



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 18, 2026 – 10:49 PM UTC

PDB ID : 4WS9 / pdb_00004ws9
Title : Crystal structure of sMAT N159G from *Sulfolobus solfataricus*
Authors : Wang, F.; Brady, E.L.; Singh, S.; Clinger, J.A.; Huber, T.D.; Thorson, J.S.;
Phillips Jr., G.N.
Deposited on : 2014-10-26
Resolution : 2.80 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
Mogul : 2022.3.0, CSD as543be (2022)
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

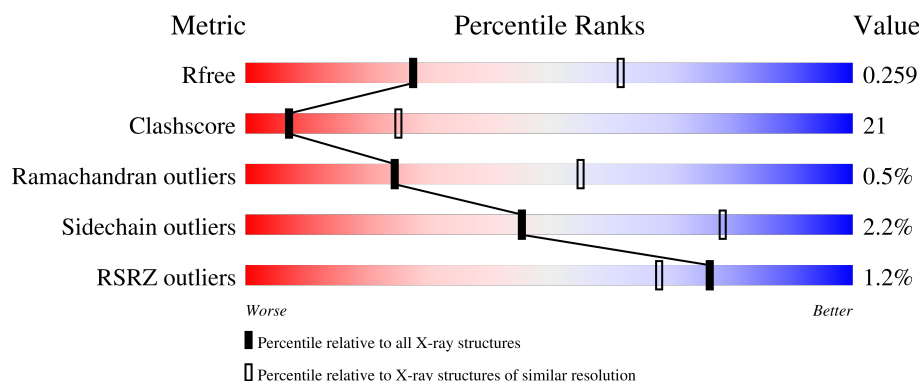
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	413	% 58% 34% • 5%
1	B	413	% 54% 38% • 5%
1	C	413	 61% 31% • 5%
1	D	413	% 51% 39% 5% 5%
1	E	413	 53% 39% • 5%

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Mol	Chain	Length	Quality of chain
1	F	413	
1	G	413	
1	H	413	
1	I	413	
1	J	413	
1	K	413	
1	L	413	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	PO4	A	502	-	-	X	-
2	PO4	D	502	-	-	-	X
2	PO4	G	501	-	-	X	-
2	PO4	L	501	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 37566 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 S-adenosylmethionine synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	391	Total	C	N	O	S	0	0	0
			3043	1928	523	585	7			
1	B	392	Total	C	N	O	S	0	0	0
			3046	1930	524	586	6			
1	C	391	Total	C	N	O	S	0	0	0
			3043	1928	523	585	7			
1	D	392	Total	C	N	O	S	0	0	0
			3046	1930	524	586	6			
1	E	391	Total	C	N	O	S	0	0	0
			3043	1928	523	585	7			
1	F	392	Total	C	N	O	S	0	0	0
			3046	1930	524	586	6			
1	G	391	Total	C	N	O	S	0	0	0
			3043	1928	523	585	7			
1	H	392	Total	C	N	O	S	0	0	0
			3046	1930	524	586	6			
1	I	391	Total	C	N	O	S	0	0	0
			3043	1928	523	585	7			
1	J	392	Total	C	N	O	S	0	0	0
			3046	1930	524	586	6			
1	K	391	Total	C	N	O	S	0	0	0
			3043	1928	523	585	7			
1	L	392	Total	C	N	O	S	0	0	0
			3046	1930	524	586	6			

There are 120 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-8	GLY	-	expression tag	UNP Q980S9
A	-7	SER	-	expression tag	UNP Q980S9
A	-6	HIS	-	expression tag	UNP Q980S9
A	-5	MET	-	expression tag	UNP Q980S9
A	-4	PHE	-	expression tag	UNP Q980S9

