



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 11:43 AM UTC

PDB ID : 5ENY / pdb_00005eny
Title : Ketosynthase from module 6 connected to acyl carrier protein from module 5 (unobservable) of the bacillaene synthase from *Bacillus subtilis* 168
Authors : Wagner, D.T.; Gay, D.C.; Keatinge-Clay, A.T.
Deposited on : 2015-11-09
Resolution : 4.00 Å(reported)

This is a Full wwPDB X-ray Structure 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/XrayValidationReportHelp>

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:

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

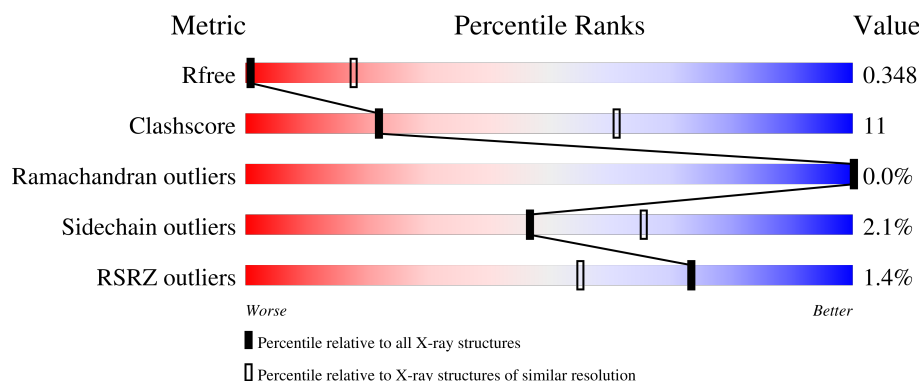
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 4.00 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)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1082 (4.20-3.80)
Clashscore	190562	1129 (4.20-3.80)
Ramachandran outliers	187476	1064 (4.20-3.80)
Sidechain outliers	187428	1055 (4.20-3.80)
RSRZ outliers	180081	1082 (4.20-3.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower 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 electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	764	2% 62% 12% . 26%
1	B	764	% 59% 10% . 30%
1	C	764	% 62% 11% 26%
1	D	764	2% 60% 9% . 30%
1	E	764	% 63% 11% 26%

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Mol	Chain	Length	Quality of chain
1	F	764	% 58% 11% 30%
1	G	764	% 62% 12% 26%
1	H	764	% 59% 10% 30%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 34416 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, 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 Polyketide synthase PksL.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	563	Total	C	N	O	S	0	0	0
			4419	2811	738	844	26			
1	B	533	Total	C	N	O	S	0	0	0
			4185	2663	703	797	22			
1	C	563	Total	C	N	O	S	0	0	0
			4419	2811	738	844	26			
1	D	533	Total	C	N	O	S	0	0	0
			4185	2663	703	797	22			
1	E	563	Total	C	N	O	S	0	0	0
			4419	2811	738	844	26			
1	F	533	Total	C	N	O	S	0	0	0
			4185	2663	703	797	22			
1	G	563	Total	C	N	O	S	0	0	0
			4419	2811	738	844	26			
1	H	533	Total	C	N	O	S	0	0	0
			4185	2663	703	797	22			

There are 160 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-172	MET	-	initiating methionine	UNP Q05470
A	-171	GLY	-	expression tag	UNP Q05470
A	-170	SER	-	expression tag	UNP Q05470
A	-169	SER	-	expression tag	UNP Q05470
A	-168	HIS	-	expression tag	UNP Q05470
A	-167	HIS	-	expression tag	UNP Q05470
A	-166	HIS	-	expression tag	UNP Q05470
A	-165	HIS	-	expression tag	UNP Q05470
A	-164	HIS	-	expression tag	UNP Q05470
A	-163	HIS	-	expression tag	UNP Q05470
A	-162	SER	-	expression tag	UNP Q05470
A	-161	SER	-	expression tag	UNP Q05470
A	-160	GLY	-	expression tag	UNP Q05470

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-159	LEU	-	expression tag	UNP Q05470
A	-158	VAL	-	expression tag	UNP Q05470
A	-157	PRO	-	expression tag	UNP Q05470
A	-156	ARG	-	expression tag	UNP Q05470
A	-155	GLY	-	expression tag	UNP Q05470
A	-154	SER	-	expression tag	UNP Q05470
A	-153	SER	-	expression tag	UNP Q05470
B	-172	MET	-	initiating methionine	UNP Q05470
B	-171	GLY	-	expression tag	UNP Q05470
B	-170	SER	-	expression tag	UNP Q05470
B	-169	SER	-	expression tag	UNP Q05470
B	-168	HIS	-	expression tag	UNP Q05470
B	-167	HIS	-	expression tag	UNP Q05470
B	-166	HIS	-	expression tag	UNP Q05470
B	-165	HIS	-	expression tag	UNP Q05470
B	-164	HIS	-	expression tag	UNP Q05470
B	-163	HIS	-	expression tag	UNP Q05470
B	-162	SER	-	expression tag	UNP Q05470
B	-161	SER	-	expression tag	UNP Q05470
B	-160	GLY	-	expression tag	UNP Q05470
B	-159	LEU	-	expression tag	UNP Q05470
B	-158	VAL	-	expression tag	UNP Q05470
B	-157	PRO	-	expression tag	UNP Q05470
B	-156	ARG	-	expression tag	UNP Q05470
B	-155	GLY	-	expression tag	UNP Q05470
B	-154	SER	-	expression tag	UNP Q05470
B	-153	SER	-	expression tag	UNP Q05470
C	-172	MET	-	initiating methionine	UNP Q05470
C	-171	GLY	-	expression tag	UNP Q05470
C	-170	SER	-	expression tag	UNP Q05470
C	-169	SER	-	expression tag	UNP Q05470
C	-168	HIS	-	expression tag	UNP Q05470
C	-167	HIS	-	expression tag	UNP Q05470
C	-166	HIS	-	expression tag	UNP Q05470
C	-165	HIS	-	expression tag	UNP Q05470
C	-164	HIS	-	expression tag	UNP Q05470
C	-163	HIS	-	expression tag	UNP Q05470
C	-162	SER	-	expression tag	UNP Q05470
C	-161	SER	-	expression tag	UNP Q05470
C	-160	GLY	-	expression tag	UNP Q05470
C	-159	LEU	-	expression tag	UNP Q05470
C	-158	VAL	-	expression tag	UNP Q05470

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-157	PRO	-	expression tag	UNP Q05470
C	-156	ARG	-	expression tag	UNP Q05470
C	-155	GLY	-	expression tag	UNP Q05470
C	-154	SER	-	expression tag	UNP Q05470
C	-153	SER	-	expression tag	UNP Q05470
D	-172	MET	-	initiating methionine	UNP Q05470
D	-171	GLY	-	expression tag	UNP Q05470
D	-170	SER	-	expression tag	UNP Q05470
D	-169	SER	-	expression tag	UNP Q05470
D	-168	HIS	-	expression tag	UNP Q05470
D	-167	HIS	-	expression tag	UNP Q05470
D	-166	HIS	-	expression tag	UNP Q05470
D	-165	HIS	-	expression tag	UNP Q05470
D	-164	HIS	-	expression tag	UNP Q05470
D	-163	HIS	-	expression tag	UNP Q05470
D	-162	SER	-	expression tag	UNP Q05470
D	-161	SER	-	expression tag	UNP Q05470
D	-160	GLY	-	expression tag	UNP Q05470
D	-159	LEU	-	expression tag	UNP Q05470
D	-158	VAL	-	expression tag	UNP Q05470
D	-157	PRO	-	expression tag	UNP Q05470
D	-156	ARG	-	expression tag	UNP Q05470
D	-155	GLY	-	expression tag	UNP Q05470
D	-154	SER	-	expression tag	UNP Q05470
D	-153	SER	-	expression tag	UNP Q05470
E	-172	MET	-	initiating methionine	UNP Q05470
E	-171	GLY	-	expression tag	UNP Q05470
E	-170	SER	-	expression tag	UNP Q05470
E	-169	SER	-	expression tag	UNP Q05470
E	-168	HIS	-	expression tag	UNP Q05470
E	-167	HIS	-	expression tag	UNP Q05470
E	-166	HIS	-	expression tag	UNP Q05470
E	-165	HIS	-	expression tag	UNP Q05470
E	-164	HIS	-	expression tag	UNP Q05470
E	-163	HIS	-	expression tag	UNP Q05470
E	-162	SER	-	expression tag	UNP Q05470
E	-161	SER	-	expression tag	UNP Q05470
E	-160	GLY	-	expression tag	UNP Q05470
E	-159	LEU	-	expression tag	UNP Q05470
E	-158	VAL	-	expression tag	UNP Q05470
E	-157	PRO	-	expression tag	UNP Q05470
E	-156	ARG	-	expression tag	UNP Q05470

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Chain	Residue	Modelled	Actual	Comment	Reference
E	-155	GLY	-	expression tag	UNP Q05470
E	-154	SER	-	expression tag	UNP Q05470
E	-153	SER	-	expression tag	UNP Q05470
F	-172	MET	-	initiating methionine	UNP Q05470
F	-171	GLY	-	expression tag	UNP Q05470
F	-170	SER	-	expression tag	UNP Q05470
F	-169	SER	-	expression tag	UNP Q05470
F	-168	HIS	-	expression tag	UNP Q05470
F	-167	HIS	-	expression tag	UNP Q05470
F	-166	HIS	-	expression tag	UNP Q05470
F	-165	HIS	-	expression tag	UNP Q05470
F	-164	HIS	-	expression tag	UNP Q05470
F	-163	HIS	-	expression tag	UNP Q05470
F	-162	SER	-	expression tag	UNP Q05470
F	-161	SER	-	expression tag	UNP Q05470
F	-160	GLY	-	expression tag	UNP Q05470
F	-159	LEU	-	expression tag	UNP Q05470
F	-158	VAL	-	expression tag	UNP Q05470
F	-157	PRO	-	expression tag	UNP Q05470
F	-156	ARG	-	expression tag	UNP Q05470
F	-155	GLY	-	expression tag	UNP Q05470
F	-154	SER	-	expression tag	UNP Q05470
F	-153	SER	-	expression tag	UNP Q05470
G	-172	MET	-	initiating methionine	UNP Q05470
G	-171	GLY	-	expression tag	UNP Q05470
G	-170	SER	-	expression tag	UNP Q05470
G	-169	SER	-	expression tag	UNP Q05470
G	-168	HIS	-	expression tag	UNP Q05470
G	-167	HIS	-	expression tag	UNP Q05470
G	-166	HIS	-	expression tag	UNP Q05470
G	-165	HIS	-	expression tag	UNP Q05470
G	-164	HIS	-	expression tag	UNP Q05470
G	-163	HIS	-	expression tag	UNP Q05470
G	-162	SER	-	expression tag	UNP Q05470
G	-161	SER	-	expression tag	UNP Q05470
G	-160	GLY	-	expression tag	UNP Q05470
G	-159	LEU	-	expression tag	UNP Q05470
G	-158	VAL	-	expression tag	UNP Q05470
G	-157	PRO	-	expression tag	UNP Q05470
G	-156	ARG	-	expression tag	UNP Q05470
G	-155	GLY	-	expression tag	UNP Q05470
G	-154	SER	-	expression tag	UNP Q05470

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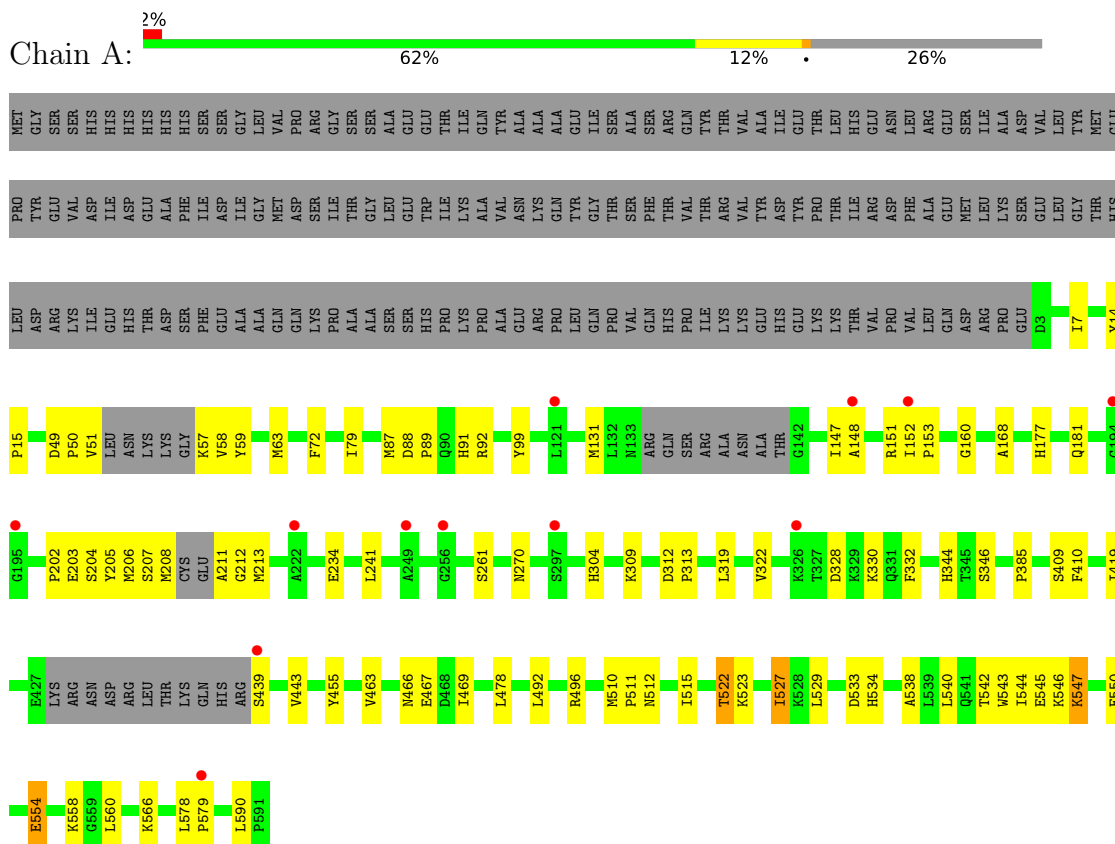
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Chain	Residue	Modelled	Actual	Comment	Reference
G	-153	SER	-	expression tag	UNP Q05470
H	-172	MET	-	initiating methionine	UNP Q05470
H	-171	GLY	-	expression tag	UNP Q05470
H	-170	SER	-	expression tag	UNP Q05470
H	-169	SER	-	expression tag	UNP Q05470
H	-168	HIS	-	expression tag	UNP Q05470
H	-167	HIS	-	expression tag	UNP Q05470
H	-166	HIS	-	expression tag	UNP Q05470
H	-165	HIS	-	expression tag	UNP Q05470
H	-164	HIS	-	expression tag	UNP Q05470
H	-163	HIS	-	expression tag	UNP Q05470
H	-162	SER	-	expression tag	UNP Q05470
H	-161	SER	-	expression tag	UNP Q05470
H	-160	GLY	-	expression tag	UNP Q05470
H	-159	LEU	-	expression tag	UNP Q05470
H	-158	VAL	-	expression tag	UNP Q05470
H	-157	PRO	-	expression tag	UNP Q05470
H	-156	ARG	-	expression tag	UNP Q05470
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H	-153	SER	-	expression tag	UNP Q05470

