:503:ILE:N	2.29	0.46
1:B:453:ASN:ND2	1:B:453:ASN:N	2.62	0.46
1:D:290:VAL:HG11	1:D:322:TYR:CD1	2.50	0.46
1:E:62:ALA:HB1	1:E:466:ARG:NE	2.30	0.46
1:E:215:LYS:CE	1:E:329:ASN:HD21	2.25	0.46
1:F:284:ARG:HA	1:F:288:ALA:O	2.14	0.46
1:F:293:VAL:HG22	1:F:294:VAL:N	2.30	0.46
1:F:379:ALA:HB2	1:F:454:TYR:CE2	2.46	0.46
1:A:27:ALA:HB2	1:A:71:PHE:CZ	2.50	0.46
1:A:68:ALA:O	1:A:69:MET:C	2.58	0.46
1:A:151:ILE:HG13	1:A:152:ALA:N	2.30	0.46
1:B:100:GLU:HG2	1:B:186:LYS:O	2.16	0.46
1:B:423:VAL:O	1:B:424:ASP:C	2.59	0.46
1:C:58:THR:H	1:C:61:THR:HB	1.79	0.46
1:C:197:GLU:CD	1:C:197:GLU:H	2.23	0.46
1:C:293:VAL:HG22	1:C:294:VAL:N	2.30	0.46
1:C:382:SER:CB	1:C:385:THR:HG22	2.42	0.46
1:D:23:SER:OG	1:D:483:ASN:HB2	2.04	0.46
1:D:62:ALA:HB1	1:D:466:ARG:NE	2.30	0.46
1:E:208:VAL:HG23	1:E:209:ASP:N	2.31	0.46
1:F:68:ALA:O	1:F:69:MET:C	2.58	0.46
1:C:37:PRO:HB2	1:C:40:GLN:HB2	1.97	0.46
1:D:56:GLN:HA	1:D:57:PRO:HD3	1.81	0.46
1:D:293:VAL:HG22	1:D:294:VAL:N	2.30	0.46
1:D:407:CYS:O	1:D:451:ASP:N	2.48	0.46
1:E:27:ALA:HB2	1:E:71:PHE:CZ	2.50	0.46
1:F:509:THR:N	1:F:510:PRO:CA	2.73	0.46
1:A:37:PRO:HB2	1:A:40:GLN:HB2	1.97	0.46
1:A:502:VAL:HG12	1:A:503:ILE:N	2.29	0.46
1:B:407:CYS:O	1:B:451:ASP:N	2.48	0.46
1:D:35:TRP:HB3	1:D:54:PHE:HA	1.98	0.46
1:E:200:ASN:O	1:E:201:ALA:C	2.59	0.46
1:E:285:ARG:O	1:E:286:ASN:HB2	2.14	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:62:ALA:HB1	1:B:466:ARG:NE	2.30	0.46
1:B:200:ASN:O	1:B:201:ALA:C	2.59	0.46
1:B:293:VAL:HG22	1:B:294:VAL:N	2.30	0.46
1:C:509:THR:CB	1:C:510:PRO:HA	2.43	0.46
1:F:37:PRO:HB2	1:F:40:GLN:HB2	1.97	0.46
1:F:208:VAL:HG23	1:F:209:ASP:N	2.31	0.46
1:F:458:TYR:CE2	1:F:460:LYS:HA	2.51	0.46
1:A:27:ALA:HB3	1:A:78:LEU:HD12	1.98	0.46
1:A:509:THR:CB	1:A:510:PRO:HA	2.43	0.46
1:B:197:GLU:CD	1:B:197:GLU:H	2.23	0.46
1:C:100:GLU:HG2	1:C:186:LYS:O	2.16	0.46
1:C:499:ILE:HD13	1:C:499:ILE:N	2.28	0.46
1:D:130:LYS:O	1:D:132:THR:HG23	2.16	0.46
1:D:453:ASN:ND2	1:D:453:ASN:N	2.62	0.46
1:E:350:VAL:HG13	1:E:354:ASP:HB2	1.97	0.46
1:F:290:VAL:HG11	1:F:322:TYR:CD1	2.50	0.46
1:F:350:VAL:HG13	1:F:354:ASP:HB2	1.97	0.46
1:A:173:SER:HA	1:A:174:SER:HB3	1.98	0.46
1:A:499:ILE:HD13	1:A:499:ILE:N	2.28	0.46
1:B:73:GLN:HB3	1:B:500:LEU:HD12	1.98	0.46
1:B:458:TYR:CE2	1:B:460:LYS:HA	2.51	0.46
1:C:208:VAL:HG23	1:C:209:ASP:N	2.31	0.46
1:D:37:PRO:HB2	1:D:40:GLN:HB2	1.97	0.46
1:D:73:GLN:HB3	1:D:500:LEU:HD12	1.98	0.46
1:D:509:THR:N	1:D:510:PRO:CA	2.73	0.46
1:E:100:GLU:HG2	1:E:186:LYS:O	2.16	0.46
1:F:381:GLU:O	1:F:382:SER:C	2.59	0.46
1:A:62:ALA:HB1	1:A:466:ARG:NE	2.30	0.46
1:A:290:VAL:HG11	1:A:322:TYR:CD1	2.50	0.46
1:D:100:GLU:HG2	1:D:186:LYS:O	2.15	0.46
1:E:381:GLU:O	1:E:382:SER:C	2.59	0.46
1:F:62:ALA:HB1	1:F:466:ARG:NE	2.30	0.46
1:F:130:LYS:O	1:F:132:THR:HG23	2.16	0.46
1:F:200:ASN:O	1:F:201:ALA:C	2.59	0.46
1:A:171:ILE:CG2	1:A:172:SER:N	2.79	0.46
1:B:27:ALA:HB3	1:B:78:LEU:HD12	1.98	0.46
1:B:178:LEU:HD23	1:B:178:LEU:N	2.22	0.46
1:C:35:TRP:HB3	1:C:54:PHE:HA	1.98	0.46
1:D:379:ALA:HB2	1:D:454:TYR:CE2	2.46	0.46
1:E:197:GLU:CD	1:E:197:GLU:H	2.23	0.46
1:E:379:ALA:HB2	1:E:454:TYR:CE2	2.46	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:211:GLN:OE1	1:F:328:GLN:HG2	2.16	0.46
1:F:215:LYS:CE	1:F:329:ASN:HD21	2.25	0.46
1:A:61:THR:O	1:A:62:ALA:C	2.59	0.45
1:A:458:TYR:CE2	1:A:460:LYS:HA	2.51	0.45
1:C:62:ALA:HB1	1:C:466:ARG:NE	2.30	0.45
1:C:458:TYR:CE2	1:C:460:LYS:HA	2.51	0.45
1:E:290:VAL:HG11	1:E:322:TYR:CD1	2.50	0.45
1:E:478:CYS:C	1:E:480:ARG:N	2.72	0.45
1:A:73:GLN:HB3	1:A:500:LEU:HD12	1.98	0.45
1:A:208:VAL:HG23	1:A:209:ASP:N	2.31	0.45
1:A:283:VAL:CG2	1:A:323:ILE:HD13	2.47	0.45
1:B:173:SER:HA	1:B:174:SER:HB3	1.98	0.45
1:C:385:THR:HA	1:C:388:THR:HG23	1.98	0.45
1:D:61:THR:O	1:D:62:ALA:C	2.59	0.45
1:D:173:SER:HA	1:D:174:SER:HB3	1.98	0.45
1:D:200:ASN:O	1:D:201:ALA:C	2.59	0.45
1:D:208:VAL:HG23	1:D:209:ASP:N	2.31	0.45
1:D:283:VAL:CG2	1:D:323:ILE:HD13	2.47	0.45
1:D:499:ILE:HD13	1:D:499:ILE:N	2.28	0.45
1:E:37:PRO:HB2	1:E:40:GLN:HB2	1.97	0.45
1:F:27:ALA:HB3	1:F:78:LEU:HD12	1.98	0.45
1:A:215:LYS:HE3	1:A:329:ASN:CG	2.42	0.45
1:A:423:VAL:O	1:A:424:ASP:C	2.59	0.45
1:B:37:PRO:HB2	1:B:40:GLN:HB2	1.97	0.45
1:B:419:VAL:HA	1:B:422:ALA:CB	2.46	0.45
1:C:73:GLN:HB3	1:C:500:LEU:HD12	1.98	0.45
1:D:385:THR:HA	1:D:388:THR:HG23	1.99	0.45
1:D:402:ASP:CG	1:D:402:ASP:O	2.59	0.45
1:D:458:TYR:CE2	1:D:460:LYS:HA	2.51	0.45
1:E:73:GLN:HB3	1:E:500:LEU:HD12	1.98	0.45
1:E:283:VAL:CG2	1:E:323:ILE:HD13	2.47	0.45
1:E:293:VAL:HG22	1:E:294:VAL:N	2.30	0.45
1:F:453:ASN:ND2	1:F:453:ASN:N	2.62	0.45
1:F:478:CYS:C	1:F:480:ARG:N	2.72	0.45
1:B:68:ALA:O	1:B:69:MET:C	2.58	0.45
1:C:50:LEU:HD12	1:C:51:VAL:N	2.32	0.45
1:C:200:ASN:O	1:C:201:ALA:C	2.59	0.45
1:C:423:VAL:O	1:C:424:ASP:C	2.59	0.45
1:E:35:TRP:HB3	1:E:54:PHE:HA	1.98	0.45
1:F:283:VAL:CG2	1:F:323:ILE:HD13	2.47	0.45
1:F:307:SER:O	1:F:309:ILE:HG23	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:215:LYS:CE	1:A:329:ASN:HD21	2.25	0.45
1:B:50:LEU:HD12	1:B:51:VAL:N	2.32	0.45
1:B:385:THR:HA	1:B:388:THR:HG23	1.98	0.45
1:B:402:ASP:O	1:B:402:ASP:CG	2.59	0.45
1:D:211:GLN:OE1	1:D:328:GLN:HG2	2.16	0.45
1:D:381:GLU:O	1:D:382:SER:C	2.59	0.45
1:E:173:SER:HA	1:E:174:SER:HB3	1.98	0.45
1:F:402:ASP:CG	1:F:402:ASP:O	2.59	0.45
1:F:419:VAL:HA	1:F:422:ALA:CB	2.46	0.45
1:B:130:LYS:O	1:B:132:THR:HG23	2.16	0.45
1:C:61:THR:O	1:C:62:ALA:C	2.59	0.45
1:C:66:MET:O	1:C:67:SER:C	2.60	0.45
1:D:66:MET:O	1:D:67:SER:C	2.60	0.45
1:D:215:LYS:HE3	1:D:329:ASN:CG	2.42	0.45
1:E:66:MET:O	1:E:67:SER:C	2.60	0.45
1:E:456:TYR:CE2	1:E:503:ILE:HD11	2.52	0.45
1:F:456:TYR:CE2	1:F:503:ILE:HD11	2.52	0.45
1:A:130:LYS:O	1:A:132:THR:HG23	2.16	0.45
1:A:197:GLU:H	1:A:197:GLU:CD	2.23	0.45
1:A:381:GLU:O	1:A:382:SER:C	2.59	0.45
1:B:215:LYS:HE3	1:B:329:ASN:CG	2.42	0.45
1:C:130:LYS:O	1:C:132:THR:HG23	2.16	0.45
1:D:456:TYR:CE2	1:D:503:ILE:HD11	2.52	0.45
1:E:61:THR:O	1:E:62:ALA:C	2.59	0.45
1:E:68:ALA:O	1:E:69:MET:C	2.58	0.45
1:F:73:GLN:HB3	1:F:500:LEU:HD12	1.98	0.45
1:F:173:SER:HA	1:F:174:SER:HB3	1.98	0.45
1:F:228:GLY:HA2	1:F:345:SER:CB	2.32	0.45
1:A:66:MET:O	1:A:67:SER:C	2.60	0.45
1:A:419:VAL:HA	1:A:422:ALA:CB	2.46	0.45
1:B:35:TRP:HB3	1:B:54:PHE:HA	1.98	0.45
1:C:68:ALA:O	1:C:69:MET:C	2.58	0.45
1:C:215:LYS:HE3	1:C:329:ASN:CG	2.42	0.45
1:D:50:LEU:HD12	1:D:51:VAL:N	2.31	0.45
1:D:304:ILE:HG13	1:D:305:TYR:HE2	1.81	0.45
1:D:307:SER:O	1:D:309:ILE:HG23	2.16	0.45
1:D:423:VAL:O	1:D:424:ASP:C	2.59	0.45
1:E:27:ALA:HB3	1:E:78:LEU:HD12	1.98	0.45
1:E:211:GLN:OE1	1:E:328:GLN:HG2	2.16	0.45
1:E:402:ASP:O	1:E:402:ASP:CG	2.59	0.45
1:F:66:MET:O	1:F:67:SER:C	2.60	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:163:LEU:HB3	1:F:164:GLY:H	1.63	0.45
1:F:423:VAL:O	1:F:424:ASP:C	2.59	0.45
1:A:385:THR:HA	1:A:388:THR:HG23	1.98	0.45
1:B:66:MET:O	1:B:67:SER:C	2.60	0.45
1:B:208:VAL:HG23	1:B:209:ASP:N	2.31	0.45
1:C:27:ALA:HB3	1:C:78:LEU:HD12	1.98	0.45
1:E:385:THR:HA	1:E:388:THR:HG23	1.98	0.45
1:F:35:TRP:HB3	1:F:54:PHE:HA	1.98	0.45
1:F:390:GLN:HE22	1:F:408:SER:H	1.65	0.45
1:A:211:GLN:OE1	1:A:328:GLN:HG2	2.16	0.45
1:A:307:SER:O	1:A:309:ILE:HG23	2.16	0.45
1:A:390:GLN:HE22	1:A:408:SER:H	1.65	0.45
1:A:456:TYR:CE2	1:A:503:ILE:HD11	2.52	0.45
1:B:61:THR:O	1:B:62:ALA:C	2.59	0.45
1:B:153:LYS:O	1:B:157:VAL:HG22	2.17	0.45
1:B:283:VAL:CG2	1:B:323:ILE:HD13	2.47	0.45
1:B:307:SER:O	1:B:309:ILE:HG23	2.16	0.45
1:C:283:VAL:CG2	1:C:323:ILE:HD13	2.47	0.45
1:C:304:ILE:HG13	1:C:305:TYR:HE2	1.81	0.45
1:D:171:ILE:CG2	1:D:172:SER:N	2.79	0.45
1:E:423:VAL:O	1:E:424:ASP:C	2.59	0.45
1:E:508:GLU:HA	1:E:509:THR:HA	1.59	0.45
1:F:61:THR:O	1:F:62:ALA:C	2.59	0.45
1:A:147:THR:O	1:A:151:ILE:HG23	2.17	0.44
1:A:200:ASN:O	1:A:201:ALA:C	2.59	0.44
1:B:51:VAL:HG13	1:B:55:GLY:C	2.42	0.44
1:C:456:TYR:CE2	1:C:503:ILE:HD11	2.52	0.44
1:E:130:LYS:O	1:E:132:THR:HG23	2.16	0.44
1:E:215:LYS:HE3	1:E:329:ASN:CG	2.42	0.44
1:E:458:TYR:CE2	1:E:460:LYS:HA	2.51	0.44
1:B:390:GLN:HE22	1:B:408:SER:H	1.65	0.44
1:B:509:THR:CB	1:B:510:PRO:HA	2.44	0.44
1:C:211:GLN:OE1	1:C:328:GLN:HG2	2.16	0.44
1:C:419:VAL:HA	1:C:422:ALA:CB	2.46	0.44
1:D:27:ALA:HB3	1:D:78:LEU:HD12	1.98	0.44
1:E:390:GLN:HE22	1:E:408:SER:H	1.65	0.44
1:F:147:THR:O	1:F:151:ILE:HG23	2.17	0.44
1:A:171:ILE:CG2	1:A:172:SER:H	2.25	0.44
1:B:381:GLU:O	1:B:382:SER:C	2.59	0.44
1:C:307:SER:O	1:C:309:ILE:HG23	2.16	0.44
1:C:381:GLU:O	1:C:382:SER:C	2.59	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:147:THR:O	1:E:151:ILE:HG23	2.17	0.44
1:F:151:ILE:HG13	1:F:152:ALA:N	2.30	0.44
1:F:153:LYS:O	1:F:157:VAL:HG22	2.17	0.44
1:A:153:LYS:O	1:A:157:VAL:HG22	2.17	0.44
1:B:211:GLN:OE1	1:B:328:GLN:HG2	2.16	0.44
1:C:153:LYS:O	1:C:157:VAL:HG22	2.17	0.44
1:C:390:GLN:HE22	1:C:408:SER:H	1.65	0.44
1:D:390:GLN:HE22	1:D:408:SER:H	1.65	0.44
1:E:229:GLU:CD	1:E:229:GLU:H	2.26	0.44
1:E:307:SER:O	1:E:309:ILE:HG23	2.16	0.44
1:E:454:TYR:O	1:E:467:TRP:CZ3	2.71	0.44
1:E:499:ILE:HD13	1:E:499:ILE:N	2.28	0.44
1:F:50:LEU:HD12	1:F:51:VAL:N	2.31	0.44
1:F:215:LYS:HE3	1:F:329:ASN:CG	2.42	0.44
1:F:454:TYR:O	1:F:467:TRP:CZ3	2.71	0.44
1:A:35:TRP:HB3	1:A:54:PHE:HA	1.98	0.44
1:B:253:PRO:HD2	1:B:336:GLY:HA2	2.00	0.44
1:C:44:VAL:HG11	1:C:50:LEU:HB3	1.99	0.44
1:C:151:ILE:HG13	1:C:152:ALA:N	2.30	0.44