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	GLU	-	expression tag	UNP Q980S9
A	-2	GLU	-	expression tag	UNP Q980S9
A	-1	PRO	-	expression tag	UNP Q980S9
A	0	ILE	-	expression tag	UNP Q980S9
A	159	GLY	ASN	conflict	UNP Q980S9
B	-8	GLY	-	expression tag	UNP Q980S9
B	-7	SER	-	expression tag	UNP Q980S9
B	-6	HIS	-	expression tag	UNP Q980S9
B	-5	MET	-	expression tag	UNP Q980S9
B	-4	PHE	-	expression tag	UNP Q980S9
B	-3	GLU	-	expression tag	UNP Q980S9
B	-2	GLU	-	expression tag	UNP Q980S9
B	-1	PRO	-	expression tag	UNP Q980S9
B	0	ILE	-	expression tag	UNP Q980S9
B	159	GLY	ASN	conflict	UNP Q980S9
C	-8	GLY	-	expression tag	UNP Q980S9
C	-7	SER	-	expression tag	UNP Q980S9
C	-6	HIS	-	expression tag	UNP Q980S9
C	-5	MET	-	expression tag	UNP Q980S9
C	-4	PHE	-	expression tag	UNP Q980S9
C	-3	GLU	-	expression tag	UNP Q980S9
C	-2	GLU	-	expression tag	UNP Q980S9
C	-1	PRO	-	expression tag	UNP Q980S9
C	0	ILE	-	expression tag	UNP Q980S9
C	159	GLY	ASN	conflict	UNP Q980S9
D	-8	GLY	-	expression tag	UNP Q980S9
D	-7	SER	-	expression tag	UNP Q980S9
D	-6	HIS	-	expression tag	UNP Q980S9
D	-5	MET	-	expression tag	UNP Q980S9
D	-4	PHE	-	expression tag	UNP Q980S9
D	-3	GLU	-	expression tag	UNP Q980S9
D	-2	GLU	-	expression tag	UNP Q980S9
D	-1	PRO	-	expression tag	UNP Q980S9
D	0	ILE	-	expression tag	UNP Q980S9
D	159	GLY	ASN	conflict	UNP Q980S9
E	-8	GLY	-	expression tag	UNP Q980S9
E	-7	SER	-	expression tag	UNP Q980S9
E	-6	HIS	-	expression tag	UNP Q980S9
E	-5	MET	-	expression tag	UNP Q980S9
E	-4	PHE	-	expression tag	UNP Q980S9
E	-3	GLU	-	expression tag	UNP Q980S9
E	-2	GLU	-	expression tag	UNP Q980S9

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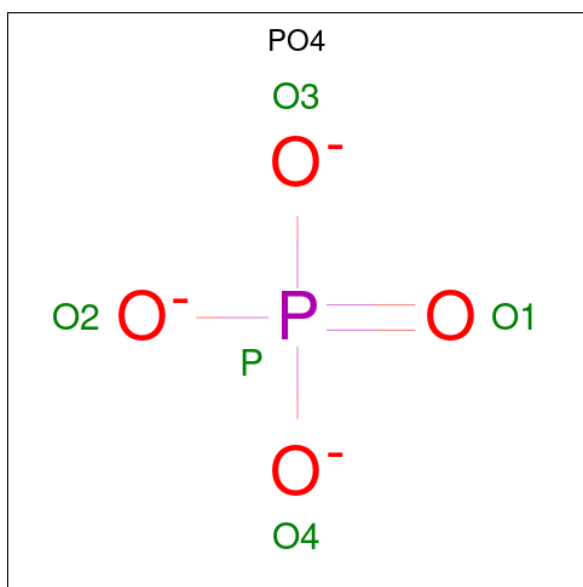
Chain	Residue	Modelled	Actual	Comment	Reference
E	-1	PRO	-	expression tag	UNP Q980S9
E	0	ILE	-	expression tag	UNP Q980S9
E	159	GLY	ASN	conflict	UNP Q980S9
F	-8	GLY	-	expression tag	UNP Q980S9
F	-7	SER	-	expression tag	UNP Q980S9
F	-6	HIS	-	expression tag	UNP Q980S9
F	-5	MET	-	expression tag	UNP Q980S9
F	-4	PHE	-	expression tag	UNP Q980S9
F	-3	GLU	-	expression tag	UNP Q980S9
F	-2	GLU	-	expression tag	UNP Q980S9
F	-1	PRO	-	expression tag	UNP Q980S9
F	0	ILE	-	expression tag	UNP Q980S9
F	159	GLY	ASN	conflict	UNP Q980S9
G	-8	GLY	-	expression tag	UNP Q980S9
G	-7	SER	-	expression tag	UNP Q980S9
G	-6	HIS	-	expression tag	UNP Q980S9
G	-5	MET	-	expression tag	UNP Q980S9
G	-4	PHE	-	expression tag	UNP Q980S9
G	-3	GLU	-	expression tag	UNP Q980S9
G	-2	GLU	-	expression tag	UNP Q980S9
G	-1	PRO	-	expression tag	UNP Q980S9
G	0	ILE	-	expression tag	UNP Q980S9
G	159	GLY	ASN	conflict	UNP Q980S9
H	-8	GLY	-	expression tag	UNP Q980S9
H	-7	SER	-	expression tag	UNP Q980S9
H	-6	HIS	-	expression tag	UNP Q980S9
H	-5	MET	-	expression tag	UNP Q980S9
H	-4	PHE	-	expression tag	UNP Q980S9
H	-3	GLU	-	expression tag	UNP Q980S9
H	-2	GLU	-	expression tag	UNP Q980S9
H	-1	PRO	-	expression tag	UNP Q980S9
H	0	ILE	-	expression tag	UNP Q980S9
H	159	GLY	ASN	conflict	UNP Q980S9
I	-8	GLY	-	expression tag	UNP Q980S9
I	-7	SER	-	expression tag	UNP Q980S9
I	-6	HIS	-	expression tag	UNP Q980S9
I	-5	MET	-	expression tag	UNP Q980S9
I	-4	PHE	-	expression tag	UNP Q980S9
I	-3	GLU	-	expression tag	UNP Q980S9
I	-2	GLU	-	expression tag	UNP Q980S9
I	-1	PRO	-	expression tag	UNP Q980S9
I	0	ILE	-	expression tag	UNP Q980S9