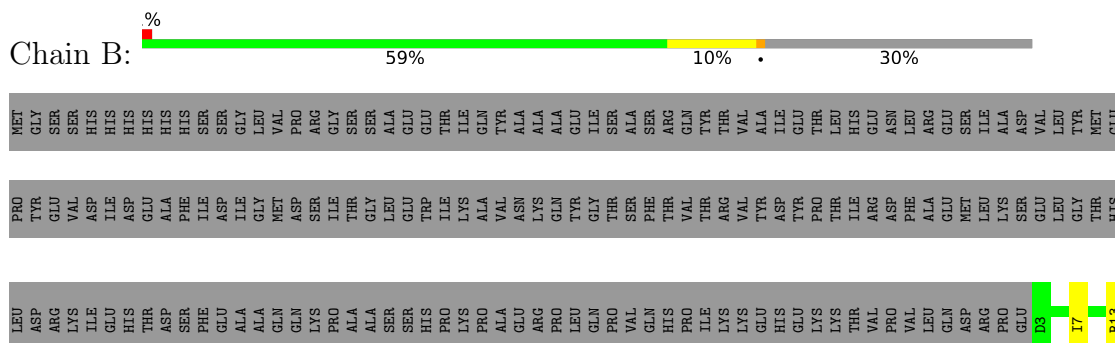
3 Residue-property plots

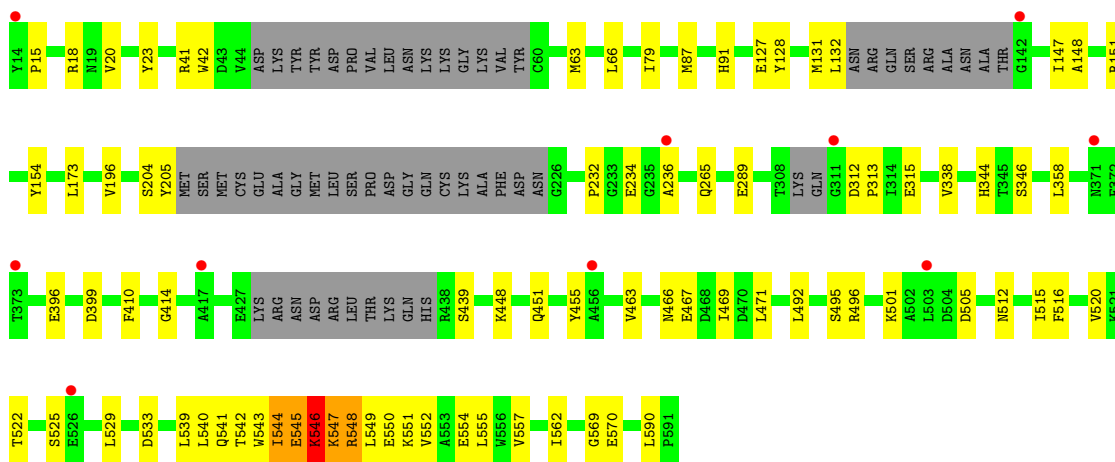
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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 dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). 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: Polyketide synthase PksL