1:C:173:SER:HA	1:C:174:SER:HB3	1.98	0.44
1:C:375:ALA:HB3	1:C:406:LEU:O	2.17	0.44
1:C:402:ASP:CG	1:C:402:ASP:O	2.59	0.44
1:D:51:VAL:HG13	1:D:55:GLY:C	2.42	0.44
1:D:375:ALA:HB3	1:D:406:LEU:O	2.17	0.44
1:F:229:GLU:H	1:F:229:GLU:CD	2.26	0.44
1:A:229:GLU:H	1:A:229:GLU:CD	2.26	0.44
1:A:402:ASP:O	1:A:402:ASP:CG	2.59	0.44
1:D:99:ILE:O	1:D:100:GLU:C	2.61	0.44
1:D:382:SER:CB	1:D:385:THR:HG22	2.42	0.44
1:E:51:VAL:HG13	1:E:55:GLY:C	2.42	0.44
1:E:252:LEU:HA	1:E:253:PRO:HD3	1.80	0.44
1:F:99:ILE:O	1:F:100:GLU:C	2.61	0.44
1:A:50:LEU:HD12	1:A:51:VAL:N	2.32	0.44
1:C:51:VAL:HG13	1:C:55:GLY:C	2.42	0.44
1:D:478:CYS:C	1:D:480:ARG:N	2.72	0.44
1:E:171:ILE:CG2	1:E:172:SER:N	2.79	0.44
1:F:385:THR:HA	1:F:388:THR:HG23	1.98	0.44
1:A:44:VAL:HG11	1:A:50:LEU:HB3	1.99	0.44
1:A:109:ASN:C	1:A:177:GLY:HA3	2.43	0.44
1:B:109:ASN:C	1:B:177:GLY:HA3	2.43	0.44
1:B:456:TYR:CE2	1:B:503:ILE:HD11	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:44:VAL:HG11	1:D:50:LEU:HB3	1.99	0.44
1:D:147:THR:O	1:D:151:ILE:HG23	2.17	0.44
1:D:151:ILE:HG13	1:D:152:ALA:N	2.30	0.44
1:D:454:TYR:O	1:D:467:TRP:CZ3	2.71	0.44
1:D:502:VAL:CG1	1:D:504:LYS:H	2.31	0.44
1:F:502:VAL:CG1	1:F:504:LYS:H	2.31	0.44
1:A:51:VAL:HG13	1:A:55:GLY:C	2.42	0.44
1:A:253:PRO:HD2	1:A:336:GLY:HA2	2.00	0.44
1:B:100:GLU:HB2	1:B:185:GLY:CA	2.48	0.44
1:C:364:ASP:N	1:C:364:ASP:OD1	2.50	0.44
1:D:153:LYS:O	1:D:157:VAL:HG22	2.17	0.44
1:D:508:GLU:HA	1:D:509:THR:HA	1.59	0.44
1:E:151:ILE:HG13	1:E:152:ALA:N	2.30	0.44
1:E:153:LYS:O	1:E:157:VAL:HG22	2.17	0.44
1:E:253:PRO:HD2	1:E:336:GLY:HA2	2.00	0.44
1:E:375:ALA:HB3	1:E:406:LEU:O	2.17	0.44
1:F:254:ILE:HB	1:F:258:GLY:O	2.18	0.44
1:F:375:ALA:HB3	1:F:406:LEU:O	2.17	0.44
1:A:228:GLY:HA2	1:A:345:SER:CB	2.32	0.43
1:A:240:SER:HB3	1:A:279:TYR:CE1	2.53	0.43
1:C:254:ILE:HB	1:C:258:GLY:O	2.18	0.43
1:C:453:ASN:HD22	1:C:453:ASN:N	2.15	0.43
1:D:407:CYS:O	1:D:451:ASP:HB3	2.14	0.43
1:E:50:LEU:HD12	1:E:51:VAL:N	2.31	0.43
1:F:51:VAL:HG13	1:F:55:GLY:C	2.42	0.43
1:F:109:ASN:C	1:F:177:GLY:HA3	2.43	0.43
1:F:178:LEU:HD23	1:F:178:LEU:N	2.22	0.43
1:F:252:LEU:HA	1:F:253:PRO:HD3	1.80	0.43
1:F:509:THR:CB	1:F:510:PRO:HA	2.44	0.43
1:A:100:GLU:HB2	1:A:185:GLY:CA	2.48	0.43
1:A:404:LEU:HD23	1:A:404:LEU:HA	1.91	0.43
1:B:147:THR:O	1:B:151:ILE:HG23	2.17	0.43
1:B:375:ALA:HB3	1:B:406:LEU:O	2.17	0.43
1:C:502:VAL:CG1	1:C:504:LYS:H	2.31	0.43
1:C:508:GLU:HA	1:C:509:THR:HA	1.59	0.43
1:E:419:VAL:HA	1:E:422:ALA:CB	2.46	0.43
1:A:375:ALA:HB3	1:A:406:LEU:O	2.17	0.43
1:B:100:GLU:HG3	1:B:186:LYS:N	2.32	0.43
1:B:254:ILE:HB	1:B:258:GLY:O	2.18	0.43
1:B:454:TYR:O	1:B:467:TRP:CZ3	2.71	0.43
1:C:109:ASN:C	1:C:177:GLY:HA3	2.43	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:48:VAL:O	1:D:48:VAL:HG12	2.19	0.43
1:D:222:VAL:HG11	1:D:236:ILE:HG12	2.00	0.43
1:E:44:VAL:HG11	1:E:50:LEU:HB3	1.99	0.43
1:A:160:TYR:HA	1:A:161:PRO:HA	1.82	0.43
1:B:295:LEU:HD11	1:B:314:PHE:CD2	2.54	0.43
1:C:147:THR:O	1:C:151:ILE:HG23	2.17	0.43
1:C:229:GLU:H	1:C:229:GLU:CD	2.26	0.43
1:D:229:GLU:H	1:D:229:GLU:CD	2.26	0.43
1:D:364:ASP:N	1:D:364:ASP:OD1	2.51	0.43
1:D:377:SER:HA	1:D:469:PRO:HG3	2.00	0.43
1:E:407:CYS:O	1:E:451:ASP:HB3	2.14	0.43
1:E:502:VAL:CG1	1:E:504:LYS:H	2.31	0.43
1:A:48:VAL:O	1:A:48:VAL:HG12	2.18	0.43
1:A:454:TYR:O	1:A:467:TRP:CZ3	2.71	0.43
1:A:508:GLU:HA	1:A:509:THR:HA	1.59	0.43
1:C:100:GLU:HB2	1:C:185:GLY:CA	2.48	0.43
1:D:62:ALA:HB1	1:D:466:ARG:HE	1.83	0.43
1:D:253:PRO:HD2	1:D:336:GLY:HA2	2.00	0.43
1:E:99:ILE:O	1:E:100:GLU:C	2.61	0.43
1:E:197:GLU:HB3	1:E:199:GLU:CD	2.44	0.43
1:E:364:ASP:OD1	1:E:364:ASP:N	2.50	0.43
1:F:194:LEU:HD23	1:F:194:LEU:C	2.43	0.43
1:F:197:GLU:HB3	1:F:199:GLU:CD	2.44	0.43
1:A:194:LEU:HD23	1:A:194:LEU:C	2.44	0.43
1:A:254:ILE:HB	1:A:258:GLY:O	2.18	0.43
1:B:44:VAL:HG11	1:B:50:LEU:HB3	1.99	0.43
1:B:48:VAL:HG12	1:B:48:VAL:O	2.19	0.43
1:C:34:GLN:OE1	1:C:230:LEU:HD11	2.19	0.43
1:C:194:LEU:HD23	1:C:194:LEU:C	2.44	0.43
1:C:197:GLU:HB3	1:C:199:GLU:CD	2.44	0.43
1:C:440:ASN:ND2	1:C:440:ASN:C	2.77	0.43
1:C:454:TYR:O	1:C:467:TRP:CZ3	2.71	0.43
1:E:97:GLY:O	1:E:98:ASN:O	2.37	0.43
1:A:478:CYS:C	1:A:480:ARG:N	2.72	0.43
1:B:99:ILE:O	1:B:100:GLU:C	2.61	0.43
1:B:151:ILE:HG13	1:B:152:ALA:N	2.30	0.43
1:B:194:LEU:HD23	1:B:194:LEU:C	2.44	0.43
1:B:197:GLU:HB3	1:B:199:GLU:CD	2.44	0.43
1:C:222:VAL:HG11	1:C:236:ILE:HG12	2.00	0.43
1:C:252:LEU:HA	1:C:253:PRO:HD3	1.80	0.43
1:C:253:PRO:HD2	1:C:336:GLY:HA2	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:295:LEU:HD11	1:C:314:PHE:CD2	2.54	0.43
1:D:171:ILE:CG2	1:D:172:SER:H	2.25	0.43
1:D:194:LEU:HD23	1:D:194:LEU:C	2.44	0.43
1:D:440:ASN:ND2	1:D:440:ASN:C	2.77	0.43
1:E:62:ALA:HB1	1:E:466:ARG:HE	1.83	0.43
1:E:240:SER:HB3	1:E:279:TYR:CE1	2.53	0.43
1:E:295:LEU:HD11	1:E:314:PHE:CD2	2.54	0.43
1:E:453:ASN:HD22	1:E:453:ASN:N	2.15	0.43
1:F:100:GLU:HB2	1:F:185:GLY:CA	2.48	0.43
1:A:62:ALA:HB1	1:A:466:ARG:HE	1.83	0.43
1:A:295:LEU:HD11	1:A:314:PHE:CD2	2.54	0.43
1:A:440:ASN:ND2	1:A:440:ASN:C	2.77	0.43
1:B:34:GLN:OE1	1:B:230:LEU:HD11	2.19	0.43
1:D:97:GLY:O	1:D:98:ASN:O	2.37	0.43
1:D:254:ILE:HB	1:D:258:GLY:O	2.18	0.43
1:D:419:VAL:HA	1:D:422:ALA:CB	2.46	0.43
1:E:109:ASN:C	1:E:177:GLY:HA3	2.43	0.43
1:E:304:ILE:HG13	1:E:305:TYR:HE2	1.81	0.43
1:E:440:ASN:ND2	1:E:440:ASN:C	2.77	0.43
1:F:44:VAL:HG11	1:F:50:LEU:HB3	1.99	0.43
1:F:48:VAL:O	1:F:48:VAL:HG12	2.19	0.43
1:F:240:SER:HB3	1:F:279:TYR:CE1	2.53	0.43
1:F:440:ASN:ND2	1:F:440:ASN:C	2.77	0.43
1:F:499:ILE:HD13	1:F:499:ILE:N	2.28	0.43
1:A:215:LYS:HE3	1:A:329:ASN:OD1	2.19	0.43
1:B:163:LEU:HB3	1:B:164:GLY:H	1.63	0.43
1:B:502:VAL:CG1	1:B:504:LYS:H	2.31	0.43
1:D:502:VAL:CG1	1:D:503:ILE:N	2.82	0.43
1:F:74:TYR:OH	1:F:473:ASP:OD2	2.35	0.43
1:F:97:GLY:O	1:F:98:ASN:O	2.37	0.43
1:F:222:VAL:HG11	1:F:236:ILE:HG12	2.00	0.43
1:A:163:LEU:HB3	1:A:164:GLY:H	1.63	0.43
1:B:312:ASP:O	1:B:316:ALA:HB2	2.19	0.43
1:C:99:ILE:O	1:C:100:GLU:C	2.61	0.43
1:C:100:GLU:HG3	1:C:186:LYS:N	2.32	0.43
1:D:197:GLU:HB3	1:D:199:GLU:CD	2.44	0.43
1:D:312:ASP:O	1:D:316:ALA:HB2	2.19	0.43
1:F:67:SER:OG	1:F:472:ALA:HB2	2.19	0.43
1:A:67:SER:OG	1:A:472:ALA:HB2	2.19	0.42
1:B:89:ALA:HB3	1:B:194:LEU:CD1	2.49	0.42
1:B:440:ASN:ND2	1:B:440:ASN:C	2.77	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:89:ALA:HB3	1:C:194:LEU:CD1	2.49	0.42
1:C:215:LYS:HE3	1:C:329:ASN:OD1	2.19	0.42
1:D:24:THR:O	1:D:26:THR:N	2.52	0.42
1:D:240:SER:HB3	1:D:279:TYR:CE1	2.53	0.42
1:E:56:GLN:HA	1:E:57:PRO:HD3	1.81	0.42
1:E:100:GLU:HG3	1:E:186:LYS:N	2.32	0.42
1:E:254:ILE:HB	1:E:258:GLY:O	2.18	0.42
1:E:377:SER:HA	1:E:469:PRO:HG3	2.00	0.42
1:F:253:PRO:HD2	1:F:336:GLY:HA2	2.00	0.42
1:A:197:GLU:HB3	1:A:199:GLU:CD	2.44	0.42
1:B:215:LYS:HE3	1:B:329:ASN:OD1	2.19	0.42
1:B:229:GLU:CD	1:B:229:GLU:H	2.26	0.42
1:C:312:ASP:O	1:C:316:ALA:HB2	2.19	0.42
1:C:377:SER:HA	1:C:469:PRO:HG3	2.00	0.42
1:C:509:THR:HG1	1:C:510:PRO:HA	1.81	0.42
1:D:295:LEU:HD11	1:D:314:PHE:CD2	2.54	0.42
1:E:312:ASP:O	1:E:316:ALA:HB2	2.19	0.42
1:F:62:ALA:HB1	1:F:466:ARG:HE	1.83	0.42
1:F:160:TYR:HA	1:F:161:PRO:HA	1.82	0.42
1:F:304:ILE:HG13	1:F:305:TYR:HE2	1.81	0.42
1:F:364:ASP:OD1	1:F:364:ASP:N	2.50	0.42
1:F:377:SER:HA	1:F:469:PRO:HG3	2.00	0.42
1:A:99:ILE:O	1:A:100:GLU:C	2.61	0.42
1:A:312:ASP:O	1:A:316:ALA:HB2	2.19	0.42
1:A:502:VAL:CG1	1:A:503:ILE:N	2.82	0.42
1:B:62:ALA:HB1	1:B:466:ARG:HE	1.83	0.42
1:C:24:THR:O	1:C:26:THR:N	2.52	0.42
1:C:240:SER:HB3	1:C:279:TYR:CE1	2.53	0.42
1:E:91:ASN:ND2	1:E:226:TYR:C	2.78	0.42
1:E:100:GLU:HB2	1:E:185:GLY:CA	2.48	0.42
1:E:161:PRO:O	1:E:162:THR:C	2.63	0.42
1:E:171:ILE:CG2	1:E:172:SER:H	2.25	0.42
1:E:194:LEU:HD23	1:E:194:LEU:C	2.43	0.42
1:E:222:VAL:HG11	1:E:236:ILE:HG12	2.00	0.42
1:A:77:ASP:OD1	1:A:77:ASP:O	2.37	0.42
1:A:100:GLU:HG3	1:A:186:LYS:N	2.32	0.42
1:A:502:VAL:CG1	1:A:504:LYS:H	2.31	0.42
1:B:24:THR:O	1:B:26:THR:N	2.52	0.42
1:B:67:SER:OG	1:B:472:ALA:HB2	2.19	0.42
1:B:364:ASP:OD1	1:B:364:ASP:N	2.50	0.42
1:B:371:GLN:O	1:B:403:CYS:HA	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:502:VAL:CG1	1:B:503:ILE:N	2.82	0.42
1:C:56:GLN:HA	1:C:57:PRO:HD3	1.81	0.42
1:C:91:ASN:ND2	1:C:226:TYR:C	2.78	0.42
1:C:161:PRO:O	1:C:162:THR:C	2.63	0.42
1:D:34:GLN:OE1	1:D:230:LEU:HD11	2.19	0.42
1:E:509:THR:CB	1:E:510:PRO:HA	2.44	0.42
1:E:509:THR:HG1	1:E:510:PRO:HA	1.83	0.42
1:F:502:VAL:CG1	1:F:503:ILE:N	2.82	0.42
1:A:24:THR:O	1:A:26:THR:N	2.52	0.42
1:C:171:ILE:CG2	1:C:172:SER:H	2.25	0.42
1:C:236:ILE:HD11	1:C:340:LEU:HD11	2.01	0.42
1:C:502:VAL:CG1	1:C:503:ILE:N	2.82	0.42
1:F:24:THR:O	1:F:26:THR:N	2.52	0.42
1:F:34:GLN:OE1	1:F:230:LEU:HD11	2.19	0.42
1:F:90:LYS:N	1:F:344:LEU:O	2.48	0.42
1:F:171:ILE:CG2	1:F:172:SER:H	2.25	0.42
1:F:295:LEU:HD11	1:F:314:PHE:CD2	2.54	0.42
1:B:240:SER:HB3	1:B:279:TYR:CE1	2.54	0.42
1:C:97:GLY:O	1:C:98:ASN:O	2.37	0.42
1:C:370:VAL:HG23	1:C:370:VAL:O	2.20	0.42
1:C:407:CYS:O	1:C:451:ASP:HB3	2.14	0.42
1:D:91:ASN:ND2	1:D:226:TYR:C	2.78	0.42
1:D:213:ASN:N	1:D:213:ASN:ND2	2.68	0.42
1:F:236:ILE:HD11	1:F:340:LEU:HD11	2.01	0.42
1:F:407:CYS:O	1:F:451:ASP:HB3	2.14	0.42
1:A:34:GLN:OE1	1:A:230:LEU:HD11	2.19	0.42
1:A:161:PRO:O	1:A:162:THR:C	2.63	0.42
1:B:304:ILE:HG13	1:B:305:TYR:HE2	1.82	0.42
1:C:62:ALA:HB1	1:C:466:ARG:HE	1.83	0.42