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Chain	Residue	Modelled	Actual	Comment	Reference
I	159	GLY	ASN	conflict	UNP Q980S9
J	-8	GLY	-	expression tag	UNP Q980S9
J	-7	SER	-	expression tag	UNP Q980S9
J	-6	HIS	-	expression tag	UNP Q980S9
J	-5	MET	-	expression tag	UNP Q980S9
J	-4	PHE	-	expression tag	UNP Q980S9
J	-3	GLU	-	expression tag	UNP Q980S9
J	-2	GLU	-	expression tag	UNP Q980S9
J	-1	PRO	-	expression tag	UNP Q980S9
J	0	ILE	-	expression tag	UNP Q980S9
J	159	GLY	ASN	conflict	UNP Q980S9
K	-8	GLY	-	expression tag	UNP Q980S9
K	-7	SER	-	expression tag	UNP Q980S9
K	-6	HIS	-	expression tag	UNP Q980S9
K	-5	MET	-	expression tag	UNP Q980S9
K	-4	PHE	-	expression tag	UNP Q980S9
K	-3	GLU	-	expression tag	UNP Q980S9
K	-2	GLU	-	expression tag	UNP Q980S9
K	-1	PRO	-	expression tag	UNP Q980S9
K	0	ILE	-	expression tag	UNP Q980S9
K	159	GLY	ASN	conflict	UNP Q980S9
L	-8	GLY	-	expression tag	UNP Q980S9
L	-7	SER	-	expression tag	UNP Q980S9
L	-6	HIS	-	expression tag	UNP Q980S9
L	-5	MET	-	expression tag	UNP Q980S9
L	-4	PHE	-	expression tag	UNP Q980S9
L	-3	GLU	-	expression tag	UNP Q980S9
L	-2	GLU	-	expression tag	UNP Q980S9
L	-1	PRO	-	expression tag	UNP Q980S9
L	0	ILE	-	expression tag	UNP Q980S9
L	159	GLY	ASN	conflict	UNP Q980S9

- Molecule 2 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	P	0	0
			5	4	1		
2	A	1	Total	O	P	0	0
			5	4	1		
2	B	1	Total	O	P	0	0
			5	4	1		
2	C	1	Total	O	P	0	0
			5	4	1		
2	D	1	Total	O	P	0	0
			5	4	1		
2	D	1	Total	O	P	0	0
			5	4	1		
2	E	1	Total	O	P	0	0
			5	4	1		
2	E	1	Total	O	P	0	0
			5	4	1		
2	F	1	Total	O	P	0	0
			5	4	1		
2	G	1	Total	O	P	0	0
			5	4	1		
2	G	1	Total	O	P	0	0
			5	4	1		
2	H	1	Total	O	P	0	0
			5	4	1		
2	I	1	Total	O	P	0	0
			5	4	1		
2	I	1	Total	O	P	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	J	1	Total	O	P	0	0
			5	4	1		
2	K	1	Total	O	P	0	0
			5	4	1		
2	K	1	Total	O	P	0	0
			5	4	1		
2	L	1	Total	O	P	0	0
			5	4	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	111	Total	O	0	0
			111	111		
3	B	96	Total	O	0	0
			96	96		
3	C	84	Total	O	0	0
			84	84		
3	D	82	Total	O	0	0
			82	82		
3	E	123	Total	O	0	0
			123	123		
3	F	110	Total	O	0	0
			110	110		
3	G	45	Total	O	0	0
			45	45		
3	H	74	Total	O	0	0
			74	74		
3	I	31	Total	O	0	0
			31	31		
3	J	43	Total	O	0	0
			43	43		
3	K	72	Total	O	0	0
			72	72		
3	L	71	Total	O	0	0
			71	71		