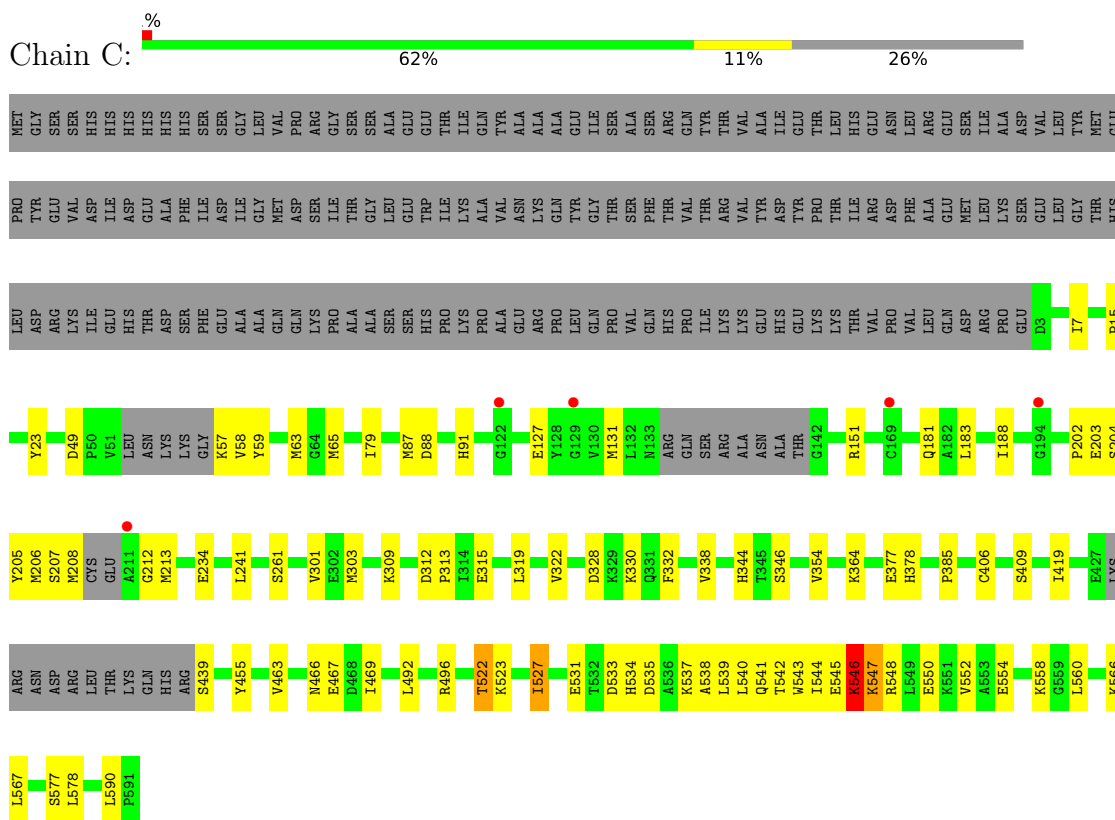


• Molecule 1: Polyketide synthase PksL

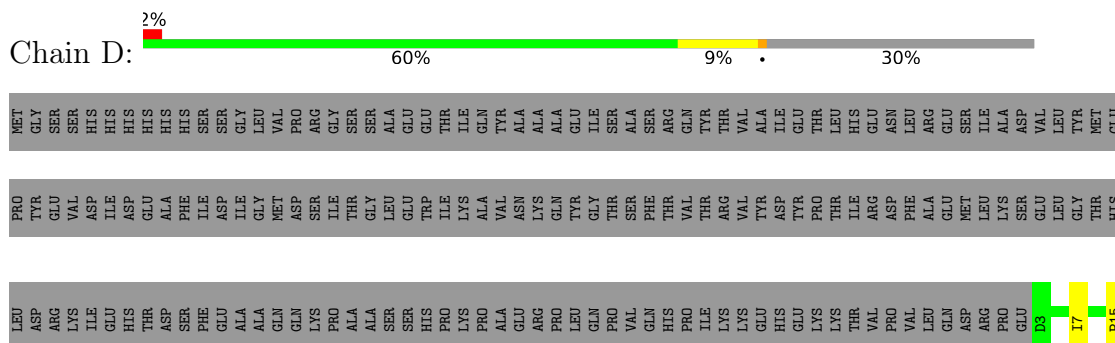


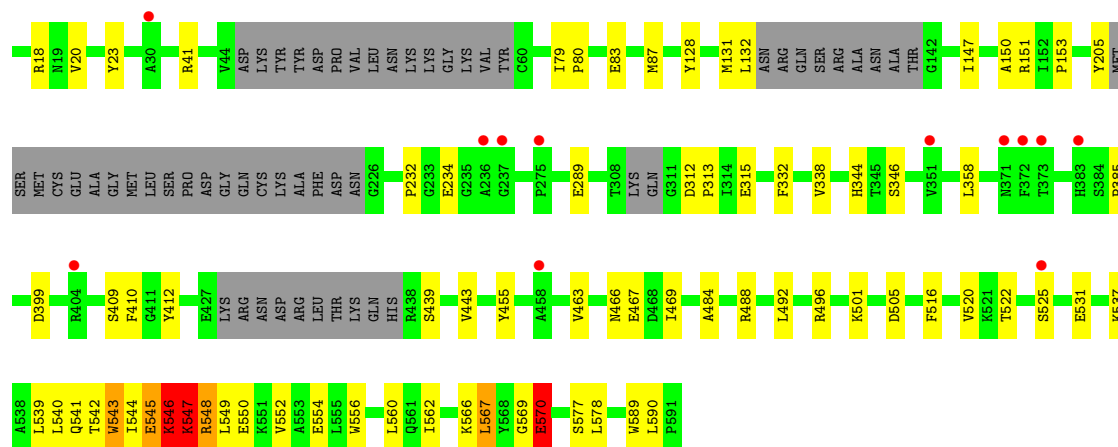


- Molecule 1: Polyketide synthase PksL

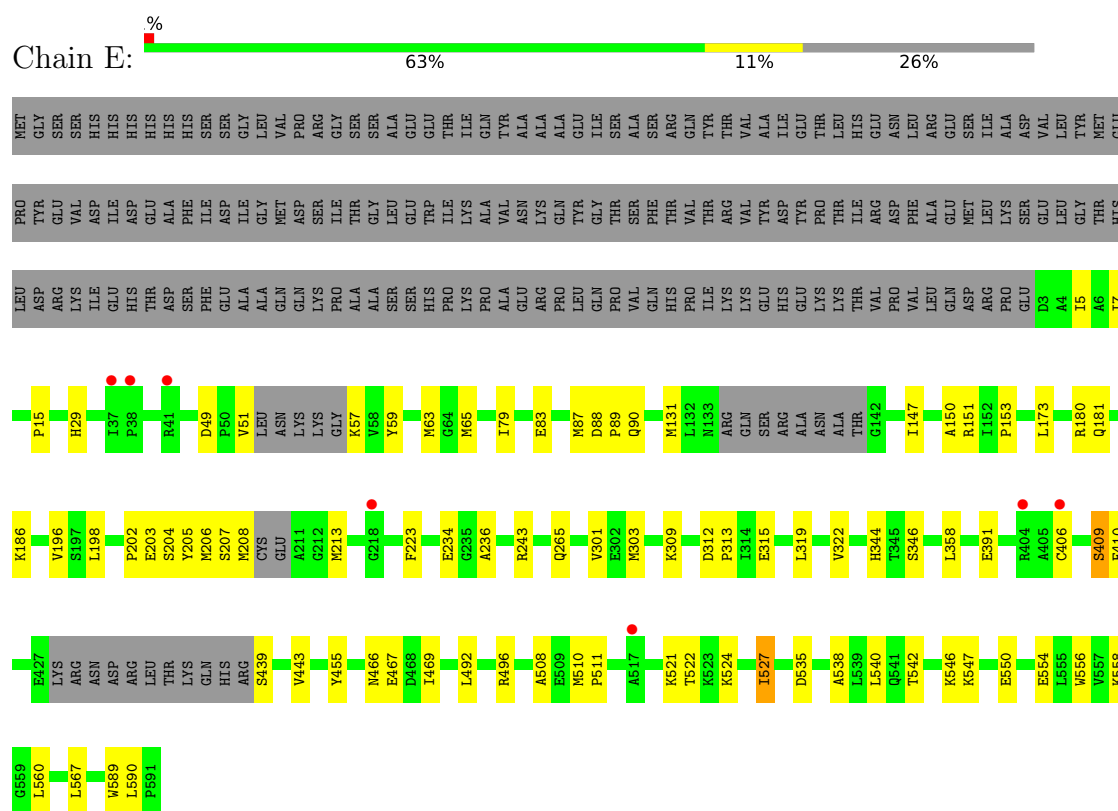


- Molecule 1: Polyketide synthase PksL

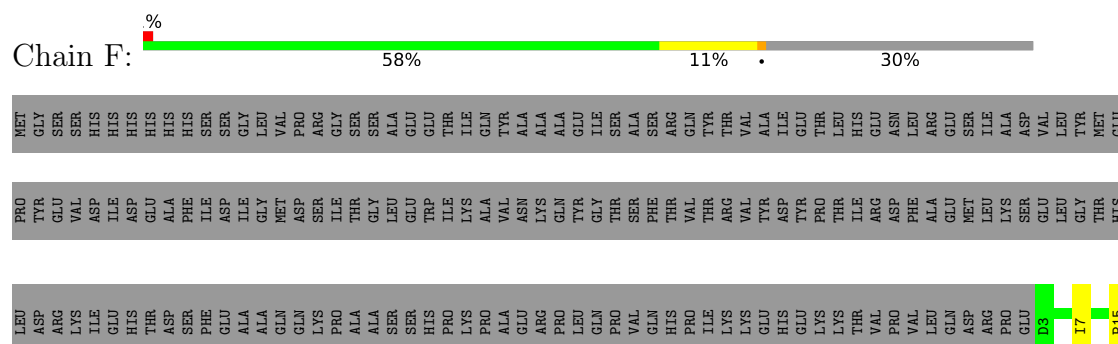


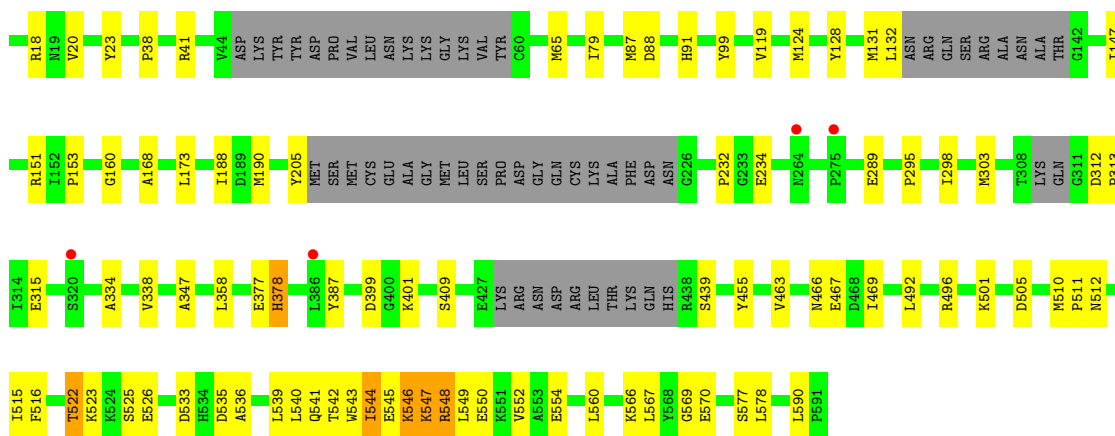


• Molecule 1: Polyketide synthase PksL

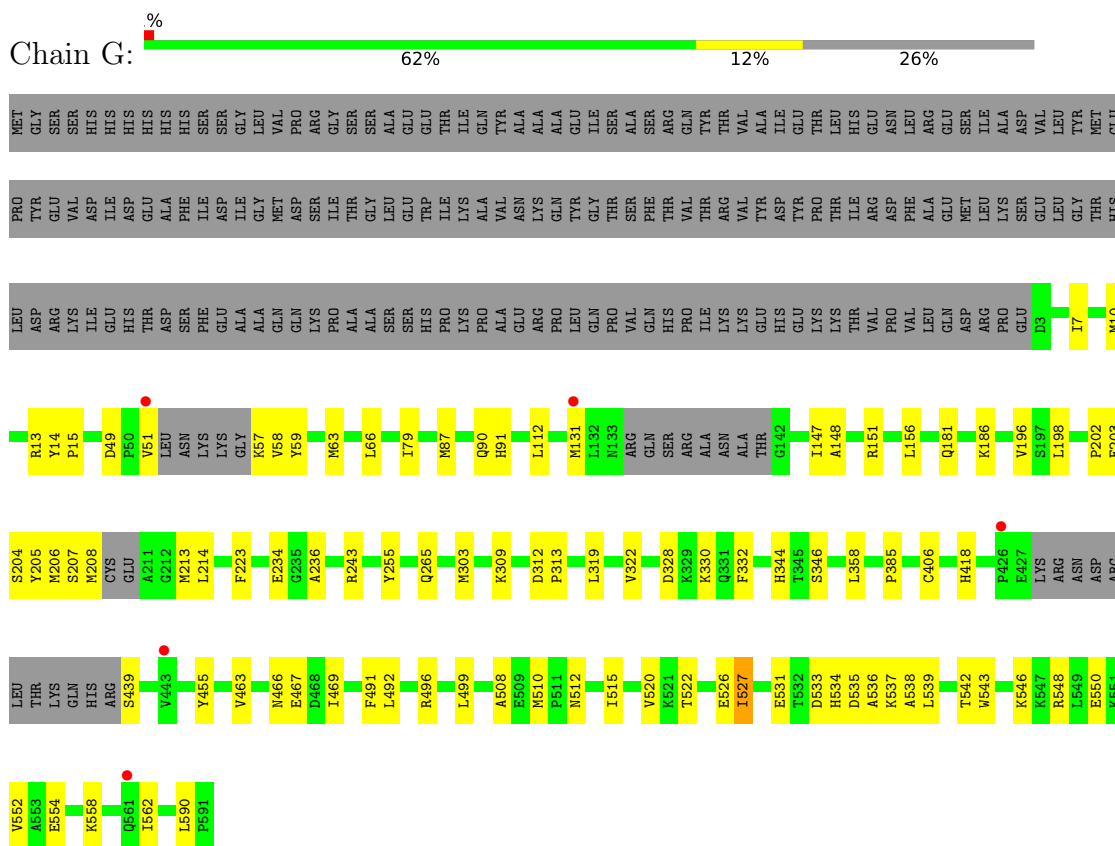


• Molecule 1: Polyketide synthase PksL

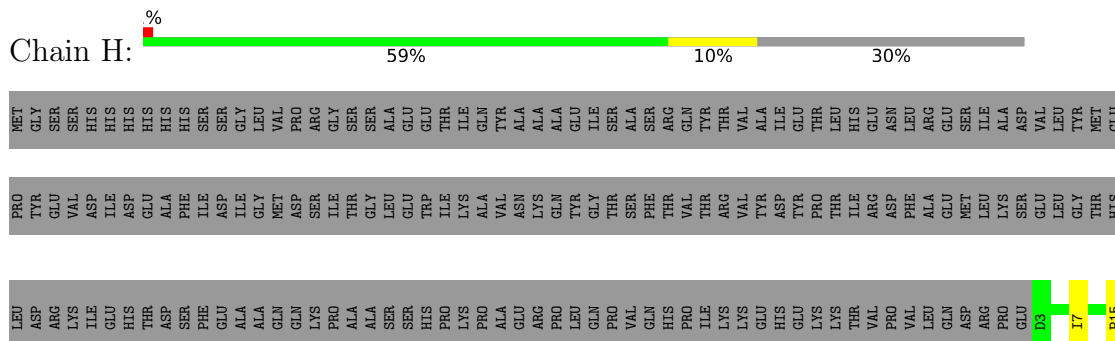


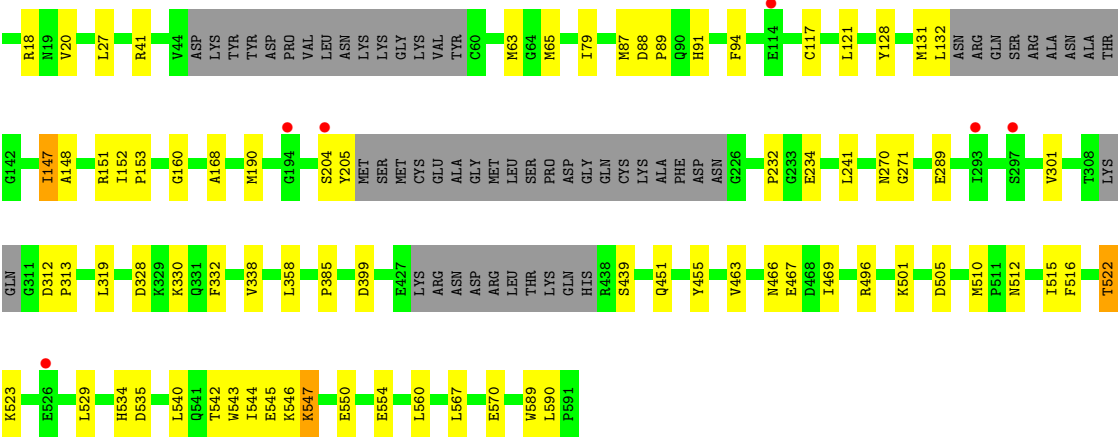


- Molecule 1: Polyketide synthase PksL



- Molecule 1: Polyketide synthase PksL





4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	63.11Å 112.73Å 211.44Å 104.96° 90.07° 106.32°	Depositor
Resolution (Å)	39.77 – 4.00 39.77 – 4.00	Depositor EDS
% Data completeness (in resolution range)	92.6 (39.77-4.00) 92.5 (39.77-4.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.12 (at 4.00Å)	Xtriage
Refinement program	REFMAC 5.8.0107	Depositor
R, R_{free}	0.333 , 0.354 0.327 , 0.348	Depositor DCC
R_{free} test set	2126 reflections (4.69%)	wwPDB-VP
Wilson B-factor (Å ²)	138.8	Xtriage
Anisotropy	0.077	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 593.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	0.347 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	34416	wwPDB-VP
Average B, all atoms (Å ²)	156.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 36.09 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 5.2714e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

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	0.52	0/4518	0.70	1/6110 (0.0%)
1	B	0.53	0/4277	0.80	7/5784 (0.1%)
1	C	0.52	0/4518	0.82	3/6110 (0.0%)
1	D	0.52	0/4277	0.85	11/5784 (0.2%)
1	E	0.52	0/4518	0.68	0/6110
1	F	0.53	1/4277 (0.0%)	0.79	11/5784 (0.2%)
1	G	0.52	0/4518	0.68	0/6110
1	H	0.52	0/4277	0.69	1/5784 (0.0%)
All	All	0.52	1/35180 (0.0%)	0.75	34/47576 (0.1%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1
1	C	0	1
1	D	0	1
1	F	0	1
1	H	0	1
All	All	0	5

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	378	HIS	CE1-NE2	5.30	1.37	1.32

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	546	LYS	CB-CA-C	-32.28	71.52	111.22
1	D	546	LYS	CB-CA-C	-25.97	71.98	111.72

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	546	LYS	N-CA-CB	-19.41	77.51	110.50
1	F	569	GLY	N-CA-C	-17.02	89.25	112.57
1	B	545	GLU	CB-CA-C	15.32	140.91	110.42
1	F	546	LYS	N-CA-C	12.89	147.08	111.00
1	D	546	LYS	N-CA-C	12.81	127.30	109.11
1	B	548	ARG	N-CA-C	12.37	132.62	109.24
1	D	549	LEU	N-CA-C	-11.36	99.24	113.55
1	B	545	GLU	N-CA-C	-11.22	86.90	110.80
1	F	548	ARG	N-CA-C	-10.90	95.09	110.35
1	D	569	GLY	N-CA-C	-9.92	95.09	112.55
1	F	547	LYS	N-CA-CB	-9.76	95.36	110.39
1	F	548	ARG	CB-CA-C	9.39	124.45	109.61
1	D	548	ARG	CB-CA-C	8.96	122.52	110.06
1	D	570	GLU	N-CA-C	8.85	129.66	110.80
1	B	548	ARG	CB-CA-C	-8.25	98.25	112.44
1	D	547	LYS	N-CA-C	7.81	127.43	110.80
1	C	546	LYS	CA-C-N	7.53	139.78	126.45
1	C	546	LYS	C-N-CA	7.53	139.78	126.45
1	B	546	LYS	N-CA-C	7.24	131.28	111.00
1	B	549	LEU	N-CA-C	6.93	125.55	110.80
1	D	570	GLU	N-CA-CB	-6.08	100.21	110.49
1	F	546	LYS	CA-C-N	6.00	130.05	121.71
1	F	546	LYS	C-N-CA	6.00	130.05	121.71
1	D	543	TRP	N-CA-C	5.87	118.81	111.24
1	F	549	LEU	N-CA-C	-5.75	98.55	110.80
1	F	546	LYS	CB-CA-C	-5.41	99.81	110.10
1	D	548	ARG	N-CA-C	-5.34	102.29	110.30
1	F	549	LEU	N-CA-CB	5.29	119.44	110.49
1	A	168	ALA	CB-CA-C	-5.28	110.51	116.63
1	D	545	GLU	CB-CA-C	5.21	118.16	109.56
1	F	168	ALA	CB-CA-C	-5.13	110.68	116.63
1	H	168	ALA	CB-CA-C	-5.02	110.81	116.63

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	570	GLU	Peptide
1	C	546	LYS	Peptide
1	D	570	GLU	Peptide
1	F	570	GLU	Peptide
1	H	570	GLU	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	4419	0	4322	99	2
1	B	4185	0	4102	81	11
1	C	4419	0	4323	93	8
1	D	4185	0	4102	79	2
1	E	4419	0	4323	61	6
1	F	4185	0	4102	200	0
1	G	4419	0	4323	150	0
1	H	4185	0	4102	90	3
All	All	34416	0	33699	726	16

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 11.