1:D:109:ASN:C	1:D:177:GLY:HA3	2.43	0.42
1:E:34:GLN:OE1	1:E:230:LEU:HD11	2.19	0.42
1:E:48:VAL:HG12	1:E:48:VAL:O	2.18	0.42
1:E:89:ALA:HB3	1:E:194:LEU:CD1	2.49	0.42
1:F:161:PRO:O	1:F:162:THR:C	2.63	0.42
1:F:171:ILE:CG2	1:F:172:SER:N	2.79	0.42
1:A:89:ALA:HB3	1:A:194:LEU:CD1	2.49	0.42
1:A:91:ASN:ND2	1:A:226:TYR:C	2.78	0.42
1:A:97:GLY:O	1:A:98:ASN:O	2.37	0.42
1:A:304:ILE:HG13	1:A:305:TYR:HE2	1.81	0.42
1:A:371:GLN:O	1:A:403:CYS:HA	2.20	0.42
1:B:97:GLY:O	1:B:98:ASN:O	2.37	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:453:ASN:HD22	1:B:453:ASN:N	2.16	0.42
1:C:90:LYS:N	1:C:344:LEU:O	2.48	0.42
1:D:100:GLU:HB2	1:D:185:GLY:CA	2.48	0.42
1:D:450:ILE:CG1	1:D:451:ASP:N	2.53	0.42
1:E:67:SER:OG	1:E:472:ALA:HB2	2.19	0.42
1:F:215:LYS:HE3	1:F:329:ASN:OD1	2.19	0.42
1:A:377:SER:HA	1:A:469:PRO:HG3	2.00	0.42
1:B:77:ASP:OD1	1:B:77:ASP:O	2.37	0.42
1:C:48:VAL:O	1:C:48:VAL:HG12	2.19	0.42
1:C:371:GLN:O	1:C:403:CYS:HA	2.20	0.42
1:D:89:ALA:HB3	1:D:194:LEU:CD1	2.49	0.42
1:D:215:LYS:HE3	1:D:329:ASN:OD1	2.19	0.42
1:F:77:ASP:OD1	1:F:77:ASP:O	2.37	0.42
1:F:371:GLN:O	1:F:403:CYS:HA	2.20	0.42
1:A:370:VAL:HG23	1:A:370:VAL:O	2.20	0.42
1:A:375:ALA:C	1:A:377:SER:H	2.28	0.42
1:B:252:LEU:HA	1:B:253:PRO:HD3	1.80	0.42
1:B:409:PRO:HG2	1:B:454:TYR:CE1	2.55	0.42
1:C:409:PRO:HG2	1:C:454:TYR:CE1	2.55	0.42
1:C:478:CYS:C	1:C:480:ARG:N	2.72	0.42
1:C:509:THR:N	1:C:510:PRO:CA	2.73	0.42
1:D:370:VAL:HG23	1:D:370:VAL:O	2.20	0.42
1:D:404:LEU:HD23	1:D:404:LEU:HA	1.91	0.42
1:E:236:ILE:HD11	1:E:340:LEU:HD11	2.01	0.42
1:F:312:ASP:O	1:F:316:ALA:HB2	2.19	0.42
1:F:370:VAL:HG23	1:F:370:VAL:O	2.20	0.42
1:A:222:VAL:HG11	1:A:236:ILE:HG12	2.00	0.41
1:B:222:VAL:HG11	1:B:236:ILE:HG12	2.00	0.41
1:C:46:ASN:HB2	1:C:49:ASP:CB	2.50	0.41
1:C:404:LEU:HD23	1:C:404:LEU:HA	1.90	0.41
1:D:409:PRO:HG2	1:D:454:TYR:CE1	2.55	0.41
1:E:215:LYS:HE3	1:E:329:ASN:OD1	2.19	0.41
1:C:375:ALA:C	1:C:377:SER:H	2.28	0.41
1:C:406:LEU:HA	1:C:449:ALA:O	2.20	0.41
1:D:236:ILE:HD11	1:D:340:LEU:HD11	2.01	0.41
1:E:24:THR:O	1:E:26:THR:N	2.52	0.41
1:E:370:VAL:HG23	1:E:370:VAL:O	2.20	0.41
1:E:371:GLN:O	1:E:403:CYS:HA	2.20	0.41
1:E:502:VAL:CG1	1:E:503:ILE:N	2.82	0.41
1:A:391:LYS:CE	1:A:440:ASN:ND2	2.84	0.41
1:A:409:PRO:HG2	1:A:454:TYR:CE1	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:207:ALA:O	1:D:208:VAL:C	2.64	0.41
1:D:406:LEU:HA	1:D:449:ALA:O	2.20	0.41
1:E:207:ALA:O	1:E:208:VAL:C	2.64	0.41
1:E:391:LYS:CE	1:E:440:ASN:ND2	2.84	0.41
1:E:409:PRO:HG2	1:E:454:TYR:CE1	2.55	0.41
1:F:89:ALA:HB3	1:F:194:LEU:CD1	2.49	0.41
1:F:119:LYS:HE3	1:F:124:ASP:OD1	2.20	0.41
1:F:391:LYS:CE	1:F:440:ASN:ND2	2.84	0.41
1:A:364:ASP:OD1	1:A:364:ASP:N	2.51	0.41
1:B:91:ASN:ND2	1:B:226:TYR:C	2.78	0.41
1:B:101:TYR:CD2	1:B:101:TYR:C	2.97	0.41
1:C:101:TYR:CD2	1:C:101:TYR:C	2.98	0.41
1:D:46:ASN:HB2	1:D:49:ASP:CB	2.50	0.41
1:D:100:GLU:HG3	1:D:186:LYS:N	2.32	0.41
1:D:390:GLN:O	1:D:391:LYS:C	2.63	0.41
1:E:406:LEU:HA	1:E:449:ALA:O	2.20	0.41
1:F:91:ASN:ND2	1:F:226:TYR:C	2.78	0.41
1:F:101:TYR:CD2	1:F:101:TYR:C	2.98	0.41
1:F:205:MET:HE3	1:F:205:MET:HB3	1.97	0.41
1:F:408:SER:OG	1:F:409:PRO:HD2	2.21	0.41
1:A:55:GLY:C	1:A:65:PHE:HE1	2.29	0.41
1:A:205:MET:HE3	1:A:205:MET:HB3	1.97	0.41
1:A:390:GLN:O	1:A:391:LYS:C	2.63	0.41
1:A:406:LEU:HA	1:A:449:ALA:O	2.20	0.41
1:B:377:SER:HA	1:B:469:PRO:HG3	2.00	0.41
1:C:50:LEU:HD11	1:C:65:PHE:CE1	2.56	0.41
1:C:408:SER:OG	1:C:409:PRO:HD2	2.21	0.41
1:D:67:SER:OG	1:D:472:ALA:HB2	2.19	0.41
1:E:213:ASN:N	1:E:213:ASN:ND2	2.68	0.41
1:E:390:GLN:O	1:E:391:LYS:C	2.63	0.41
1:F:213:ASN:N	1:F:213:ASN:ND2	2.68	0.41
1:F:406:LEU:HA	1:F:449:ALA:O	2.20	0.41
1:F:409:PRO:HG2	1:F:454:TYR:CE1	2.55	0.41
1:A:50:LEU:HD11	1:A:65:PHE:CE1	2.56	0.41
1:B:23:SER:OG	1:B:483:ASN:CA	2.69	0.41
1:B:50:LEU:HD11	1:B:65:PHE:CE1	2.56	0.41
1:B:478:CYS:C	1:B:480:ARG:N	2.72	0.41
1:D:161:PRO:O	1:D:162:THR:C	2.63	0.41
1:E:119:LYS:HE3	1:E:124:ASP:OD1	2.20	0.41
1:E:404:LEU:HD23	1:E:404:LEU:HA	1.91	0.41
1:F:404:LEU:HD23	1:F:404:LEU:HA	1.90	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:119:LYS:HE3	1:A:124:ASP:OD1	2.20	0.41
1:A:375:ALA:C	1:A:377:SER:N	2.79	0.41
1:A:460:LYS:HB3	1:A:460:LYS:HE3	1.82	0.41
1:B:370:VAL:O	1:B:370:VAL:HG23	2.20	0.41
1:B:390:GLN:O	1:B:391:LYS:C	2.63	0.41
1:C:77:ASP:OD1	1:C:77:ASP:O	2.38	0.41
1:C:390:GLN:O	1:C:391:LYS:C	2.63	0.41
1:D:23:SER:HG	1:D:483:ASN:HB3	1.78	0.41
1:D:101:TYR:CD2	1:D:101:TYR:C	2.98	0.41
1:D:408:SER:OG	1:D:409:PRO:HD2	2.21	0.41
1:E:77:ASP:OD1	1:E:77:ASP:O	2.37	0.41
1:F:508:GLU:HA	1:F:509:THR:HA	1.59	0.41
1:A:236:ILE:HD11	1:A:340:LEU:HD11	2.01	0.41
1:B:74:TYR:OH	1:B:473:ASP:OD2	2.35	0.41
1:B:341:SER:C	1:B:343:GLY:H	2.29	0.41
1:B:375:ALA:C	1:B:377:SER:H	2.28	0.41
1:B:408:SER:OG	1:B:409:PRO:HD2	2.21	0.41
1:B:459:ASP:C	1:B:459:ASP:OD2	2.64	0.41
1:C:119:LYS:HE3	1:C:124:ASP:OD1	2.20	0.41
1:D:100:GLU:HB2	1:D:185:GLY:HA3	2.03	0.41
1:D:252:LEU:HA	1:D:253:PRO:HD3	1.80	0.41
1:D:371:GLN:O	1:D:403:CYS:HA	2.20	0.41
1:D:391:LYS:CE	1:D:440:ASN:ND2	2.84	0.41
1:F:50:LEU:HD11	1:F:65:PHE:CE1	2.56	0.41
1:A:194:LEU:C	1:A:194:LEU:CD2	2.94	0.41
1:A:290:VAL:HG11	1:A:322:TYR:CE1	2.56	0.41
1:A:308:ASN:HD21	1:A:313:ASP:CB	2.34	0.41
1:A:482:ASP:HB3	1:A:483:ASN:H	1.78	0.41
1:B:90:LYS:N	1:B:344:LEU:O	2.48	0.41
1:B:150:ILE:H	1:B:150:ILE:HG12	1.63	0.41
1:B:161:PRO:O	1:B:162:THR:C	2.63	0.41
1:B:236:ILE:HD11	1:B:340:LEU:HD11	2.01	0.41
1:B:308:ASN:HD21	1:B:313:ASP:CB	2.34	0.41
1:B:391:LYS:CE	1:B:440:ASN:ND2	2.84	0.41
1:C:67:SER:OG	1:C:472:ALA:HB2	2.19	0.41
1:C:194:LEU:C	1:C:194:LEU:CD2	2.94	0.41
1:C:383:LEU:O	1:C:384:GLU:C	2.64	0.41
1:D:73:GLN:HB3	1:D:500:LEU:CD1	2.51	0.41
1:D:194:LEU:C	1:D:194:LEU:CD2	2.94	0.41
1:D:375:ALA:C	1:D:377:SER:H	2.28	0.41
1:D:383:LEU:O	1:D:384:GLU:C	2.64	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:46:ASN:HB2	1:E:49:ASP:CB	2.50	0.41
1:E:101:TYR:CD2	1:E:101:TYR:C	2.98	0.41
1:F:100:GLU:HB2	1:F:185:GLY:HA3	2.03	0.41
1:F:375:ALA:C	1:F:377:SER:H	2.28	0.41
1:F:375:ALA:C	1:F:377:SER:N	2.79	0.41
1:F:459:ASP:C	1:F:459:ASP:OD2	2.64	0.41
1:A:73:GLN:HB3	1:A:500:LEU:CD1	2.51	0.41
1:A:90:LYS:N	1:A:344:LEU:O	2.48	0.41
1:A:341:SER:C	1:A:343:GLY:H	2.29	0.41
1:B:406:LEU:HA	1:B:449:ALA:O	2.20	0.41
1:D:50:LEU:HD11	1:D:65:PHE:CE1	2.56	0.41
1:D:55:GLY:C	1:D:65:PHE:HE1	2.28	0.41
1:E:245:ALA:C	1:E:247:GLY:N	2.79	0.41
1:E:383:LEU:O	1:E:384:GLU:C	2.64	0.41
1:E:408:SER:OG	1:E:409:PRO:HD2	2.21	0.41
1:F:308:ASN:HD21	1:F:313:ASP:CB	2.34	0.41
1:A:92:SER:OG	1:A:343:GLY:HA3	2.21	0.40
1:A:101:TYR:CD2	1:A:101:TYR:C	2.98	0.40
1:A:459:ASP:OD2	1:A:459:ASP:C	2.64	0.40
1:D:92:SER:OG	1:D:343:GLY:HA3	2.21	0.40
1:D:459:ASP:OD2	1:D:459:ASP:C	2.64	0.40
1:E:92:SER:OG	1:E:343:GLY:HA3	2.22	0.40
1:E:100:GLU:HB2	1:E:185:GLY:HA3	2.03	0.40
1:E:194:LEU:C	1:E:194:LEU:CD2	2.94	0.40
1:E:303:ASP:C	1:E:305:TYR:H	2.29	0.40
1:E:459:ASP:OD2	1:E:459:ASP:C	2.64	0.40
1:F:73:GLN:HB3	1:F:500:LEU:CD1	2.51	0.40
1:F:290:VAL:HG11	1:F:322:TYR:CE1	2.56	0.40
1:F:410:PRO:O	1:F:411:ARG:C	2.65	0.40
1:A:23:SER:OG	1:A:483:ASN:CA	2.69	0.40
1:A:408:SER:OG	1:A:409:PRO:HD2	2.21	0.40
1:B:290:VAL:HG11	1:B:322:TYR:CE1	2.56	0.40
1:B:383:LEU:O	1:B:384:GLU:C	2.64	0.40
1:C:163:LEU:HB3	1:C:164:GLY:H	1.63	0.40
1:C:207:ALA:O	1:C:208:VAL:C	2.64	0.40
1:C:391:LYS:CE	1:C:440:ASN:ND2	2.84	0.40
1:D:308:ASN:HD21	1:D:313:ASP:CB	2.34	0.40
1:E:73:GLN:HB3	1:E:500:LEU:CD1	2.51	0.40
1:F:100:GLU:HG3	1:F:186:LYS:N	2.32	0.40
1:A:303:ASP:C	1:A:305:TYR:H	2.29	0.40
1:A:317:LYS:HE3	1:A:317:LYS:HB2	1.97	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:56:GLN:HA	1:B:57:PRO:HD3	1.81	0.40
1:B:119:LYS:HE3	1:B:124:ASP:OD1	2.20	0.40
1:B:194:LEU:C	1:B:194:LEU:CD2	2.94	0.40
1:B:249:SER:O	1:B:250:ALA:C	2.65	0.40
1:C:92:SER:OG	1:C:343:GLY:HA3	2.22	0.40
1:C:375:ALA:C	1:C:377:SER:N	2.79	0.40
1:C:459:ASP:OD2	1:C:459:ASP:C	2.64	0.40
1:D:21:ASN:O	1:D:21:ASN:CG	2.65	0.40
1:D:77:ASP:OD1	1:D:77:ASP:O	2.37	0.40
1:D:499:ILE:H	1:D:499:ILE:CD1	2.24	0.40
1:E:341:SER:C	1:E:343:GLY:H	2.28	0.40
1:F:92:SER:OG	1:F:343:GLY:HA3	2.21	0.40
1:F:207:ALA:O	1:F:208:VAL:C	2.64	0.40
1:B:21:ASN:O	1:B:21:ASN:CG	2.65	0.40
1:C:74:TYR:OH	1:C:473:ASP:OD2	2.35	0.40
1:C:100:GLU:HB2	1:C:185:GLY:HA3	2.03	0.40
1:D:43:GLN:NE2	1:D:77:ASP:HB2	2.36	0.40
1:D:341:SER:C	1:D:343:GLY:H	2.28	0.40
1:E:308:ASN:HD21	1:E:313:ASP:CB	2.34	0.40
1:E:317:LYS:HE3	1:E:317:LYS:HB2	1.97	0.40
1:E:410:PRO:O	1:E:411:ARG:C	2.65	0.40
1:B:22:ASN:O	1:B:23:SER:CB	2.70	0.40
1:B:347:ASN:HA	1:B:350:VAL:HG23	2.04	0.40
1:C:43:GLN:NE2	1:C:77:ASP:HB2	2.36	0.40
1:C:73:GLN:HB3	1:C:500:LEU:CD1	2.51	0.40
1:D:119:LYS:HE3	1:D:124:ASP:OD1	2.20	0.40
1:D:375:ALA:C	1:D:377:SER:N	2.79	0.40
1:D:427:VAL:O	1:D:431:THR:CG2	2.68	0.40
1:E:43:GLN:NE2	1:E:77:ASP:HB2	2.37	0.40
1:F:104:SER:HB2	1:F:181:VAL:HG12	2.04	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	A	475/510 (93%)	371 (78%)	87 (18%)	17 (4%)	2	20
1	B	475/510 (93%)	371 (78%)	87 (18%)	17 (4%)	2	20
1	C	475/510 (93%)	371 (78%)	87 (18%)	17 (4%)	2	20
1	D	475/510 (93%)	371 (78%)	87 (18%)	17 (4%)	2	20
1	E	475/510 (93%)	371 (78%)	87 (18%)	17 (4%)	2	20
1	F	475/510 (93%)	371 (78%)	87 (18%)	17 (4%)	2	20
All	All	2850/3060 (93%)	2226 (78%)	522 (18%)	102 (4%)	4	20