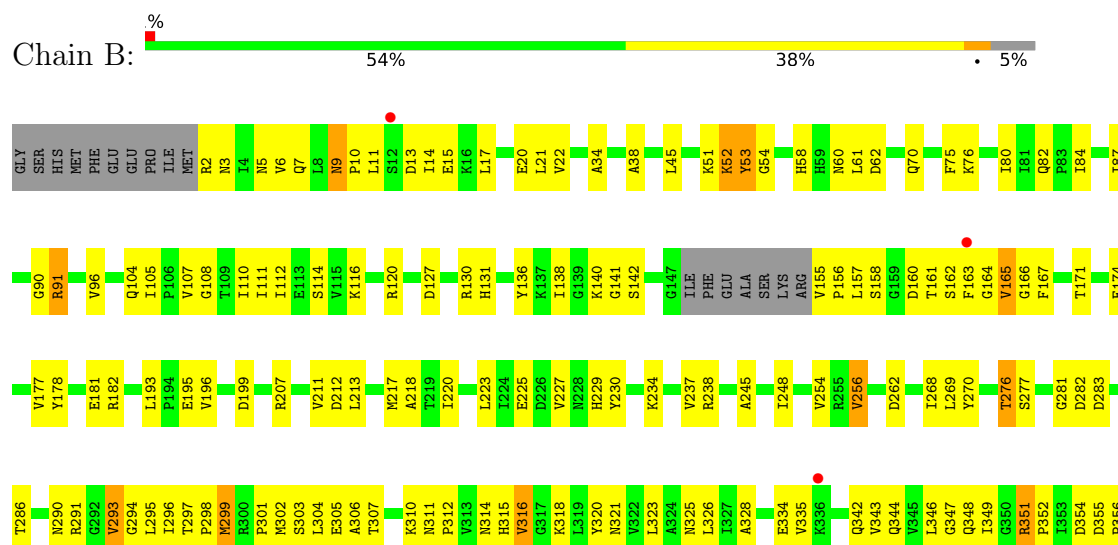
3 Residue-property plots

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: S-adenosylmethionine synthase



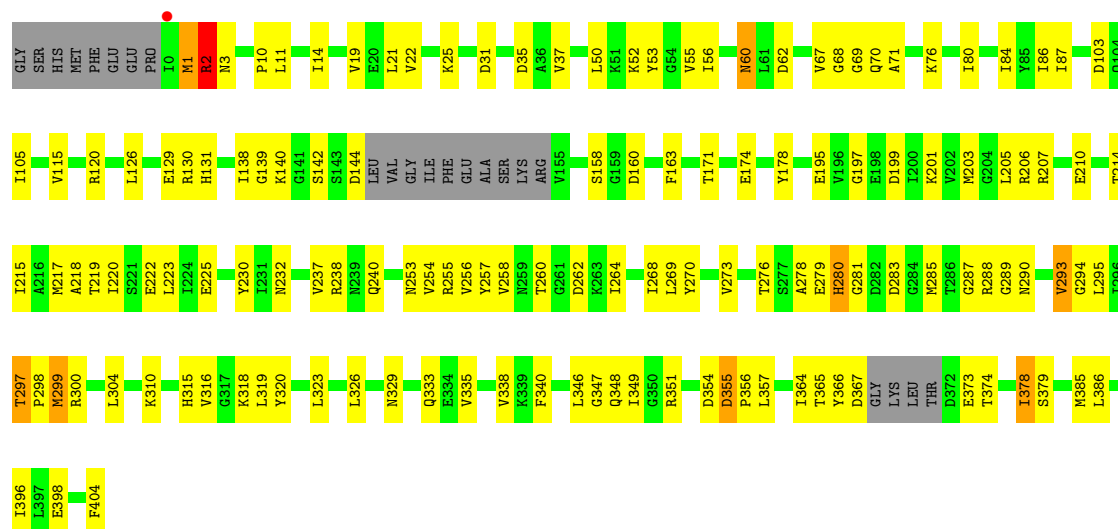
• Molecule 1: S-adenosylmethionine synthase





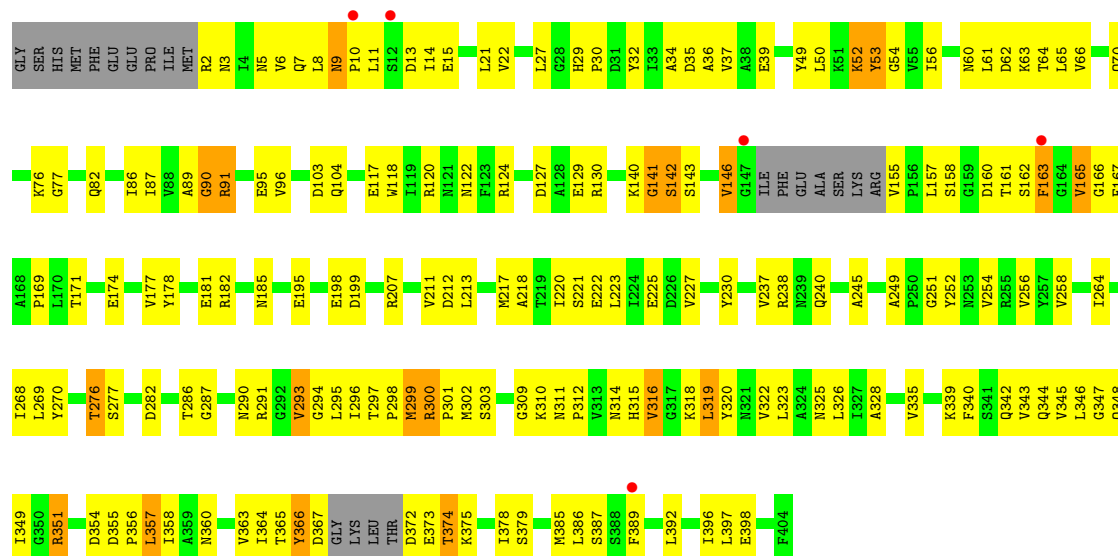
• Molecule 1: S-adenosylmethionine synthase