All (726) 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:F:535:ASP:CB	1:G:539:LEU:HB2	1.33	1.56
1:F:539:LEU:CD1	1:G:533:ASP:CG	1.80	1.55
1:F:535:ASP:HB2	1:G:539:LEU:CB	1.28	1.52
1:F:545:GLU:CA	1:F:546:LYS:HG3	1.36	1.51
1:H:542:THR:C	1:H:546:LYS:HG2	1.32	1.51
1:A:270:ASN:HD21	1:F:378:HIS:CA	1.23	1.51
1:A:270:ASN:CG	1:F:378:HIS:HA	1.35	1.50
1:F:544:ILE:C	1:F:546:LYS:HG2	1.34	1.50
1:F:545:GLU:HA	1:F:546:LYS:CG	1.38	1.49
1:C:378:HIS:C	1:H:270:ASN:HD21	1.24	1.43
1:H:542:THR:O	1:H:546:LYS:CG	1.68	1.40
1:A:270:ASN:ND2	1:F:378:HIS:HA	1.12	1.40
1:A:270:ASN:OD1	1:F:377:GLU:C	1.65	1.38
1:H:542:THR:CG2	1:H:546:LYS:HD3	1.53	1.38
1:D:540:LEU:HD11	1:D:544:ILE:CD1	1.53	1.38
1:A:270:ASN:HD21	1:F:378:HIS:C	1.34	1.33
1:F:539:LEU:CD2	1:G:536:ALA:HB2	1.57	1.32
1:F:546:LYS:CD	1:F:566:LYS:O	1.77	1.31
1:C:378:HIS:HA	1:H:270:ASN:ND2	1.43	1.31

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:546:LYS:O	1:H:547:LYS:HG3	1.19	1.31
1:F:539:LEU:HD11	1:G:533:ASP:CG	0.88	1.30
1:F:539:LEU:HD22	1:G:536:ALA:CB	1.58	1.30
1:A:270:ASN:ND2	1:F:378:HIS:CA	1.86	1.27
1:F:544:ILE:O	1:F:546:LYS:CG	1.82	1.26
1:D:541:GLN:O	1:D:545:GLU:HB2	1.35	1.26
1:F:546:LYS:HD3	1:F:566:LYS:O	1.14	1.25
1:F:533:ASP:OD2	1:G:539:LEU:CD1	1.85	1.23
1:C:378:HIS:HA	1:H:270:ASN:CG	1.52	1.22
1:H:542:THR:CG2	1:H:546:LYS:CD	2.18	1.20
1:C:378:HIS:CA	1:H:270:ASN:HD21	1.55	1.20
1:H:542:THR:HG22	1:H:546:LYS:CG	1.72	1.19
1:C:378:HIS:CA	1:H:270:ASN:ND2	2.06	1.18
1:F:544:ILE:O	1:F:546:LYS:HG2	1.02	1.17
1:F:539:LEU:HD11	1:G:533:ASP:OD2	1.44	1.17
1:B:540:LEU:O	1:B:544:ILE:HD12	1.44	1.16
1:B:540:LEU:CG	1:B:544:ILE:HD11	1.76	1.16
1:F:539:LEU:HD11	1:G:533:ASP:CB	1.76	1.15
1:H:542:THR:HG23	1:H:546:LYS:CD	1.76	1.13
1:D:540:LEU:CD1	1:D:544:ILE:CD1	2.25	1.12
1:F:539:LEU:CD1	1:G:533:ASP:OD2	1.95	1.12
1:B:544:ILE:C	1:B:546:LYS:HB3	1.72	1.11
1:D:541:GLN:NE2	1:D:545:GLU:OE2	1.84	1.11
1:F:545:GLU:CA	1:F:546:LYS:CG	2.08	1.09
1:B:540:LEU:O	1:B:544:ILE:CD1	2.01	1.09
1:F:539:LEU:CA	1:G:535:ASP:HB2	1.82	1.08
1:B:540:LEU:HG	1:B:544:ILE:CD1	1.82	1.08
1:F:542:THR:CB	1:G:535:ASP:OD2	2.02	1.08
1:H:542:THR:HG22	1:H:546:LYS:HG3	1.33	1.07
1:C:378:HIS:C	1:H:270:ASN:ND2	2.10	1.07
1:F:540:LEU:HD11	1:F:544:ILE:HD11	1.36	1.06
1:A:270:ASN:OD1	1:F:378:HIS:N	1.89	1.05
1:H:542:THR:C	1:H:546:LYS:CG	2.20	1.05
1:A:270:ASN:OD1	1:F:377:GLU:O	1.71	1.05
1:F:539:LEU:HD13	1:G:536:ALA:H	1.20	1.05
1:F:535:ASP:HB2	1:G:539:LEU:CG	1.87	1.04
1:C:541:GLN:O	1:C:545:GLU:HG2	1.55	1.04
1:F:539:LEU:HA	1:G:535:ASP:HB2	1.07	1.03
1:F:543:TRP:CZ2	1:G:533:ASP:HB2	1.92	1.03
1:B:546:LYS:H	1:B:546:LYS:CD	1.67	1.03
1:F:533:ASP:OD2	1:G:539:LEU:HD11	1.57	1.03

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:540:LEU:HG	1:B:544:ILE:HD11	1.06	1.03
1:F:535:ASP:CB	1:G:539:LEU:CB	2.10	1.02
1:B:546:LYS:HD3	1:B:546:LYS:N	1.72	1.02
1:D:540:LEU:CD1	1:D:544:ILE:HD11	1.89	1.02
1:F:544:ILE:C	1:F:546:LYS:CG	2.28	1.02
1:B:546:LYS:H	1:B:546:LYS:HD3	0.88	1.01
1:D:540:LEU:CD1	1:D:544:ILE:HD12	1.87	1.01
1:E:208:MET:C	1:E:213:MET:H	1.69	1.01
1:A:270:ASN:OD1	1:F:378:HIS:HA	1.60	1.00
1:D:544:ILE:HG12	1:D:552:VAL:HG13	1.42	1.00
1:F:540:LEU:CD1	1:F:544:ILE:HD11	1.90	1.00
1:G:208:MET:C	1:G:213:MET:H	1.68	1.00
1:H:546:LYS:O	1:H:547:LYS:CG	2.10	1.00
1:F:542:THR:HB	1:G:535:ASP:OD2	1.59	0.99
1:F:539:LEU:HA	1:G:535:ASP:CB	1.92	0.99
1:H:542:THR:HG22	1:H:546:LYS:CD	1.84	0.99
1:D:540:LEU:HD11	1:D:544:ILE:HD11	1.01	0.98
1:H:542:THR:O	1:H:546:LYS:HG2	0.82	0.98
1:C:208:MET:C	1:C:213:MET:H	1.71	0.98
1:A:270:ASN:ND2	1:F:378:HIS:C	2.15	0.98
1:F:536:ALA:N	1:G:539:LEU:HD22	1.79	0.98
1:C:378:HIS:CA	1:H:270:ASN:CG	2.34	0.98
1:A:270:ASN:CG	1:F:378:HIS:CA	2.26	0.97
1:A:208:MET:C	1:A:213:MET:H	1.73	0.97
1:F:533:ASP:CG	1:G:543:TRP:CZ2	2.43	0.97
1:F:533:ASP:OD2	1:G:539:LEU:HD13	1.65	0.96
1:F:535:ASP:OD2	1:G:539:LEU:HD12	1.65	0.96
1:F:539:LEU:HD11	1:G:533:ASP:OD1	1.64	0.96
1:B:540:LEU:C	1:B:544:ILE:HD12	1.90	0.95
1:C:377:GLU:OE1	1:H:271:GLY:HA2	1.67	0.95
1:F:543:TRP:CE2	1:G:533:ASP:OD2	2.21	0.94
1:A:270:ASN:CG	1:F:377:GLU:O	2.09	0.94
1:D:541:GLN:HE21	1:D:545:GLU:CD	1.74	0.94
1:F:535:ASP:HB3	1:G:539:LEU:HB2	1.47	0.94
1:A:270:ASN:OD1	1:F:378:HIS:CA	2.13	0.94
1:F:205:TYR:CE1	1:F:232:PRO:HG2	2.03	0.93
1:H:205:TYR:CE1	1:H:232:PRO:HG2	2.04	0.93
1:F:545:GLU:N	1:F:546:LYS:HG2	1.83	0.92
1:D:542:THR:O	1:D:546:LYS:CG	2.18	0.92
1:A:542:THR:O	1:A:546:LYS:N	2.03	0.91
1:D:541:GLN:NE2	1:D:545:GLU:CD	2.29	0.91