All (102) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	98	ASN
1	A	201	ALA
1	A	432	ALA
1	B	98	ASN
1	B	201	ALA
1	B	432	ALA
1	C	98	ASN
1	C	201	ALA
1	C	432	ALA
1	D	98	ASN
1	D	201	ALA
1	D	432	ALA
1	E	98	ASN
1	E	201	ALA
1	E	432	ALA
1	F	98	ASN
1	F	201	ALA
1	F	432	ALA
1	A	403	CYS
1	A	480	ARG
1	A	504	LYS
1	B	403	CYS
1	B	480	ARG
1	B	504	LYS
1	C	403	CYS
1	C	480	ARG
1	C	504	LYS

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Mol	Chain	Res	Type
1	D	403	CYS
1	D	480	ARG
1	D	504	LYS
1	E	403	CYS
1	E	480	ARG
1	E	504	LYS
1	F	403	CYS
1	F	480	ARG
1	F	504	LYS
1	A	62	ALA
1	A	311	ILE
1	A	348	ALA
1	A	482	ASP
1	B	62	ALA
1	B	311	ILE
1	B	348	ALA
1	B	482	ASP
1	C	62	ALA
1	C	311	ILE
1	C	348	ALA
1	C	482	ASP
1	D	62	ALA
1	D	311	ILE
1	D	348	ALA
1	D	482	ASP
1	E	62	ALA
1	E	311	ILE
1	E	348	ALA
1	E	482	ASP
1	F	62	ALA
1	F	311	ILE
1	F	348	ALA
1	F	482	ASP
1	A	509	THR
1	B	509	THR
1	C	509	THR
1	D	509	THR
1	E	509	THR
1	F	509	THR
1	A	25	GLY
1	A	304	ILE
1	B	25	GLY