Chain C: 61% 31% 5%



• Molecule 1: S-adenosylmethionine synthase

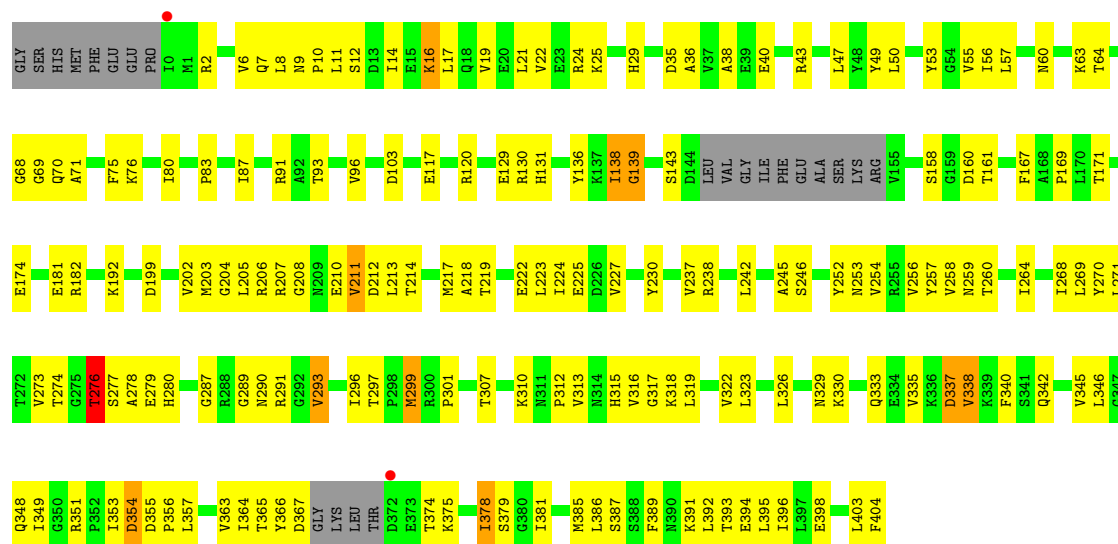
Chain D: 51% 39% 5% 5%



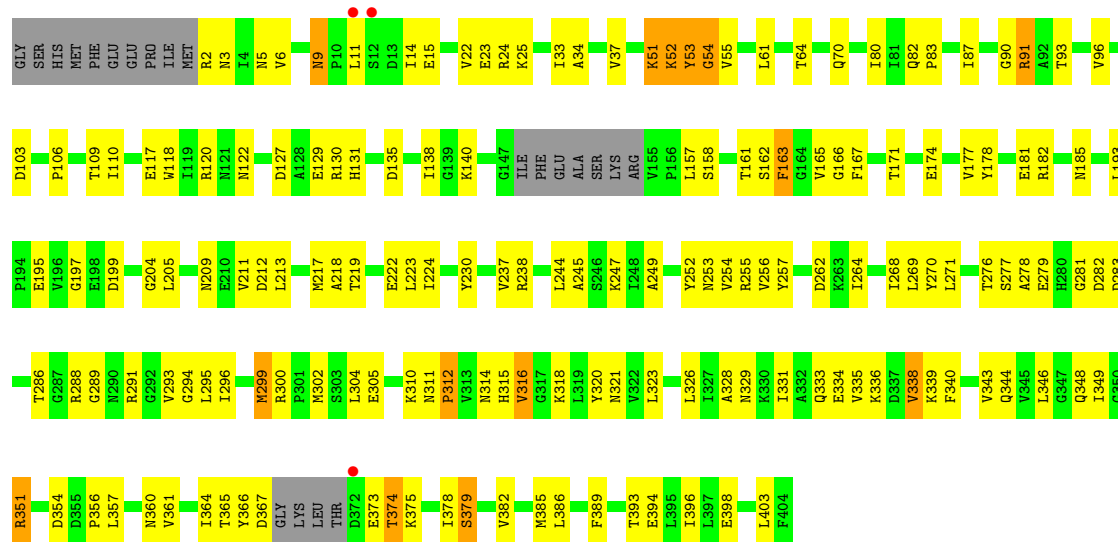
• Molecule 1: S-adenosylmethionine synthase