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:539:LEU:CG	1:G:533:ASP:OD2	2.18	0.91
1:A:270:ASN:HA	1:F:377:GLU:HG2	1.53	0.91
1:D:542:THR:O	1:D:546:LYS:HG2	1.72	0.90
1:A:59:TYR:CE2	1:A:208:MET:HB3	2.06	0.90
1:E:206:MET:O	1:E:208:MET:HE2	1.70	0.90
1:F:533:ASP:CG	1:G:539:LEU:HD11	1.95	0.90
1:G:206:MET:O	1:G:208:MET:HE2	1.71	0.90
1:F:466:ASN:O	1:F:469:ILE:HG13	1.73	0.89
1:C:542:THR:O	1:C:546:LYS:N	2.05	0.88
1:F:539:LEU:CD1	1:G:533:ASP:OD1	2.20	0.87
1:G:202:PRO:O	1:G:206:MET:HB2	1.75	0.86
1:B:540:LEU:CD1	1:B:544:ILE:HD11	2.05	0.86
1:C:377:GLU:CD	1:H:271:GLY:HA2	1.99	0.86
1:B:154:TYR:OH	1:F:377:GLU:CD	2.18	0.86
1:F:546:LYS:CE	1:F:566:LYS:O	2.23	0.85
1:A:58:VAL:HA	1:A:208:MET:HE3	1.58	0.85
1:F:545:GLU:N	1:F:546:LYS:CG	2.39	0.85
1:A:202:PRO:O	1:A:206:MET:HB2	1.76	0.85
1:H:542:THR:HG23	1:H:546:LYS:HD3	0.87	0.84
1:F:535:ASP:CB	1:G:539:LEU:CA	2.56	0.84
1:F:535:ASP:HB2	1:G:539:LEU:CD1	2.08	0.83
1:F:546:LYS:HD2	1:F:567:LEU:HA	1.61	0.83
1:F:539:LEU:HD13	1:G:536:ALA:N	1.93	0.83
1:F:541:GLN:HB3	1:F:545:GLU:OE2	1.79	0.83
1:B:544:ILE:O	1:B:545:GLU:C	2.20	0.82
1:F:533:ASP:OD2	1:G:539:LEU:CD2	2.27	0.82
1:C:59:TYR:CE2	1:C:208:MET:HB3	2.14	0.82
1:E:466:ASN:O	1:E:469:ILE:HG13	1.77	0.82
1:F:535:ASP:HB2	1:G:539:LEU:CA	2.10	0.82
1:H:205:TYR:HE1	1:H:232:PRO:HG2	1.44	0.82
1:H:546:LYS:C	1:H:547:LYS:HG3	2.04	0.82
1:H:128:TYR:O	1:H:132:LEU:HD13	1.80	0.81
1:F:539:LEU:HG	1:G:533:ASP:OD2	1.81	0.81
1:E:202:PRO:O	1:E:206:MET:HB2	1.82	0.80
1:F:543:TRP:HZ2	1:G:533:ASP:HB2	1.44	0.80
1:C:202:PRO:O	1:C:206:MET:HB2	1.81	0.80
1:D:542:THR:O	1:D:546:LYS:CB	2.30	0.79
1:G:550:GLU:O	1:G:554:GLU:HG3	1.82	0.79
1:C:466:ASN:O	1:C:469:ILE:HG13	1.83	0.79
1:F:205:TYR:HE1	1:F:232:PRO:HG2	1.47	0.79
1:D:544:ILE:HG12	1:D:552:VAL:CG1	2.11	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:544:ILE:O	1:F:546:LYS:CD	2.30	0.78
1:G:59:TYR:CE2	1:G:208:MET:HB3	2.19	0.78
1:B:544:ILE:CA	1:B:546:LYS:HB3	2.14	0.78
1:F:535:ASP:OD2	1:G:539:LEU:CD1	2.31	0.78
1:F:546:LYS:HD3	1:F:566:LYS:C	2.06	0.78
1:B:466:ASN:O	1:B:469:ILE:HG13	1.83	0.78
1:A:527:ILE:HD11	1:A:558:LYS:HB2	1.65	0.78
1:D:466:ASN:O	1:D:469:ILE:HG13	1.84	0.78
1:E:550:GLU:O	1:E:554:GLU:HG3	1.84	0.77
1:C:364:LYS:HB2	1:F:511:PRO:HD2	1.65	0.77
1:D:540:LEU:CG	1:D:544:ILE:HD12	2.13	0.77
1:G:466:ASN:O	1:G:469:ILE:HG13	1.84	0.77
1:C:550:GLU:O	1:C:554:GLU:HG3	1.83	0.76
1:F:547:LYS:O	1:F:548:ARG:C	2.27	0.76
1:E:59:TYR:CE2	1:E:208:MET:HB3	2.22	0.75
1:C:206:MET:O	1:C:208:MET:HE2	1.87	0.75
1:H:91:HIS:CE1	1:H:148:ALA:HB2	2.22	0.75
1:H:543:TRP:HA	1:H:546:LYS:HB2	1.68	0.74
1:C:377:GLU:OE1	1:H:271:GLY:CA	2.35	0.74
1:C:539:LEU:HD12	1:C:539:LEU:O	1.87	0.74
1:B:546:LYS:HG2	1:B:548:ARG:H	1.51	0.74
1:E:466:ASN:HB2	1:E:469:ILE:HD11	1.71	0.73
1:F:533:ASP:OD1	1:G:543:TRP:NE1	2.21	0.73
1:G:208:MET:C	1:G:213:MET:N	2.46	0.73
1:B:205:TYR:CE1	1:B:232:PRO:HG2	2.23	0.73
1:C:377:GLU:CD	1:H:271:GLY:CA	2.61	0.73
1:B:128:TYR:CE2	1:B:132:LEU:HD11	2.25	0.72
1:E:79:ILE:HD13	1:E:87:MET:HE1	1.71	0.72
1:F:501:LYS:HE2	1:F:505:ASP:OD2	1.88	0.72
1:D:544:ILE:CD1	1:D:562:ILE:HD12	2.20	0.72
1:D:542:THR:O	1:D:546:LYS:HB2	1.89	0.72
1:A:49:ASP:HB2	1:A:57:LYS:HD3	1.72	0.71
1:A:550:GLU:O	1:A:554:GLU:HG3	1.90	0.71
1:B:131:MET:HE2	1:B:590:LEU:HD11	1.72	0.71
1:H:550:GLU:O	1:H:554:GLU:HG3	1.90	0.71
1:D:466:ASN:HB2	1:D:469:ILE:HD11	1.72	0.71
1:E:527:ILE:HD13	1:E:558:LYS:HB3	1.71	0.71
1:F:545:GLU:HG3	1:F:545:GLU:O	1.89	0.71
1:B:154:TYR:OH	1:F:377:GLU:OE1	2.07	0.70
1:E:206:MET:O	1:E:207:SER:C	2.33	0.70
1:C:206:MET:O	1:C:207:SER:C	2.34	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:535:ASP:CB	1:G:539:LEU:HA	2.20	0.70
1:D:79:ILE:HD13	1:D:87:MET:HE1	1.73	0.70
1:D:540:LEU:HG	1:D:544:ILE:HD12	1.74	0.70
1:H:542:THR:O	1:H:546:LYS:CD	2.38	0.70
1:B:540:LEU:C	1:B:544:ILE:CD1	2.56	0.70
1:F:533:ASP:CG	1:G:543:TRP:HZ2	1.98	0.70
1:C:208:MET:C	1:C:213:MET:N	2.49	0.69
1:F:535:ASP:CA	1:G:539:LEU:HB2	2.20	0.69
1:B:512:ASN:HB3	1:B:515:ILE:HD12	1.73	0.69
1:F:535:ASP:H	1:G:539:LEU:HD13	1.57	0.69
1:F:543:TRP:CE2	1:G:533:ASP:CG	2.70	0.69
1:H:128:TYR:CE2	1:H:132:LEU:HD11	2.27	0.69
1:C:58:VAL:HA	1:C:208:MET:HE3	1.76	0.68
1:D:205:TYR:CE1	1:D:232:PRO:HG2	2.28	0.68
1:E:527:ILE:HD11	1:E:558:LYS:HB2	1.76	0.68
1:D:131:MET:HE2	1:D:590:LEU:HD11	1.76	0.68
1:F:543:TRP:CZ2	1:G:533:ASP:CB	2.75	0.68
1:F:542:THR:HA	1:F:545:GLU:CG	2.24	0.68
1:B:466:ASN:HB2	1:B:469:ILE:HD11	1.75	0.68
1:D:128:TYR:O	1:D:132:LEU:HD13	1.92	0.68
1:D:541:GLN:O	1:D:545:GLU:CB	2.29	0.68
1:E:208:MET:C	1:E:213:MET:N	2.48	0.67
1:F:542:THR:OG1	1:G:535:ASP:OD2	2.12	0.67
1:F:205:TYR:CD1	1:F:232:PRO:HG2	2.29	0.67
1:D:132:LEU:HD12	1:D:590:LEU:CD2	2.24	0.67
1:H:205:TYR:CD1	1:H:232:PRO:HG2	2.29	0.67
1:H:546:LYS:C	1:H:547:LYS:CG	2.66	0.67
1:A:206:MET:O	1:A:207:SER:C	2.37	0.66
1:D:550:GLU:O	1:D:554:GLU:HG3	1.96	0.66
1:E:131:MET:HE2	1:E:590:LEU:HD11	1.78	0.66
1:B:128:TYR:O	1:B:132:LEU:HD13	1.95	0.66
1:C:539:LEU:HD11	1:C:543:TRP:CE2	2.31	0.66
1:G:466:ASN:HB2	1:G:469:ILE:HD11	1.78	0.66
1:F:463:VAL:HA	1:F:469:ILE:CD1	2.26	0.66
1:G:49:ASP:OD1	1:G:51:VAL:HG22	1.95	0.66
1:H:79:ILE:HD13	1:H:87:MET:HE1	1.77	0.66
1:A:208:MET:C	1:A:213:MET:N	2.52	0.65
1:B:546:LYS:HG2	1:B:547:LYS:N	2.11	0.65
1:H:542:THR:CG2	1:H:546:LYS:CE	2.74	0.65
1:G:206:MET:O	1:G:207:SER:C	2.38	0.65
1:A:206:MET:O	1:A:208:MET:HE2	1.96	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:516:PHE:CE1	1:B:554:GLU:HG2	2.31	0.65
1:B:540:LEU:O	1:B:544:ILE:CG1	2.44	0.65
1:F:533:ASP:OD1	1:G:543:TRP:CE2	2.48	0.65
1:D:544:ILE:HD11	1:D:562:ILE:HD12	1.77	0.65
1:F:131:MET:HE2	1:F:590:LEU:HD11	1.79	0.65
1:G:538:ALA:O	1:G:542:THR:HG23	1.97	0.65
1:D:546:LYS:C	1:D:547:LYS:HG3	2.20	0.65
1:F:544:ILE:HG12	1:F:552:VAL:HG22	1.77	0.65
1:F:535:ASP:CB	1:G:539:LEU:CD1	2.75	0.65
1:F:550:GLU:O	1:F:554:GLU:HG3	1.96	0.65
1:A:527:ILE:HD13	1:A:558:LYS:HB3	1.79	0.65
1:F:541:GLN:HB3	1:F:545:GLU:CD	2.21	0.65
1:A:49:ASP:OD1	1:A:51:VAL:HG22	1.97	0.64
1:C:7:ILE:HG12	1:C:241:LEU:CD2	2.28	0.64
1:B:544:ILE:O	1:B:545:GLU:O	2.14	0.64
1:E:527:ILE:CD1	1:E:558:LYS:HB3	2.27	0.64
1:H:131:MET:HE2	1:H:590:LEU:HD11	1.79	0.64
1:B:546:LYS:O	1:B:547:LYS:HB2	1.96	0.64
1:G:527:ILE:HD13	1:G:558:LYS:HB3	1.78	0.64
1:C:79:ILE:HD13	1:C:87:MET:HE1	1.78	0.64
1:A:58:VAL:CA	1:A:208:MET:HE3	2.27	0.64
1:F:533:ASP:OD1	1:G:543:TRP:CZ2	2.51	0.64
1:A:527:ILE:CD1	1:A:558:LYS:CB	2.76	0.63
1:C:364:LYS:HE3	1:F:510:MET:SD	2.38	0.63
1:B:544:ILE:C	1:B:546:LYS:CB	2.62	0.63
1:G:527:ILE:HD11	1:G:558:LYS:HB2	1.80	0.63
1:H:466:ASN:HB2	1:H:469:ILE:HD11	1.79	0.63
1:F:535:ASP:CB	1:G:539:LEU:HD13	2.29	0.63
1:H:516:PHE:CE1	1:H:554:GLU:HG2	2.34	0.63
1:F:533:ASP:OD2	1:G:539:LEU:CG	2.47	0.62
1:F:542:THR:HA	1:F:545:GLU:HG3	1.80	0.62
1:C:364:LYS:CB	1:F:511:PRO:HD2	2.30	0.62
1:E:315:GLU:HG3	1:E:410:PHE:HE2	1.64	0.62
1:H:542:THR:CG2	1:H:546:LYS:CG	2.54	0.62
1:H:542:THR:O	1:H:546:LYS:N	2.32	0.62
1:D:544:ILE:O	1:D:567:LEU:HD12	2.00	0.62
1:F:128:TYR:CE2	1:F:132:LEU:HD11	2.34	0.62
1:F:535:ASP:HB2	1:G:539:LEU:HB2	0.80	0.62
1:C:301:VAL:HG22	1:C:406:CYS:HB2	1.82	0.62
1:G:49:ASP:HB2	1:G:57:LYS:HD3	1.81	0.62
1:D:501:LYS:HE2	1:D:505:ASP:OD2	1.99	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:7:ILE:HD13	1:F:358:LEU:HD11	1.82	0.61
1:G:7:ILE:HD13	1:G:358:LEU:HD11	1.81	0.61
1:A:131:MET:HE2	1:A:590:LEU:HD11	1.81	0.61
1:F:535:ASP:CG	1:G:539:LEU:HA	2.25	0.61
1:F:312:ASP:N	1:F:313:PRO:HD2	2.15	0.61
1:A:91:HIS:CE1	1:A:148:ALA:HB2	2.35	0.61
1:G:63:MET:HE2	1:G:205:TYR:CD2	2.36	0.61
1:G:79:ILE:HD13	1:G:87:MET:HE1	1.83	0.61
1:G:527:ILE:CD1	1:G:558:LYS:CB	2.79	0.61
1:G:508:ALA:HB1	1:G:510:MET:HE3	1.81	0.61
1:C:527:ILE:HD13	1:C:558:LYS:HB3	1.82	0.60
1:F:533:ASP:CB	1:G:543:TRP:HZ2	2.13	0.60
1:F:541:GLN:CB	1:F:545:GLU:OE2	2.48	0.60
1:H:543:TRP:N	1:H:546:LYS:HG2	2.06	0.60
1:B:545:GLU:N	1:B:546:LYS:HB3	2.15	0.60
1:C:541:GLN:O	1:C:545:GLU:CG	2.41	0.60
1:D:544:ILE:CG1	1:D:552:VAL:HG22	2.31	0.60
1:C:463:VAL:HA	1:C:469:ILE:CD1	2.32	0.60
1:A:58:VAL:HB	1:A:208:MET:HE1	1.84	0.60
1:C:202:PRO:O	1:C:206:MET:CB	2.49	0.60
1:F:536:ALA:H	1:G:539:LEU:HD22	1.63	0.60
1:F:539:LEU:HD12	1:G:533:ASP:CG	2.14	0.60
1:G:550:GLU:O	1:G:554:GLU:CG	2.50	0.60
1:A:527:ILE:CD1	1:A:558:LYS:HB3	2.32	0.60
1:C:131:MET:HE2	1:C:590:LEU:HD11	1.84	0.60
1:B:541:GLN:HA	1:B:544:ILE:HD12	1.84	0.60
1:A:466:ASN:HB2	1:A:469:ILE:HD11	1.84	0.59
1:A:543:TRP:HA	1:A:546:LYS:HB3	1.84	0.59
1:E:527:ILE:CD1	1:E:558:LYS:CB	2.80	0.59
1:A:202:PRO:O	1:A:206:MET:CB	2.51	0.59
1:E:202:PRO:O	1:E:206:MET:CB	2.51	0.59
1:F:533:ASP:HB2	1:G:543:TRP:HZ2	1.67	0.59
1:A:550:GLU:O	1:A:554:GLU:CG	2.51	0.59
1:G:527:ILE:CD1	1:G:558:LYS:HB3	2.33	0.59
1:F:533:ASP:OD2	1:G:539:LEU:HD21	2.01	0.59
1:G:202:PRO:O	1:G:206:MET:CB	2.49	0.59
1:F:542:THR:OG1	1:G:535:ASP:CG	2.46	0.59
1:E:547:LYS:HE3	1:E:567:LEU:O	2.03	0.59
1:A:49:ASP:CB	1:A:57:LYS:HD3	2.33	0.58
1:A:319:LEU:O	1:A:322:VAL:HG22	2.04	0.58
1:C:542:THR:O	1:C:546:LYS:HG3	2.03	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:543:TRP:NE1	1:G:533:ASP:CG	2.62	0.58
1:A:546:LYS:O	1:A:547:LYS:HB2	2.03	0.58
1:B:128:TYR:O	1:B:132:LEU:CD1	2.51	0.58
1:A:203:GLU:O	1:A:206:MET:HB3	2.04	0.58
1:C:312:ASP:N	1:C:313:PRO:HD2	2.19	0.58
1:F:128:TYR:O	1:F:132:LEU:HD13	2.04	0.58
1:F:463:VAL:HA	1:F:469:ILE:HD11	1.85	0.58
1:F:492:LEU:N	1:F:492:LEU:HD12	2.20	0.57
1:C:466:ASN:HB2	1:C:469:ILE:HD11	1.85	0.57
1:H:501:LYS:HE2	1:H:505:ASP:OD2	2.04	0.57
1:A:527:ILE:CD1	1:A:558:LYS:HB2	2.33	0.57
1:C:319:LEU:O	1:C:322:VAL:HG22	2.05	0.57
1:F:533:ASP:OD1	1:G:539:LEU:HD11	2.04	0.57
1:C:550:GLU:O	1:C:554:GLU:CG	2.53	0.57
1:G:196:VAL:HG23	1:G:236:ALA:HB2	1.86	0.57
1:H:91:HIS:NE2	1:H:148:ALA:HB2	2.19	0.57
1:F:542:THR:CG2	1:G:535:ASP:OD2	2.52	0.57
1:A:7:ILE:HG12	1:A:241:LEU:CD2	2.34	0.57
1:B:463:VAL:HA	1:B:469:ILE:CD1	2.35	0.57
1:F:535:ASP:HB3	1:G:539:LEU:CA	2.35	0.57
1:G:492:LEU:HD12	1:G:492:LEU:N	2.19	0.57
1:F:540:LEU:HD22	1:F:560:LEU:HD21	1.87	0.56
1:G:131:MET:HE2	1:G:590:LEU:HD11	1.86	0.56
1:G:344:HIS:HD2	1:G:346:SER:H	1.53	0.56
1:A:58:VAL:CB	1:A:208:MET:HE3	2.35	0.56
1:B:544:ILE:CA	1:B:546:LYS:CB	2.83	0.56
1:B:550:GLU:O	1:B:554:GLU:HG3	2.06	0.56
1:D:545:GLU:HG2	1:D:566:LYS:HB3	1.87	0.56
1:A:58:VAL:HB	1:A:208:MET:CE	2.36	0.56
1:D:205:TYR:CD1	1:D:232:PRO:HG2	2.40	0.56
1:G:552:VAL:HG13	1:G:562:ILE:HD12	1.88	0.56
1:F:539:LEU:HB2	1:G:536:ALA:N	2.20	0.56
1:G:512:ASN:HB3	1:G:515:ILE:HD12	1.87	0.56
1:H:128:TYR:O	1:H:132:LEU:CD1	2.53	0.56
1:E:550:GLU:O	1:E:554:GLU:CG	2.53	0.56
1:F:540:LEU:O	1:F:544:ILE:HG13	2.05	0.56
1:G:548:ARG:NH2	1:G:550:GLU:OE1	2.39	0.56
1:A:79:ILE:HD13	1:A:87:MET:HE1	1.87	0.56
1:A:270:ASN:CA	1:F:377:GLU:HG2	2.32	0.56
1:C:303:MET:HE3	1:C:315:GLU:HG2	1.86	0.56
1:D:544:ILE:HD13	1:D:562:ILE:HG23	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:543:TRP:CD2	1:G:533:ASP:OD2	2.59	0.56
1:A:466:ASN:O	1:A:469:ILE:HG13	2.06	0.56
1:C:49:ASP:HB2	1:C:57:LYS:HD3	1.88	0.56
1:D:128:TYR:CE2	1:D:132:LEU:HD11	2.41	0.56
1:F:466:ASN:HB2	1:F:469:ILE:HD11	1.86	0.56
1:F:546:LYS:HD2	1:F:567:LEU:CA	2.35	0.56
1:F:539:LEU:CB	1:G:535:ASP:HB2	2.35	0.55
1:C:181:GLN:OE1	1:C:181:GLN:HA	2.06	0.55
1:C:492:LEU:HD12	1:C:492:LEU:N	2.22	0.55
1:D:544:ILE:HG13	1:D:552:VAL:HG22	1.87	0.55
1:B:205:TYR:CD1	1:B:232:PRO:HG2	2.41	0.55
1:B:205:TYR:HE1	1:B:232:PRO:HG2	1.66	0.55
1:H:205:TYR:HE1	1:H:232:PRO:CG	2.16	0.55
1:F:546:LYS:HE3	1:F:566:LYS:O	2.05	0.55
1:G:344:HIS:CD2	1:G:346:SER:H	2.24	0.55
1:C:377:GLU:OE1	1:H:271:GLY:N	2.40	0.55
1:E:49:ASP:HB2	1:E:57:LYS:HD3	1.88	0.55
1:F:132:LEU:HD12	1:F:590:LEU:CD2	2.36	0.55
1:H:312:ASP:N	1:H:313:PRO:HD2	2.22	0.55
1:B:79:ILE:HD13	1:B:87:MET:HE1	1.87	0.55
1:E:7:ILE:HD13	1:E:358:LEU:HD11	1.87	0.55
1:E:319:LEU:O	1:E:322:VAL:HG22	2.07	0.55
1:A:15:PRO:HD3	1:A:234:GLU:O	2.07	0.54
1:C:88:ASP:HB3	1:C:91:HIS:CD2	2.41	0.54
1:E:63:MET:SD	1:E:65:MET:HG2	2.47	0.54
1:A:204:SER:OG	1:A:205:TYR:N	2.39	0.54
1:F:79:ILE:HD13	1:F:87:MET:HE1	1.88	0.54
1:H:522:THR:HG22	1:H:523:LYS:HG2	1.89	0.54
1:F:303:MET:HE3	1:F:315:GLU:CD	2.33	0.54
1:D:132:LEU:HD12	1:D:590:LEU:HD21	1.90	0.54
1:H:332:PHE:CE2	1:H:385:PRO:HB3	2.43	0.54
1:F:87:MET:HE2	1:F:151:ARG:NH2	2.23	0.54
1:F:545:GLU:O	1:F:545:GLU:CG	2.56	0.54
1:G:467:GLU:HA	1:G:496:ARG:NE	2.23	0.54
1:C:542:THR:HA	1:C:545:GLU:HB2	1.90	0.54
1:G:203:GLU:O	1:G:206:MET:HB3	2.08	0.54
1:G:319:LEU:O	1:G:322:VAL:HG22	2.07	0.54
1:G:539:LEU:HD11	1:G:543:TRP:CE2	2.42	0.54
1:G:527:ILE:CD1	1:G:558:LYS:HB2	2.37	0.54
1:A:205:TYR:CD1	1:A:205:TYR:C	2.86	0.53
1:D:205:TYR:HE1	1:D:232:PRO:HG2	1.73	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:546:LYS:O	1:E:546:LYS:HG3	2.07	0.53
1:F:541:GLN:C	1:F:545:GLU:OE2	2.50	0.53
1:H:466:ASN:O	1:H:469:ILE:HG13	2.08	0.53
1:H:543:TRP:HA	1:H:546:LYS:CG	2.38	0.53
1:C:538:ALA:O	1:C:542:THR:HG23	2.09	0.53
1:H:543:TRP:HA	1:H:546:LYS:CB	2.35	0.53
1:F:539:LEU:HD12	1:G:533:ASP:OD1	2.06	0.53
1:E:204:SER:OG	1:E:205:TYR:N	2.40	0.53
1:H:542:THR:O	1:H:546:LYS:CB	2.52	0.53
1:D:15:PRO:HD3	1:D:234:GLU:O	2.09	0.53
1:F:15:PRO:HD3	1:F:234:GLU:O	2.09	0.53
1:C:204:SER:OG	1:C:205:TYR:N	2.42	0.53
1:C:522:THR:HG22	1:C:523:LYS:HG2	1.90	0.53
1:B:545:GLU:N	1:B:546:LYS:CA	2.72	0.52
1:F:205:TYR:HE1	1:F:232:PRO:CG	2.17	0.52
1:C:58:VAL:HB	1:C:208:MET:HE1	1.91	0.52
1:F:516:PHE:CE1	1:F:554:GLU:HG2	2.44	0.52
1:F:535:ASP:C	1:G:539:LEU:HB2	2.34	0.52
1:H:7:ILE:HD13	1:H:358:LEU:HD11	1.92	0.52
1:H:87:MET:HE2	1:H:151:ARG:NH2	2.24	0.52
1:F:536:ALA:CA	1:G:539:LEU:HD22	2.38	0.52
1:F:542:THR:CA	1:F:545:GLU:HG2	2.35	0.52
1:A:542:THR:O	1:A:546:LYS:HB2	2.10	0.52
1:C:58:VAL:HB	1:C:208:MET:CE	2.39	0.52
1:D:467:GLU:HA	1:D:496:ARG:NE	2.25	0.52