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Mol	Chain	Res	Type
1	B	304	ILE
1	C	25	GLY
1	C	304	ILE
1	D	25	GLY
1	D	304	ILE
1	E	25	GLY
1	E	304	ILE
1	F	25	GLY
1	F	304	ILE
1	A	247	GLY
1	A	434	GLY
1	B	247	GLY
1	B	434	GLY
1	C	247	GLY
1	C	434	GLY
1	D	247	GLY
1	D	434	GLY
1	E	247	GLY
1	E	434	GLY
1	F	247	GLY
1	F	434	GLY
1	A	121	VAL
1	A	318	GLY
1	B	121	VAL
1	B	318	GLY
1	C	121	VAL
1	C	318	GLY
1	D	121	VAL
1	D	318	GLY
1	E	121	VAL
1	E	318	GLY
1	F	121	VAL
1	F	318	GLY

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	383/411 (93%)	335 (88%)	48 (12%)	4	16
1	B	383/411 (93%)	335 (88%)	48 (12%)	4	16
1	C	383/411 (93%)	335 (88%)	48 (12%)	4	16
1	D	383/411 (93%)	335 (88%)	48 (12%)	4	16
1	E	383/411 (93%)	335 (88%)	48 (12%)	4	16
1	F	383/411 (93%)	334 (87%)	49 (13%)	4	15
All	All	2298/2466 (93%)	2009 (87%)	289 (13%)	6	16