Chain E: 53% 39% 5%

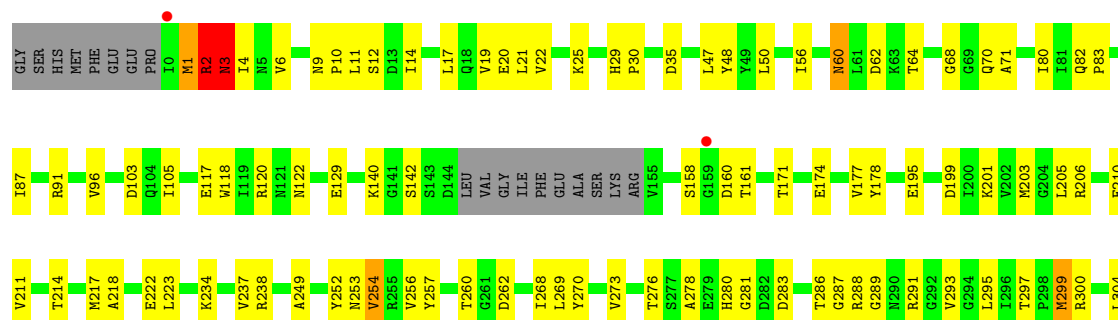


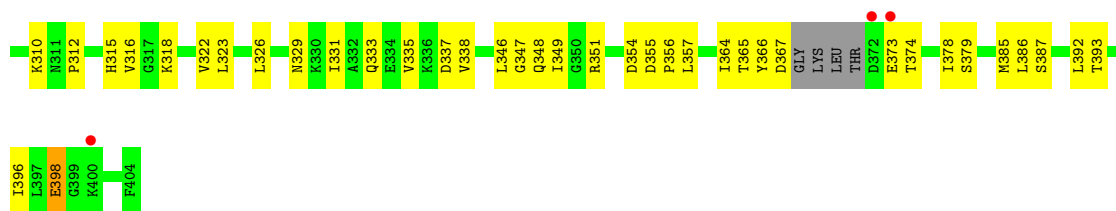


• Molecule 1: S-adenosylmethionine synthase

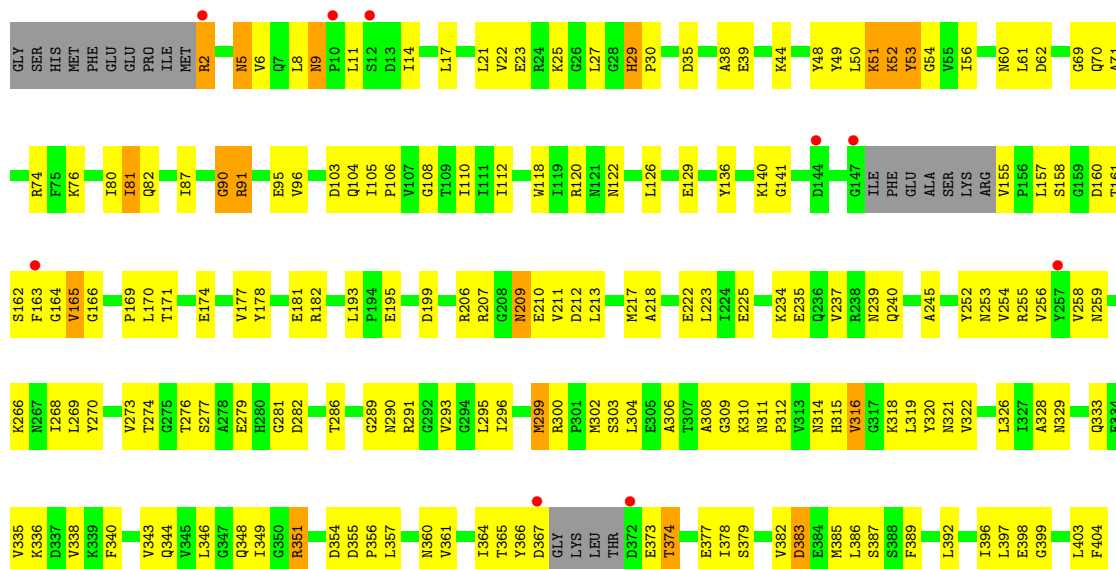


• Molecule 1: S-adenosylmethionine synthase