1:C:49:ASP:CB	1:C:57:LYS:HD3	2.39	0.52
1:D:540:LEU:HD22	1:D:560:LEU:HD21	1.91	0.52
1:B:41:ARG:NH1	1:B:234:GLU:OE2	2.36	0.52
1:H:328:ASP:O	1:H:330:LYS:HE2	2.09	0.52
1:A:63:MET:HE2	1:A:205:TYR:CD2	2.45	0.52
1:F:295:PRO:HA	1:F:298:ILE:HD12	1.92	0.52
1:C:364:LYS:NZ	1:F:510:MET:HB3	2.25	0.51
1:D:315:GLU:HG3	1:D:410:PHE:HE2	1.75	0.51
1:D:516:PHE:CE1	1:D:554:GLU:HG2	2.45	0.51
1:E:467:GLU:HA	1:E:496:ARG:NE	2.24	0.51
1:B:552:VAL:HG13	1:B:562:ILE:HD12	1.92	0.51
1:C:467:GLU:HA	1:C:496:ARG:NE	2.25	0.51
1:E:527:ILE:HD11	1:E:558:LYS:CB	2.39	0.51
1:F:88:ASP:HB3	1:F:91:HIS:CD2	2.44	0.51
1:C:205:TYR:CD1	1:C:205:TYR:C	2.88	0.51
1:C:548:ARG:O	1:C:552:VAL:HG23	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:540:LEU:HD22	1:E:560:LEU:HD21	1.90	0.51
1:E:547:LYS:CE	1:E:567:LEU:O	2.59	0.51
1:B:545:GLU:N	1:B:546:LYS:CB	2.73	0.51
1:F:41:ARG:NH1	1:F:234:GLU:OE2	2.33	0.51
1:H:15:PRO:HD3	1:H:234:GLU:O	2.10	0.51
1:E:301:VAL:HG22	1:E:406:CYS:HB2	1.92	0.51
1:G:205:TYR:CD1	1:G:205:TYR:C	2.88	0.51
1:B:128:TYR:CD2	1:B:132:LEU:HD11	2.46	0.51
1:B:492:LEU:HD12	1:B:492:LEU:N	2.25	0.51
1:D:132:LEU:CD1	1:D:590:LEU:CD2	2.88	0.51
1:F:540:LEU:CG	1:F:544:ILE:HD11	2.41	0.51
1:G:91:HIS:CE1	1:G:148:ALA:HB2	2.46	0.51
1:B:23:TYR:OH	1:B:338:VAL:HG21	2.11	0.50
1:E:546:LYS:O	1:E:546:LYS:CG	2.60	0.50
1:D:23:TYR:OH	1:D:338:VAL:HG21	2.12	0.50
1:G:204:SER:OG	1:G:205:TYR:N	2.42	0.50
1:H:132:LEU:HD12	1:H:590:LEU:CD2	2.40	0.50
1:A:59:TYR:CE2	1:A:208:MET:CB	2.89	0.50
1:B:541:GLN:CA	1:B:544:ILE:HD12	2.42	0.50
1:C:539:LEU:HD12	1:C:539:LEU:C	2.36	0.50
1:G:49:ASP:CB	1:G:57:LYS:HD3	2.41	0.50
1:B:546:LYS:HE2	1:B:547:LYS:H	1.76	0.50
1:D:312:ASP:N	1:D:313:PRO:HD2	2.27	0.50
1:E:49:ASP:OD1	1:E:51:VAL:HG22	2.11	0.50
1:E:492:LEU:HD12	1:E:492:LEU:N	2.26	0.50
1:G:58:VAL:HA	1:G:208:MET:HE3	1.94	0.50
1:G:546:LYS:CG	1:G:546:LYS:O	2.59	0.50
1:F:539:LEU:HB2	1:G:535:ASP:C	2.37	0.50
1:A:312:ASP:N	1:A:313:PRO:HD2	2.26	0.50
1:B:344:HIS:CD2	1:B:346:SER:H	2.30	0.50
1:B:467:GLU:HA	1:B:496:ARG:NE	2.26	0.50
1:G:10:MET:HE3	1:G:255:TYR:CD1	2.46	0.50
1:B:540:LEU:O	1:B:544:ILE:HG13	2.12	0.50
1:D:147:ILE:O	1:D:151:ARG:HG2	2.12	0.50
1:E:196:VAL:HG23	1:E:236:ALA:HB2	1.94	0.50
1:C:58:VAL:CB	1:C:208:MET:HE3	2.41	0.49
1:C:58:VAL:CA	1:C:208:MET:HE3	2.42	0.49
1:G:463:VAL:HA	1:G:469:ILE:CD1	2.42	0.49
1:C:328:ASP:O	1:C:330:LYS:HE2	2.12	0.49
1:F:542:THR:HA	1:F:545:GLU:HG2	1.93	0.49
1:B:312:ASP:N	1:B:313:PRO:HD2	2.26	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:203:GLU:O	1:C:206:MET:HB3	2.12	0.49
1:D:545:GLU:CG	1:D:566:LYS:HB3	2.42	0.49
1:F:544:ILE:HG12	1:F:552:VAL:CG2	2.42	0.49
1:H:512:ASN:HB3	1:H:515:ILE:HD12	1.94	0.49
1:A:538:ALA:O	1:A:542:THR:HG23	2.12	0.49
1:D:492:LEU:N	1:D:492:LEU:HD12	2.27	0.49
1:D:544:ILE:HA	1:D:552:VAL:HG21	1.95	0.49
1:A:304:HIS:CE1	1:A:410:PHE:O	2.66	0.49
1:B:463:VAL:HA	1:B:469:ILE:HD12	1.94	0.49
1:C:183:LEU:HD23	1:C:188:ILE:HG13	1.93	0.49
1:F:526:GLU:HG3	1:G:526:GLU:HG2	1.95	0.49
1:C:15:PRO:HD3	1:C:234:GLU:O	2.13	0.49
1:E:150:ALA:C	1:E:153:PRO:HD2	2.38	0.49
1:F:467:GLU:HA	1:F:496:ARG:NE	2.27	0.49
1:C:87:MET:HE2	1:C:151:ARG:NH2	2.28	0.49
1:C:527:ILE:HD11	1:C:558:LYS:HB2	1.94	0.49
1:E:147:ILE:O	1:E:151:ARG:HG2	2.13	0.49
1:F:533:ASP:CB	1:G:543:TRP:CZ2	2.92	0.49
1:A:463:VAL:HA	1:A:469:ILE:CD1	2.42	0.49
1:C:545:GLU:OE2	1:C:566:LYS:HB3	2.13	0.49
1:F:540:LEU:HG	1:F:544:ILE:CD1	2.43	0.49
1:G:181:GLN:HA	1:G:181:GLN:OE1	2.12	0.48
1:H:79:ILE:CD1	1:H:87:MET:HE1	2.44	0.48
1:E:15:PRO:HD3	1:E:234:GLU:O	2.13	0.48
1:C:463:VAL:HA	1:C:469:ILE:HD11	1.94	0.48
1:E:344:HIS:CD2	1:E:346:SER:H	2.31	0.48
1:D:544:ILE:HD13	1:D:562:ILE:HD12	1.92	0.48
1:F:540:LEU:CD1	1:F:544:ILE:CD1	2.79	0.48
1:G:15:PRO:HD3	1:G:234:GLU:O	2.13	0.48
1:G:546:LYS:O	1:G:546:LYS:HG3	2.12	0.48
1:F:23:TYR:OH	1:F:338:VAL:HG21	2.14	0.48
1:F:544:ILE:HG23	1:F:567:LEU:HD13	1.96	0.48
1:G:147:ILE:O	1:G:151:ARG:HG2	2.13	0.48
1:F:132:LEU:HD12	1:F:590:LEU:HD21	1.95	0.48
1:B:265:GLN:NE2	1:B:414:GLY:C	2.72	0.48
1:A:332:PHE:CE2	1:A:385:PRO:HB3	2.49	0.47
1:B:7:ILE:HD13	1:B:358:LEU:HD11	1.94	0.47
1:C:577:SER:O	1:C:578:LEU:HD23	2.14	0.47
1:H:463:VAL:HA	1:H:469:ILE:CD1	2.44	0.47
1:A:540:LEU:HD22	1:A:560:LEU:HD21	1.96	0.47
1:A:542:THR:O	1:A:546:LYS:CB	2.63	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:119:VAL:HG21	1:F:153:PRO:HG3	1.96	0.47
1:F:535:ASP:CG	1:G:539:LEU:CD1	2.86	0.47
1:F:547:LYS:C	1:F:548:ARG:O	2.51	0.47
1:G:91:HIS:NE2	1:G:148:ALA:HB2	2.29	0.47
1:G:527:ILE:HD13	1:G:558:LYS:CB	2.43	0.47
1:A:58:VAL:CB	1:A:208:MET:CE	2.93	0.47
1:A:63:MET:HE2	1:A:205:TYR:HD2	1.80	0.47
1:C:545:GLU:OE2	1:C:566:LYS:HD2	2.14	0.47
1:B:15:PRO:HD3	1:B:234:GLU:O	2.14	0.47
1:F:540:LEU:HG	1:F:544:ILE:HD12	1.97	0.47
1:B:546:LYS:CG	1:B:547:LYS:N	2.76	0.47
1:G:332:PHE:CE2	1:G:385:PRO:HB3	2.50	0.47
1:H:301:VAL:HG21	1:H:319:LEU:HD21	1.95	0.47
1:F:539:LEU:CA	1:G:535:ASP:CB	2.69	0.47
1:F:541:GLN:O	1:F:545:GLU:N	2.48	0.47
1:A:72:PHE:CD1	1:A:92:ARG:HB3	2.50	0.47
1:A:208:MET:O	1:A:211:ALA:HB3	2.15	0.47
1:A:533:ASP:OD1	1:A:534:HIS:N	2.47	0.47
1:A:544:ILE:HG21	1:A:566:LYS:HB2	1.97	0.47
1:F:544:ILE:HG23	1:F:567:LEU:CD1	2.45	0.47
1:G:186:LYS:HD2	1:G:243:ARG:CZ	2.45	0.47
1:E:205:TYR:CD1	1:E:205:TYR:C	2.93	0.46
1:E:223:PHE:CZ	1:E:303:MET:HG3	2.51	0.46
1:A:152:ILE:HB	1:A:153:PRO:HD3	1.97	0.46
1:C:547:LYS:HA	1:C:547:LYS:HD3	1.64	0.46
1:A:328:ASP:O	1:A:330:LYS:HE2	2.15	0.46
1:B:501:LYS:HE2	1:B:505:ASP:OD2	2.15	0.46
1:E:49:ASP:CB	1:E:57:LYS:HD3	2.44	0.46
1:E:88:ASP:OD1	1:E:89:PRO:HD2	2.16	0.46
1:G:208:MET:HA	1:G:214:LEU:HG	1.98	0.46
1:B:147:ILE:O	1:B:151:ARG:HG2	2.15	0.46
1:D:87:MET:HE2	1:D:151:ARG:NH2	2.30	0.46
1:E:173:LEU:HD22	1:E:409:SER:HB2	1.98	0.46
1:B:315:GLU:HG3	1:B:410:PHE:HE2	1.81	0.46
1:F:535:ASP:HB2	1:G:539:LEU:HD13	1.88	0.46
1:F:542:THR:HG21	1:G:535:ASP:OD2	2.16	0.46
1:F:545:GLU:N	1:F:546:LYS:CA	2.79	0.46
1:B:543:TRP:CE2	1:B:551:LYS:HG3	2.51	0.46
1:D:540:LEU:HD12	1:D:544:ILE:CD1	2.34	0.46
1:F:533:ASP:CG	1:G:543:TRP:CE2	2.92	0.46
1:A:14:TYR:HB3	1:A:15:PRO:HD2	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:91:HIS:NE2	1:A:148:ALA:HB2	2.31	0.45
1:A:545:GLU:C	1:A:547:LYS:H	2.24	0.45
1:E:186:LYS:HD2	1:E:243:ARG:CZ	2.46	0.45
1:D:488:ARG:HB3	1:D:556:TRP:CZ3	2.52	0.45
1:D:544:ILE:HA	1:D:552:VAL:CG2	2.46	0.45
1:F:539:LEU:CD1	1:G:533:ASP:CB	2.64	0.45
1:A:87:MET:HE2	1:A:151:ARG:NH2	2.32	0.45
1:C:303:MET:CE	1:C:315:GLU:HG2	2.46	0.45
1:C:533:ASP:OD1	1:C:534:HIS:N	2.49	0.45
1:H:7:ILE:HG12	1:H:241:LEU:CD2	2.46	0.45
1:A:99:TYR:CD1	1:A:99:TYR:C	2.94	0.45
1:G:312:ASP:N	1:G:313:PRO:HD2	2.30	0.45
1:C:540:LEU:HD22	1:C:560:LEU:HD21	1.99	0.45
1:A:50:PRO:HG3	1:E:83:GLU:OE2	2.17	0.45
1:B:196:VAL:HG23	1:B:236:ALA:HB2	1.98	0.45
1:D:312:ASP:N	1:D:313:PRO:CD	2.80	0.45
1:F:535:ASP:HB3	1:G:539:LEU:CB	2.19	0.45
1:E:203:GLU:O	1:E:206:MET:HB3	2.17	0.45
1:C:63:MET:HE2	1:C:205:TYR:CD2	2.51	0.44
1:D:7:ILE:HD13	1:D:358:LEU:HD11	1.98	0.44
1:D:150:ALA:C	1:D:153:PRO:HD2	2.42	0.44
1:D:344:HIS:CD2	1:D:346:SER:H	2.35	0.44
1:E:508:ALA:HB1	1:E:510:MET:HE3	1.99	0.44
1:F:312:ASP:N	1:F:313:PRO:CD	2.78	0.44
1:B:543:TRP:HZ3	1:B:555:LEU:HD11	1.80	0.44
1:G:328:ASP:O	1:G:330:LYS:HE2	2.18	0.44
1:A:545:GLU:C	1:A:547:LYS:N	2.75	0.44
1:C:364:LYS:O	1:F:511:PRO:HG2	2.17	0.44
1:D:332:PHE:CE2	1:D:385:PRO:HB3	2.52	0.44
1:D:516:PHE:HZ	1:D:550:GLU:HG3	1.82	0.44
1:F:522:THR:HG22	1:F:523:LYS:HG2	1.99	0.44
1:E:344:HIS:HD2	1:E:346:SER:H	1.66	0.44
1:G:13:ARG:HG2	1:G:66:LEU:HD22	2.00	0.44
1:F:188:ILE:HD12	1:F:190:MET:O	2.18	0.44
1:A:177:HIS:CE1	1:A:181:GLN:HE21	2.34	0.44
1:A:208:MET:C	1:A:212:GLY:N	2.75	0.44
1:B:42:TRP:CH2	1:B:63:MET:HE1	2.53	0.44
1:B:344:HIS:HD2	1:B:346:SER:H	1.66	0.44
1:B:539:LEU:HD11	1:B:543:TRP:CE2	2.53	0.44
1:E:538:ALA:O	1:E:542:THR:HG23	2.18	0.44
1:F:124:MET:HG2	1:F:347:ALA:HB2	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:463:VAL:HA	1:H:469:ILE:HD12	2.00	0.44
1:A:88:ASP:OD1	1:A:89:PRO:HD2	2.17	0.44
1:G:265:GLN:HB2	1:H:160:GLY:O	2.18	0.44
1:F:501:LYS:HE2	1:F:505:ASP:CG	2.42	0.44
1:H:63:MET:SD	1:H:65:MET:HG2	2.58	0.44
1:D:577:SER:O	1:D:578:LEU:HD23	2.18	0.43
1:F:536:ALA:HB2	1:G:539:LEU:CD2	2.48	0.43
1:A:467:GLU:HA	1:A:496:ARG:NE	2.32	0.43
1:B:91:HIS:CE1	1:B:148:ALA:HB2	2.53	0.43
1:C:344:HIS:CD2	1:C:346:SER:H	2.35	0.43
1:F:535:ASP:OD2	1:G:542:THR:OG1	2.36	0.43
1:B:541:GLN:N	1:B:544:ILE:HD12	2.32	0.43
1:C:354:VAL:HG22	1:C:419:ILE:HD11	2.01	0.43
1:C:377:GLU:CD	1:H:271:GLY:N	2.76	0.43
1:E:90:GLN:HG3	1:E:198:LEU:HD12	2.01	0.43
1:F:535:ASP:N	1:G:539:LEU:HD13	2.27	0.43
1:B:541:GLN:O	1:B:545:GLU:HB2	2.18	0.43
1:A:543:TRP:HD1	1:A:546:LYS:HD3	1.84	0.43
1:A:547:LYS:HD3	1:A:547:LYS:HA	1.67	0.43
1:C:539:LEU:HD21	1:C:543:TRP:CZ2	2.53	0.43
1:D:80:PRO:HD2	1:D:83:GLU:OE1	2.19	0.43
1:D:463:VAL:HA	1:D:469:ILE:CD1	2.49	0.43
1:A:160:GLY:O	1:B:265:GLN:HB2	2.18	0.43
1:F:334:ALA:HB2	1:F:387:TYR:CZ	2.54	0.43
1:D:544:ILE:HG12	1:D:552:VAL:HG22	2.00	0.43
1:E:65:MET:HE3	1:E:65:MET:HB2	1.88	0.43
1:E:443:VAL:HB	1:E:556:TRP:CH2	2.54	0.43
1:F:401:LYS:HB3	1:F:401:LYS:HE2	1.85	0.43
1:A:147:ILE:O	1:A:151:ARG:HG2	2.19	0.43
1:B:128:TYR:CD2	1:B:132:LEU:CD1	3.02	0.43
1:F:544:ILE:O	1:F:546:LYS:HD3	2.15	0.43
1:F:544:ILE:CG2	1:F:567:LEU:HD13	2.49	0.43
1:E:87:MET:HE2	1:E:151:ARG:NH2	2.33	0.43
1:F:516:PHE:HZ	1:F:550:GLU:HG3	1.84	0.43
1:H:27:LEU:HD21	1:H:338:VAL:HG23	2.01	0.43
1:H:516:PHE:HZ	1:H:550:GLU:HG3	1.83	0.43
1:A:59:TYR:HE2	1:A:208:MET:HB3	1.74	0.42
1:E:206:MET:O	1:E:208:MET:N	2.51	0.42
1:F:543:TRP:NE1	1:G:533:ASP:OD2	2.50	0.42
1:G:491:PHE:CE1	1:G:499:LEU:HD12	2.54	0.42
1:D:589:TRP:CD1	1:D:590:LEU:H	2.37	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:112:LEU:HD13	1:G:156:LEU:HD13	2.01	0.42
1:H:152:ILE:HB	1:H:153:PRO:HD3	2.00	0.42
1:H:467:GLU:HA	1:H:496:ARG:NE	2.34	0.42
1:A:181:GLN:OE1	1:A:181:GLN:HA	2.19	0.42
1:A:522:THR:HG22	1:A:523:LYS:HG2	2.01	0.42
1:H:41:ARG:NH1	1:H:234:GLU:OE2	2.33	0.42
1:C:23:TYR:OH	1:C:338:VAL:HG21	2.19	0.42
1:H:544:ILE:O	1:H:567:LEU:HD12	2.19	0.42
1:B:173:LEU:HD12	1:B:173:LEU:HA	1.93	0.42
1:H:94:PHE:CE2	1:H:121:LEU:HD13	2.54	0.42
1:B:91:HIS:NE2	1:B:148:ALA:HB2	2.34	0.42
1:E:5:ILE:HG13	1:E:180:ARG:HG3	2.01	0.42
1:F:541:GLN:HA	1:F:541:GLN:NE2	2.34	0.42
1:H:540:LEU:HD22	1:H:560:LEU:HD21	2.00	0.42
1:A:58:VAL:HG12	1:A:208:MET:HE3	2.01	0.42
1:D:544:ILE:HD11	1:D:562:ILE:CD1	2.47	0.42
1:G:90:GLN:HG3	1:G:198:LEU:HD12	2.01	0.42
1:A:344:HIS:CD2	1:A:346:SER:H	2.38	0.42
1:C:531:GLU:C	1:C:537:LYS:HE3	2.45	0.42
1:D:463:VAL:HA	1:D:469:ILE:HD12	2.01	0.42
1:G:14:TYR:HB3	1:G:15:PRO:HD2	2.00	0.42
1:G:223:PHE:CZ	1:G:303:MET:HG3	2.55	0.42
1:A:79:ILE:CD1	1:A:87:MET:HE1	2.50	0.42
1:A:510:MET:HA	1:A:511:PRO:HD3	1.90	0.42
1:F:463:VAL:HA	1:F:469:ILE:HD12	1.99	0.42
1:G:463:VAL:HA	1:G:469:ILE:HD12	2.01	0.42
1:G:533:ASP:OD1	1:G:534:HIS:N	2.52	0.42
1:C:208:MET:C	1:C:212:GLY:N	2.78	0.42
1:D:443:VAL:HB	1:D:556:TRP:CH2	2.55	0.42
1:F:38:PRO:HD3	1:F:65:MET:HE1	2.01	0.42
1:H:543:TRP:CA	1:H:546:LYS:HG2	2.50	0.42
1:H:589:TRP:CG	1:H:590:LEU:H	2.38	0.42
1:D:589:TRP:CG	1:D:590:LEU:H	2.38	0.41
1:E:510:MET:HA	1:E:511:PRO:HD3	1.90	0.41
1:F:512:ASN:HB3	1:F:515:ILE:HD12	2.00	0.41
1:F:535:ASP:CG	1:G:539:LEU:HD13	2.45	0.41
1:H:589:TRP:CD1	1:H:590:LEU:H	2.37	0.41
1:B:41:ARG:HH12	1:B:234:GLU:CD	2.26	0.41
1:B:542:THR:O	1:B:546:LYS:HA	2.20	0.41
1:C:59:TYR:CE2	1:C:208:MET:CB	2.96	0.41
1:C:544:ILE:HG23	1:C:567:LEU:HD13	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:512:ASN:HB3	1:A:515:ILE:HD12	2.02	0.41
1:B:471:LEU:HD11	1:B:495:SER:C	2.46	0.41
1:F:577:SER:O	1:F:578:LEU:HD23	2.20	0.41
1:B:540:LEU:HD12	1:B:544:ILE:HD11	1.98	0.41
1:C:63:MET:SD	1:C:65:MET:HG2	2.60	0.41
1:D:484:ALA:HB1	1:D:488:ARG:CZ	2.50	0.41
1:E:312:ASP:N	1:E:313:PRO:HD2	2.36	0.41
1:F:99:TYR:CD1	1:F:99:TYR:C	2.98	0.41
1:H:88:ASP:OD1	1:H:89:PRO:HD2	2.21	0.41
1:B:127:GLU:OE2	1:B:204:SER:HB3	2.20	0.41
1:C:332:PHE:CE2	1:C:385:PRO:HB3	2.55	0.41
1:C:527:ILE:CD1	1:C:558:LYS:HB3	2.49	0.41
1:B:13:ARG:HG2	1:B:66:LEU:HD22	2.02	0.41
1:D:41:ARG:NH1	1:D:234:GLU:OE2	2.37	0.41
1:F:173:LEU:HD12	1:F:173:LEU:HA	1.94	0.41
1:H:147:ILE:O	1:H:151:ARG:HG2	2.20	0.41
1:H:204:SER:O	1:H:204:SER:OG	2.33	0.41
1:A:540:LEU:O	1:A:544:ILE:HG13	2.21	0.41
1:C:364:LYS:HZ2	1:F:510:MET:HB3	1.86	0.41
1:D:501:LYS:HE2	1:D:505:ASP:CG	2.46	0.41
1:D:544:ILE:CG1	1:D:552:VAL:HG13	2.31	0.41
1:E:589:TRP:CD1	1:E:590:LEU:H	2.39	0.41
1:H:312:ASP:N	1:H:313:PRO:CD	2.83	0.41
1:H:542:THR:HG22	1:H:546:LYS:CE	2.45	0.41
1:A:208:MET:C	1:A:212:GLY:H	2.29	0.41
1:A:443:VAL:O	1:A:478:LEU:HB3	2.21	0.41
1:C:261:SER:HB2	1:C:419:ILE:HG22	2.02	0.41
1:D:531:GLU:O	1:D:537:LYS:HE3	2.20	0.41
1:G:406:CYS:HB3	1:G:418:HIS:CE1	2.56	0.41
1:G:531:GLU:C	1:G:537:LYS:HE3	2.46	0.41
1:H:128:TYR:CD2	1:H:132:LEU:HD11	2.54	0.41
1:D:412:TYR:CD1	1:D:412:TYR:C	2.97	0.41
1:D:546:LYS:HA	1:D:546:LYS:HE3	2.03	0.41
1:E:521:LYS:NZ	1:E:524:LYS:NZ	2.69	0.41
1:H:542:THR:CG2	1:H:546:LYS:HE2	2.50	0.41
1:A:578:LEU:HB3	1:A:579:PRO:HD2	2.03	0.41
1:C:364:LYS:HB2	1:F:511:PRO:CD	2.42	0.41
1:E:265:GLN:HB2	1:F:160:GLY:O	2.20	0.41
1:F:543:TRP:CZ2	1:G:533:ASP:OD2	2.72	0.41
1:E:181:GLN:OE1	1:E:181:GLN:HA	2.21	0.40
1:H:543:TRP:C	1:H:545:GLU:N	2.78	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:261:SER:HB2	1:A:419:ILE:HG22	2.03	0.40
1:A:492:LEU:N	1:A:492:LEU:HD12	2.37	0.40
1:H:117:CYS:HA	1:H:190:MET:O	2.21	0.40
1:H:132:LEU:HD12	1:H:590:LEU:HD21	2.02	0.40
1:A:270:ASN:ND2	1:F:378:HIS:O	2.47	0.40
1:C:58:VAL:HG12	1:C:208:MET:HE3	2.03	0.40
1:C:127:GLU:OE2	1:C:204:SER:OG	2.29	0.40
1:D:547:LYS:O	1:D:548:ARG:C	2.65	0.40
1:B:448:LYS:HB2	1:B:451:GLN:OE1	2.21	0.40
1:B:554:GLU:O	1:B:557:VAL:HG22	2.22	0.40