All (289) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	41	ILE
1	A	50	LEU
1	A	58	THR
1	A	66	MET
1	A	67	SER
1	A	71	PHE
1	A	74	TYR
1	A	80	VAL
1	A	84	VAL
1	A	86	ARG
1	A	95	ILE
1	A	102	THR
1	A	150	ILE
1	A	162	THR
1	A	188	ILE
1	A	189	THR
1	A	202	GLU
1	A	213	ASN
1	A	290	VAL
1	A	304	ILE
1	A	347	ASN
1	A	349	GLU
1	A	367	SER
1	A	368	VAL
1	A	371	GLN
1	A	382	SER
1	A	383	LEU
1	A	388	THR
1	A	395	SER
1	A	404	LEU

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Mol	Chain	Res	Type
1	A	406	LEU
1	A	414	VAL
1	A	417	ILE
1	A	426	LEU
1	A	430	ARG
1	A	431	THR
1	A	436	TYR
1	A	438	ASP
1	A	440	ASN
1	A	446	THR
1	A	450	ILE
1	A	453	ASN
1	A	460	LYS
1	A	464	VAL
1	A	483	ASN
1	A	499	ILE
1	A	501	ASN
1	A	507	ILE
1	B	41	ILE
1	B	50	LEU
1	B	58	THR
1	B	66	MET
1	B	67	SER
1	B	71	PHE
1	B	74	TYR
1	B	80	VAL
1	B	84	VAL
1	B	86	ARG
1	B	95	ILE
1	B	102	THR
1	B	150	ILE
1	B	162	THR
1	B	188	ILE
1	B	189	THR
1	B	202	GLU
1	B	213	ASN
1	B	290	VAL
1	B	304	ILE
1	B	347	ASN
1	B	349	GLU
1	B	367	SER
1	B	368	VAL

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Mol	Chain	Res	Type
1	B	371	GLN
1	B	382	SER
1	B	383	LEU
1	B	388	THR
1	B	395	SER
1	B	404	LEU
1	B	406	LEU
1	B	414	VAL
1	B	417	ILE
1	B	426	LEU
1	B	430	ARG
1	B	431	THR
1	B	436	TYR
1	B	438	ASP
1	B	440	ASN
1	B	446	THR
1	B	450	ILE
1	B	453	ASN
1	B	460	LYS
1	B	464	VAL
1	B	483	ASN
1	B	499	ILE
1	B	501	ASN
1	B	507	ILE
1	C	41	ILE
1	C	50	LEU
1	C	58	THR
1	C	66	MET
1	C	67	SER
1	C	71	PHE
1	C	74	TYR
1	C	80	VAL
1	C	84	VAL
1	C	86	ARG
1	C	95	ILE
1	C	102	THR
1	C	150	ILE
1	C	162	THR
1	C	188	ILE
1	C	189	THR
1	C	202	GLU
1	C	213	ASN

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Mol	Chain	Res	Type
1	C	290	VAL
1	C	304	ILE
1	C	347	ASN
1	C	349	GLU
1	C	367	SER
1	C	368	VAL
1	C	371	GLN
1	C	382	SER
1	C	383	LEU
1	C	388	THR
1	C	395	SER
1	C	404	LEU
1	C	406	LEU
1	C	414	VAL
1	C	417	ILE
1	C	426	LEU
1	C	430	ARG
1	C	431	THR
1	C	436	TYR
1	C	438	ASP
1	C	440	ASN
1	C	446	THR
1	C	450	ILE
1	C	453	ASN
1	C	460	LYS
1	C	464	VAL
1	C	483	ASN
1	C	499	ILE
1	C	501	ASN
1	C	507	ILE
1	D	41	ILE
1	D	50	LEU
1	D	58	THR
1	D	66	MET
1	D	67	SER
1	D	71	PHE
1	D	74	TYR
1	D	80	VAL
1	D	84	VAL
1	D	86	ARG
1	D	95	ILE
1	D	102	THR

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Mol	Chain	Res	Type
1	D	150	ILE
1	D	162	THR
1	D	188	ILE
1	D	189	THR
1	D	202	GLU
1	D	213	ASN
1	D	290	VAL
1	D	304	ILE
1	D	347	ASN
1	D	349	GLU
1	D	367	SER
1	D	368	VAL
1	D	371	GLN
1	D	382	SER
1	D	383	LEU
1	D	388	THR
1	D	395	SER
1	D	404	LEU
1	D	406	LEU
1	D	414	VAL
1	D	417	ILE
1	D	426	LEU
1	D	430	ARG
1	D	431	THR
1	D	436	TYR
1	D	438	ASP
1	D	440	ASN
1	D	446	THR
1	D	450	ILE
1	D	453	ASN
1	D	460	LYS
1	D	464	VAL
1	D	483	ASN
1	D	499	ILE
1	D	501	ASN
1	D	507	ILE
1	E	41	ILE
1	E	50	LEU
1	E	58	THR
1	E	66	MET
1	E	67	SER
1	E	71	PHE

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Mol	Chain	Res	Type
1	E	74	TYR
1	E	80	VAL
1	E	84	VAL
1	E	86	ARG
1	E	95	ILE
1	E	102	THR
1	E	150	ILE
1	E	162	THR
1	E	188	ILE
1	E	189	THR
1	E	202	GLU
1	E	213	ASN
1	E	290	VAL
1	E	304	ILE
1	E	347	ASN
1	E	349	GLU
1	E	367	SER
1	E	368	VAL
1	E	371	GLN
1	E	382	SER
1	E	383	LEU
1	E	388	THR
1	E	395	SER
1	E	404	LEU
1	E	406	LEU
1	E	414	VAL
1	E	417	ILE
1	E	426	LEU
1	E	430	ARG
1	E	431	THR
1	E	436	TYR
1	E	438	ASP
1	E	440	ASN
1	E	446	THR
1	E	450	ILE
1	E	453	ASN
1	E	460	LYS
1	E	464	VAL
1	E	483	ASN
1	E	499	ILE
1	E	501	ASN
1	E	507	ILE

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Mol	Chain	Res	Type
1	F	41	ILE
1	F	50	LEU
1	F	58	THR
1	F	66	MET
1	F	67	SER
1	F	71	PHE
1	F	74	TYR
1	F	80	VAL
1	F	84	VAL
1	F	86	ARG
1	F	95	ILE
1	F	102	THR
1	F	150	ILE
1	F	162	THR
1	F	188	ILE
1	F	189	THR
1	F	202	GLU
1	F	213	ASN
1	F	290	VAL
1	F	304	ILE
1	F	347	ASN
1	F	349	GLU
1	F	367	SER
1	F	368	VAL
1	F	371	GLN
1	F	382	SER
1	F	383	LEU
1	F	388	THR
1	F	395	SER
1	F	404	LEU
1	F	406	LEU
1	F	414	VAL
1	F	417	ILE
1	F	426	LEU
1	F	430	ARG
1	F	431	THR
1	F	436	TYR
1	F	437	THR
1	F	438	ASP
1	F	440	ASN
1	F	446	THR
1	F	450	ILE

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Mol	Chain	Res	Type
1	F	453	ASN
1	F	460	LYS
1	F	464	VAL
1	F	483	ASN
1	F	499	ILE
1	F	501	ASN
1	F	507	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (57) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	43	GLN
1	A	70	ASN
1	A	166	ASN
1	A	213	ASN
1	A	278	GLN
1	A	390	GLN
1	A	440	ASN
1	A	453	ASN
1	A	457	GLN
1	B	43	GLN
1	B	70	ASN
1	B	166	ASN
1	B	213	ASN
1	B	278	GLN
1	B	347	ASN
1	B	390	GLN
1	B	440	ASN
1	B	453	ASN
1	B	457	GLN
1	C	43	GLN
1	C	70	ASN
1	C	166	ASN
1	C	213	ASN
1	C	278	GLN
1	C	347	ASN
1	C	390	GLN
1	C	440	ASN
1	C	453	ASN
1	C	457	GLN
1	D	43	GLN
1	D	70	ASN

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Mol	Chain	Res	Type
1	D	166	ASN
1	D	213	ASN
1	D	278	GLN
1	D	390	GLN
1	D	440	ASN
1	D	453	ASN
1	D	457	GLN
1	E	43	GLN
1	E	70	ASN
1	E	166	ASN
1	E	213	ASN
1	E	278	GLN
1	E	390	GLN
1	E	440	ASN
1	E	453	ASN
1	E	457	GLN
1	F	43	GLN
1	F	70	ASN
1	F	166	ASN
1	F	213	ASN
1	F	278	GLN
1	F	329	ASN
1	F	390	GLN
1	F	440	ASN
1	F	453	ASN
1	F	457	GLN

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 [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

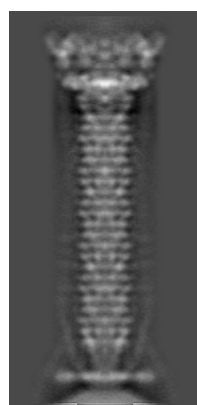
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-1126. These allow visual inspection of the internal detail of the map and identification of artifacts.