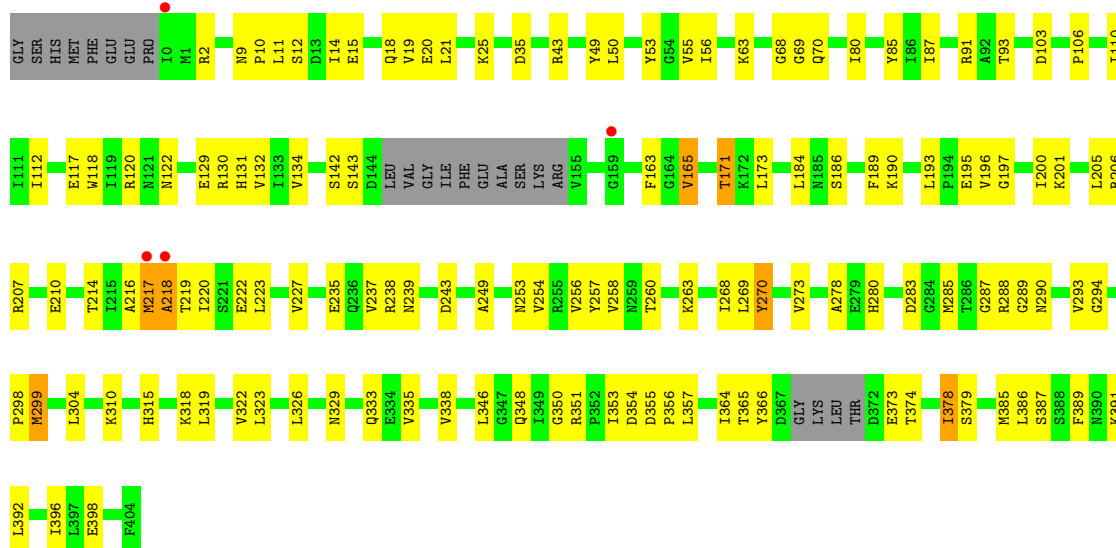




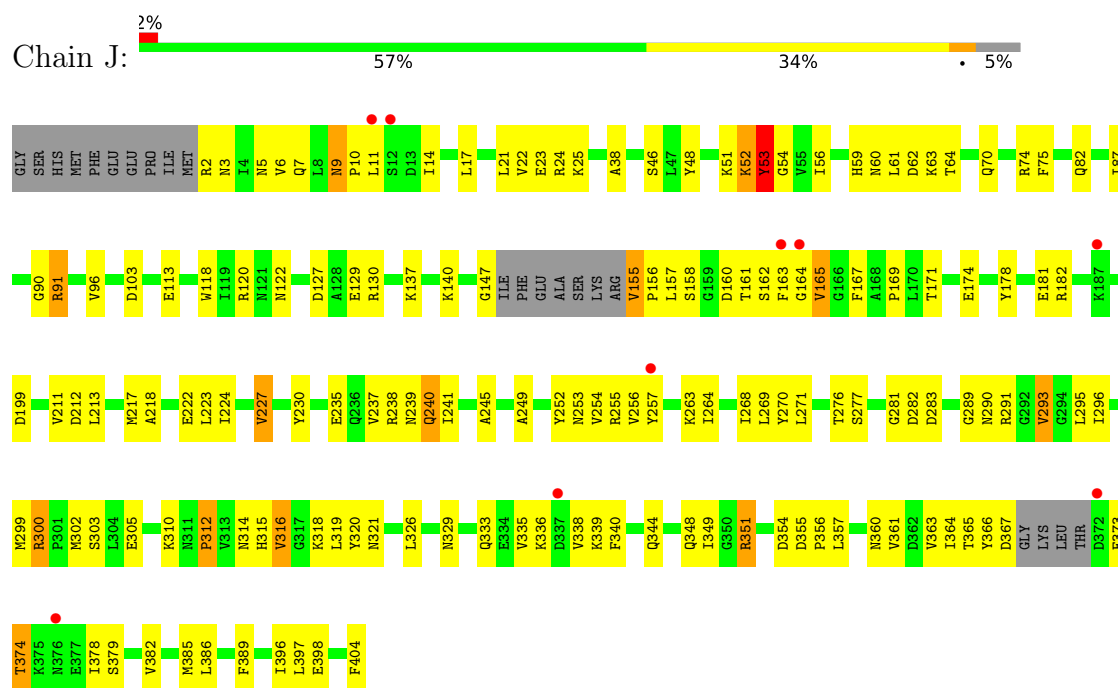
• Molecule 1: S-adenosylmethionine synthase



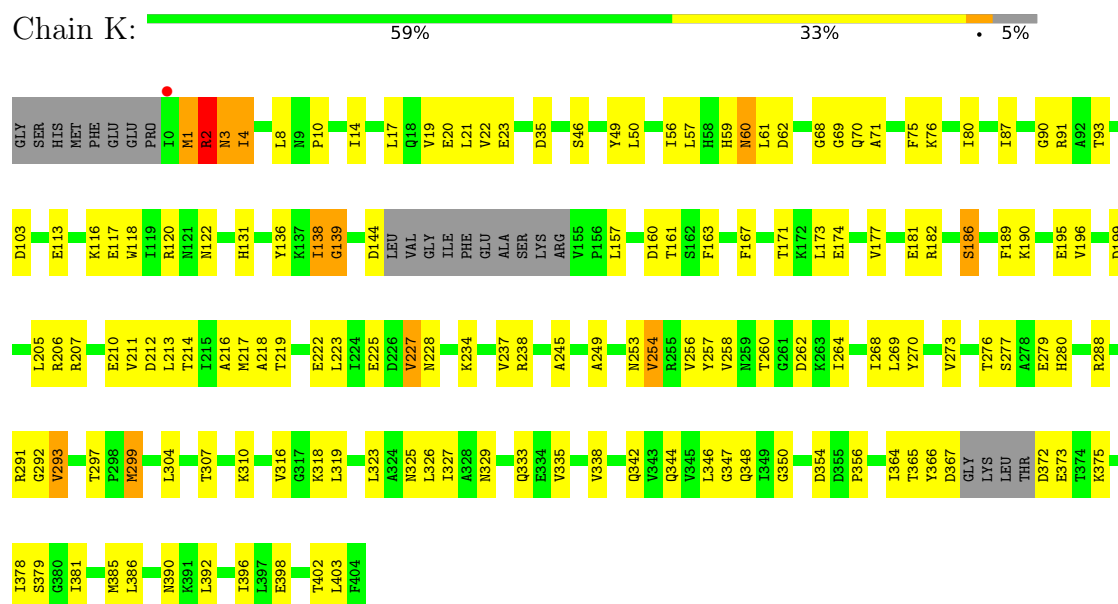
• Molecule 1: S-adenosylmethionine synthase