All (16) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:547:LYS:O	1:E:29:HIS:NE2[1_665]	1.53	0.67
1:B:547:LYS:O	1:E:29:HIS:CD2[1_665]	1.78	0.42
1:B:539:LEU:CB	1:C:535:ASP:CB[1_655]	1.88	0.32
1:B:539:LEU:CD2	1:C:533:ASP:OD2[1_655]	1.93	0.27
1:E:546:LYS:NZ	1:H:535:ASP:OD2[1_656]	1.96	0.24
1:B:539:LEU:CD1	1:C:533:ASP:OD1[1_655]	2.00	0.20
1:B:543:TRP:CZ2	1:C:533:ASP:CG[1_655]	2.04	0.16
1:B:543:TRP:CH2	1:C:533:ASP:OD2[1_655]	2.05	0.15
1:C:537:LYS:NZ	1:E:391:GLU:OE2[1_565]	2.06	0.14
1:B:547:LYS:C	1:E:29:HIS:NE2[1_665]	2.08	0.12
1:B:543:TRP:CZ2	1:C:533:ASP:OD2[1_655]	2.09	0.11
1:A:529:LEU:CD2	1:D:539:LEU:CD2[1_556]	2.10	0.10
1:B:533:ASP:OD1	1:C:539:LEU:CD1[1_655]	2.11	0.09
1:E:535:ASP:CG	1:H:543:TRP:CZ2[1_656]	2.11	0.09
1:A:529:LEU:CD1	1:D:543:TRP:CZ2[1_556]	2.16	0.04
1:B:396:GLU:OE2	1:H:534:HIS:CD2[1_666]	2.18	0.02