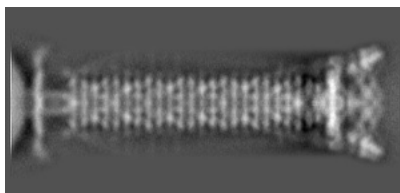
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

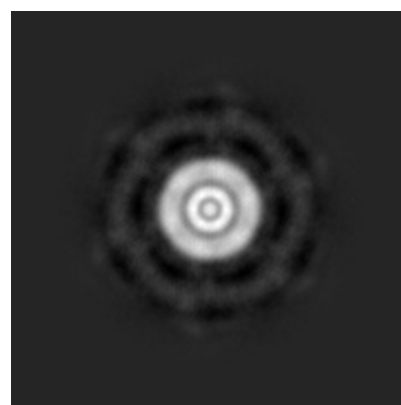
6.1.1 Primary map



X



Y

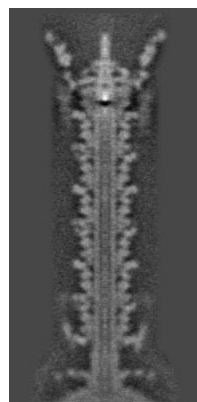


Z

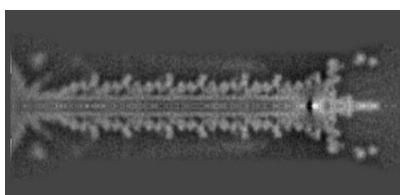
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

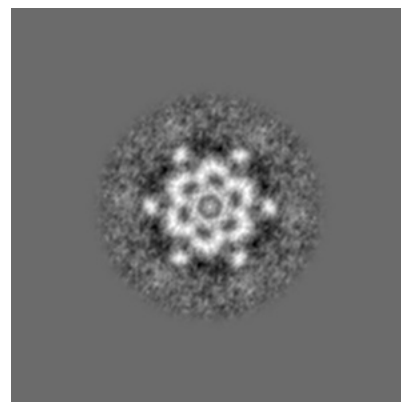
6.2.1 Primary map



X Index: 90



Y Index: 90

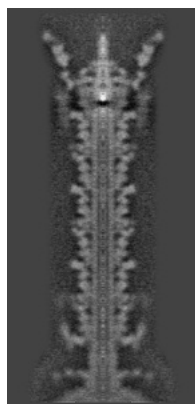


Z Index: 190

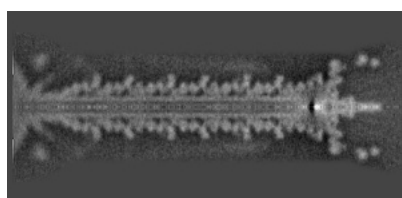
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

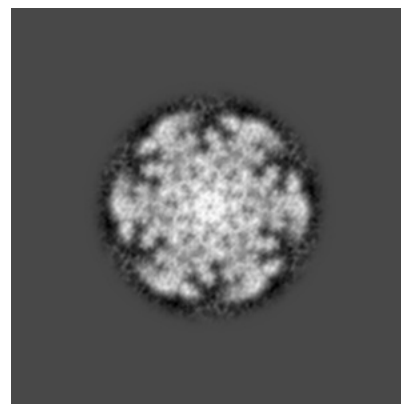
6.3.1 Primary map



X Index: 89



Y Index: 90

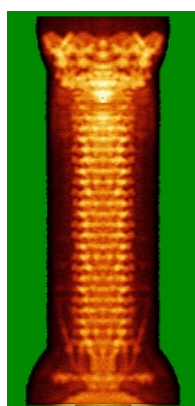


Z Index: 310

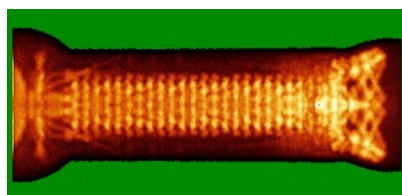
The images above show the largest variance slices of the map in three orthogonal directions.

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

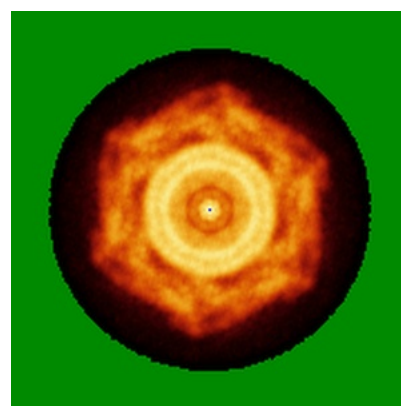
6.4.1 Primary map



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

This section was not generated.