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

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	553/764 (72%)	541 (98%)	12 (2%)	0	100	100
1	B	521/764 (68%)	503 (96%)	17 (3%)	1 (0%)	43	75
1	C	553/764 (72%)	539 (98%)	14 (2%)	0	100	100
1	D	521/764 (68%)	508 (98%)	12 (2%)	1 (0%)	43	75
1	E	553/764 (72%)	541 (98%)	12 (2%)	0	100	100
1	F	521/764 (68%)	508 (98%)	13 (2%)	0	100	100
1	G	553/764 (72%)	540 (98%)	13 (2%)	0	100	100
1	H	521/764 (68%)	507 (97%)	14 (3%)	0	100	100
All	All	4296/6112 (70%)	4187 (98%)	107 (2%)	2 (0%)	100	100

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	570	GLU
1	B	569	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 X-ray entries followed by that with respect to entries of similar resolution.

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	468/644 (73%)	460 (98%)	8 (2%)	53	68
1	B	442/644 (69%)	429 (97%)	13 (3%)	37	58
1	C	468/644 (73%)	461 (98%)	7 (2%)	57	70
1	D	442/644 (69%)	429 (97%)	13 (3%)	37	58
1	E	468/644 (73%)	462 (99%)	6 (1%)	61	72
1	F	442/644 (69%)	431 (98%)	11 (2%)	42	62
1	G	468/644 (73%)	462 (99%)	6 (1%)	61	72
1	H	442/644 (69%)	430 (97%)	12 (3%)	39	60

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	3640/5152 (71%)	3564 (98%)	76 (2%)	47 65

All (76) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	309	LYS
1	A	409	SER
1	A	439	SER
1	A	455	TYR
1	A	522	THR
1	A	527	ILE
1	A	547	LYS
1	A	554	GLU
1	B	18	ARG
1	B	20	VAL
1	B	289	GLU
1	B	399	ASP
1	B	439	SER
1	B	455	TYR
1	B	520	VAL
1	B	522	THR
1	B	525	SER
1	B	529	LEU
1	B	544	ILE
1	B	546	LYS
1	B	547	LYS
1	C	309	LYS
1	C	409	SER
1	C	439	SER
1	C	455	TYR
1	C	522	THR
1	C	527	ILE
1	C	547	LYS
1	D	18	ARG
1	D	20	VAL
1	D	289	GLU
1	D	399	ASP
1	D	409	SER
1	D	439	SER
1	D	455	TYR
1	D	520	VAL
1	D	522	THR

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Mol	Chain	Res	Type
1	D	525	SER
1	D	546	LYS
1	D	547	LYS
1	D	567	LEU
1	E	309	LYS
1	E	409	SER
1	E	439	SER
1	E	455	TYR
1	E	522	THR
1	E	527	ILE
1	F	18	ARG
1	F	20	VAL
1	F	147	ILE
1	F	289	GLU
1	F	399	ASP
1	F	409	SER
1	F	439	SER
1	F	455	TYR
1	F	522	THR
1	F	525	SER
1	F	544	ILE
1	G	309	LYS
1	G	439	SER
1	G	455	TYR
1	G	520	VAL
1	G	522	THR
1	G	527	ILE
1	H	18	ARG
1	H	20	VAL
1	H	147	ILE
1	H	289	GLU
1	H	399	ASP
1	H	439	SER
1	H	451	GLN
1	H	455	TYR
1	H	510	MET
1	H	522	THR
1	H	529	LEU
1	H	547	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	91	HIS
1	A	126	ASN
1	A	378	HIS
1	A	416	ASN
1	B	304	HIS
1	B	331	GLN
1	B	344	HIS
1	B	416	ASN
1	B	534	HIS
1	C	91	HIS
1	C	185	ASN
1	D	331	GLN
1	D	416	ASN
1	D	541	GLN
1	E	91	HIS
1	E	378	HIS
1	E	416	ASN
1	F	91	HIS
1	F	331	GLN
1	F	341	ASN
1	F	534	HIS
1	F	541	GLN
1	G	344	HIS
1	G	389	ASN
1	G	416	ASN
1	H	416	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	563/764 (73%)	-0.24	12 (2%) 63 47	24, 154, 186, 226	0
1	B	533/764 (69%)	-0.24	10 (1%) 66 49	24, 156, 195, 222	0
1	C	563/764 (73%)	-0.32	5 (0%) 81 64	24, 153, 185, 226	0
1	D	533/764 (69%)	-0.23	12 (2%) 61 45	24, 156, 191, 216	0
1	E	563/764 (73%)	-0.33	7 (1%) 76 59	100, 153, 187, 213	0
1	F	533/764 (69%)	-0.29	4 (0%) 82 65	24, 156, 189, 216	0
1	G	563/764 (73%)	-0.30	5 (0%) 81 64	101, 153, 186, 206	0
1	H	533/764 (69%)	-0.30	6 (1%) 78 60	24, 157, 193, 225	0
All	All	4384/6112 (71%)	-0.28	61 (1%) 73 55	24, 155, 189, 226	0

All (61) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	256	GLY	5.0
1	B	371	ASN	5.0
1	B	526	GLU	4.9
1	A	249	ALA	4.9
1	D	372	PHE	4.2
1	E	37	ILE	3.9
1	D	525	SER	3.9
1	D	275	PRO	3.8
1	D	371	ASN	3.7
1	A	194	GLY	3.7
1	D	404	ARG	3.4
1	C	129	GLY	3.4
1	B	373	THR	3.2
1	E	38	PRO	3.2
1	B	503	LEU	3.1
1	F	386	LEU	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	439	SER	3.0
1	G	131	MET	2.9
1	D	30	ALA	2.9
1	D	373	THR	2.9
1	B	142	GLY	2.9
1	B	14	TYR	2.8
1	B	456	ALA	2.8
1	B	417	ALA	2.7
1	A	326	LYS	2.6
1	H	297	SER	2.6
1	G	561	GLN	2.6
1	D	383	HIS	2.6
1	E	406	CYS	2.6
1	E	218	GLY	2.6
1	C	211	ALA	2.5
1	A	148	ALA	2.5
1	E	517	ALA	2.4
1	C	169	CYS	2.4
1	G	443	VAL	2.4
1	A	195	GLY	2.4
1	B	236	ALA	2.4
1	H	293	ILE	2.4
1	F	275	PRO	2.3
1	G	426	PRO	2.3
1	C	194	GLY	2.3
1	G	51	VAL	2.3
1	D	237	GLY	2.3
1	E	404	ARG	2.3
1	D	236	ALA	2.3
1	A	152	ILE	2.2
1	H	114	GLU	2.2
1	D	351	VAL	2.2
1	C	122	GLY	2.2
1	F	264	ASN	2.2
1	F	320	SER	2.2
1	A	121	LEU	2.1
1	A	222	ALA	2.1
1	E	41	ARG	2.1
1	A	579	PRO	2.1
1	H	526	GLU	2.1
1	B	311	GLY	2.1
1	D	458	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	297	SER	2.1
1	H	194	GLY	2.1
1	H	204	SER	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.