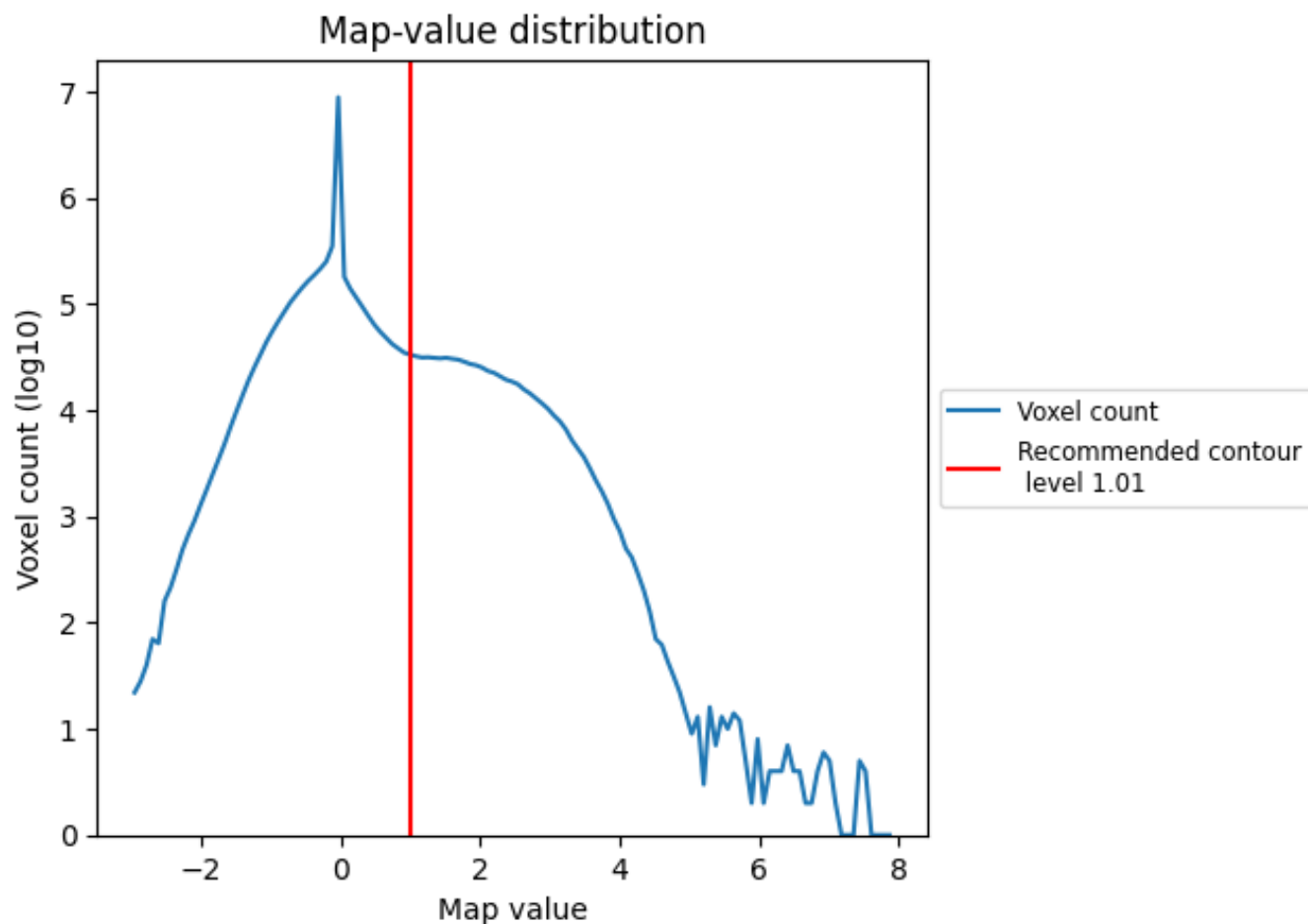
6.6 Mask visualisation

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

7 Map analysis [i](#)

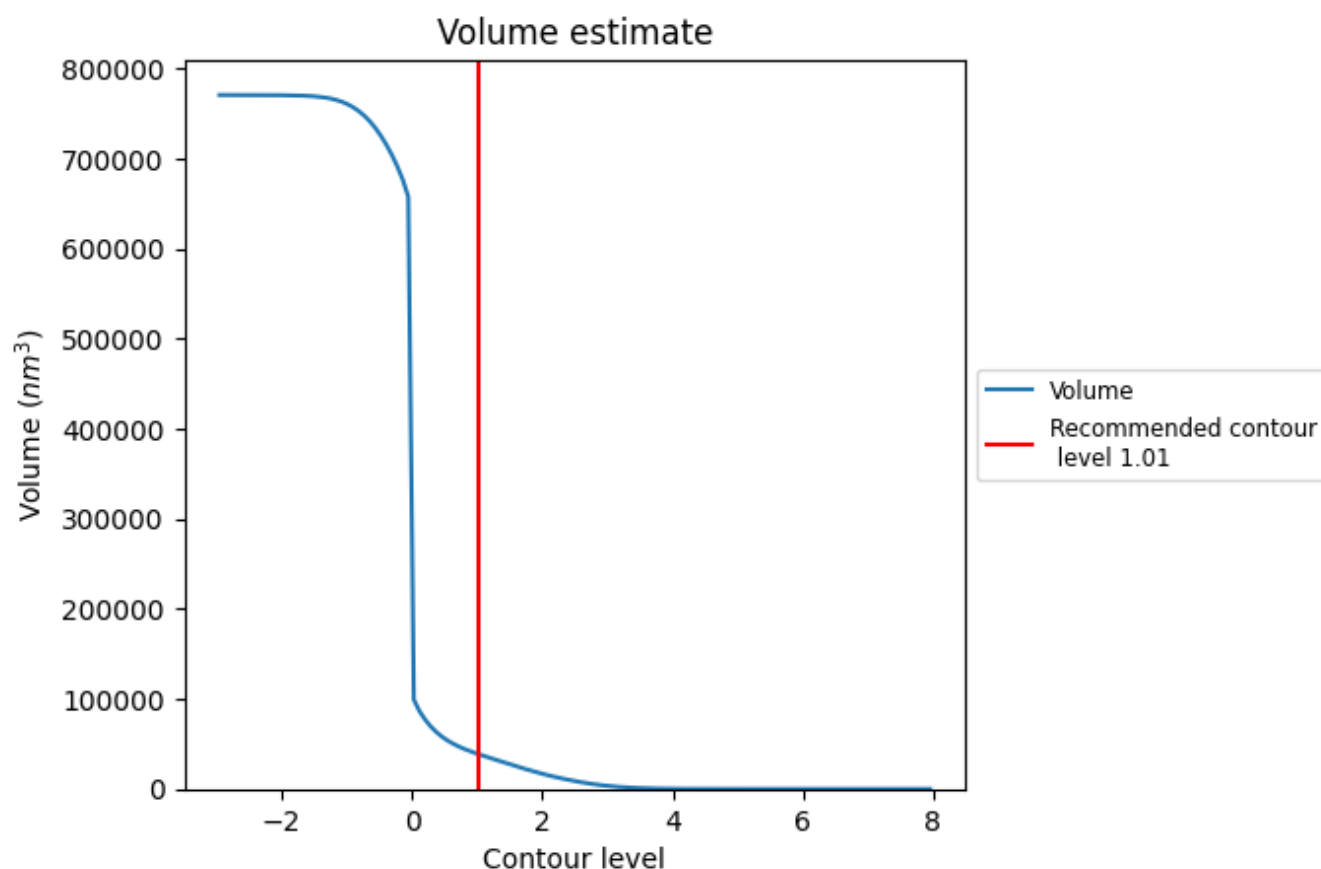
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



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

7.2 Volume estimate [i](#)



The volume at the recommended contour level is 38892 nm³; this corresponds to an approximate mass of 35132 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](#)

This section was not generated. The rotationally averaged power spectrum is only generated for cubic maps.

8 Fourier-Shell correlation

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

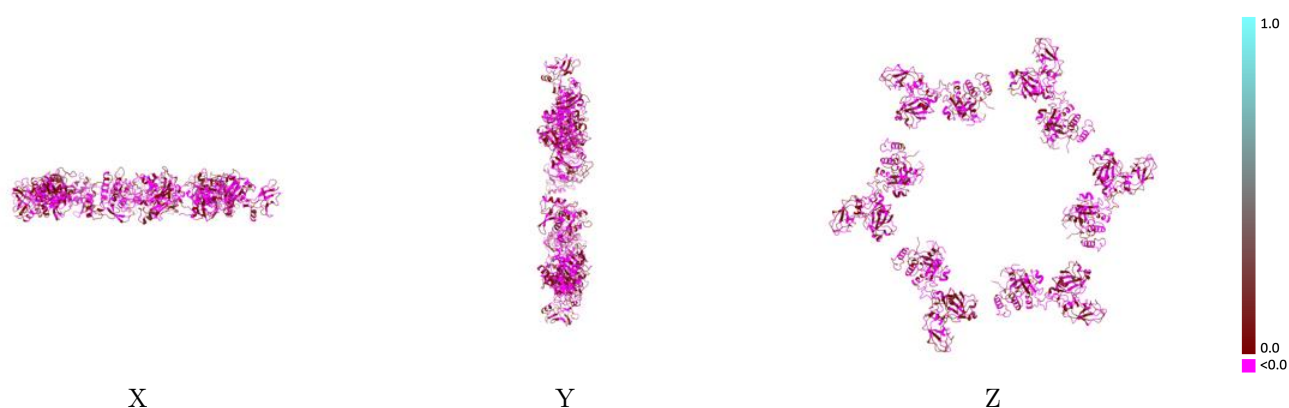
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-1126 and PDB model 3FOH. Per-residue inclusion information can be found in section 3 on page 6.

9.1 Map-model overlay [i](#)

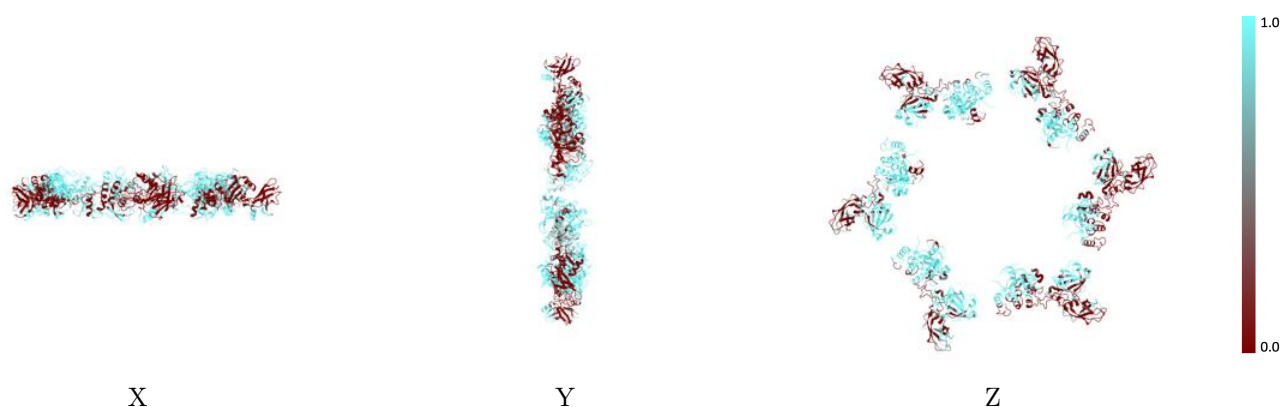
This section was not generated.

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



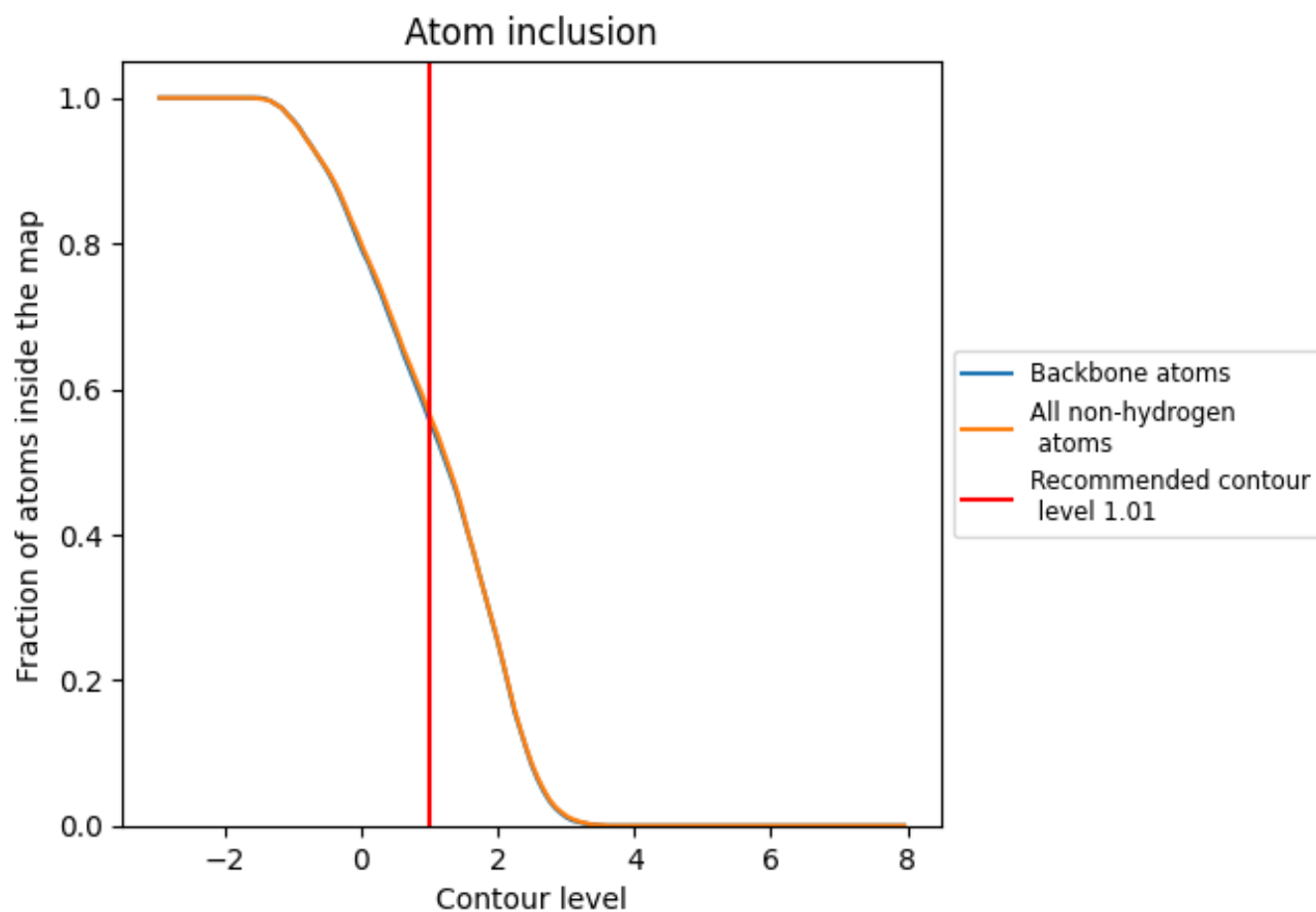
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

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



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (1.01).

9.4 Atom inclusion [i](#)



At the recommended contour level, 55% of all backbone atoms, 56% 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 (1.01) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.5620	0.0010
A	0.4620	-0.0020
B	0.6060	0.0050
C	0.6970	0.0100
D	0.6950	0.0020
E	0.5150	0.0010
F	0.3990	-0.0080

1.0
0.0
-0.0