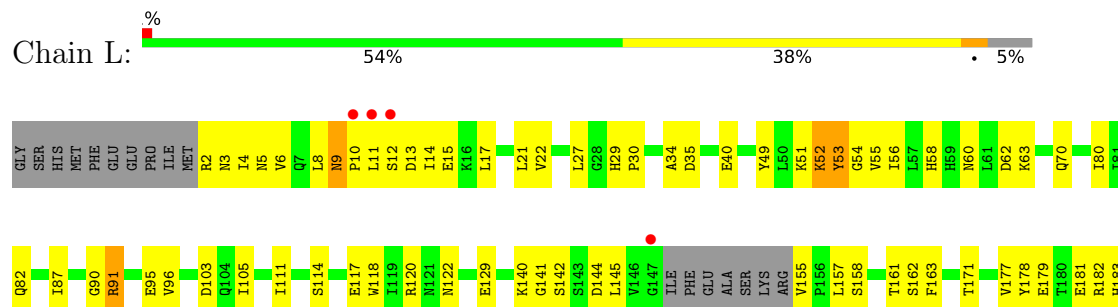
• Molecule 1: S-adenosylmethionine synthase



• Molecule 1: S-adenosylmethionine synthase



• Molecule 1: S-adenosylmethionine synthase





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	74.72Å 289.20Å 174.69Å 90.00° 99.85° 90.00°	Depositor
Resolution (Å)	30.00 – 2.80 30.00 – 2.80	Depositor EDS
% Data completeness (in resolution range)	75.6 (30.00-2.80) 67.0 (30.00-2.80)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	0.16	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.15 (at 2.80Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.221 , 0.256 0.227 , 0.259	Depositor DCC
R_{free} test set	1530 reflections (0.85%)	wwPDB-VP
Wilson B-factor (Å ²)	37.8	Xtriage
Anisotropy	1.226	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 47.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.031 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	37566	wwPDB-VP
Average B, all atoms (Å ²)	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.52% of the height of the origin peak. No significant pseudotranslation is detected.*

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

Bond lengths and bond angles in the following residue types are not validated in this section: PO4

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	1.07	0/3086	1.16	14/4172 (0.3%)
1	B	0.99	0/3089	1.19	14/4177 (0.3%)
1	C	0.98	0/3086	1.13	13/4172 (0.3%)
1	D	0.95	0/3089	1.21	25/4177 (0.6%)
1	E	1.15	2/3086 (0.1%)	1.19	13/4172 (0.3%)
1	F	1.02	0/3089	1.22	20/4177 (0.5%)
1	G	0.97	0/3086	1.13	12/4172 (0.3%)
1	H	1.10	1/3089 (0.0%)	1.26	27/4177 (0.6%)
1	I	0.95	4/3086 (0.1%)	1.14	12/4172 (0.3%)
1	J	0.96	0/3089	1.19	19/4177 (0.5%)
1	K	1.00	0/3086	1.16	12/4172 (0.3%)
1	L	0.97	1/3089 (0.0%)	1.18	17/4177 (0.4%)
All	All	1.01	8/37050 (0.0%)	1.18	198/50094 (0.4%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	C	0	1
1	G	0	1
1	K	0	1
All	All	0	3

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	I	218	ALA	CA-CB	-7.56	1.41	1.53
1	I	270	TYR	CA-C	7.02	1.61	1.53
1	H	2	ARG	CZ-NH1	-6.85	1.23	1.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	L	286	THR	CA-CB	-5.57	1.45	1.53
1	I	270	TYR	N-CA	-5.42	1.39	1.46
1	E	345	VAL	C-O	-5.42	1.18	1.24
1	I	220	ILE	CA-C	-5.06	1.48	1.53
1	E	276	THR	C-O	-5.04	1.18	1.23

All (198) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	I	218	ALA	N-CA-C	11.50	127.54	110.52
1	F	9	ASN	CA-C-N	9.53	130.70	120.12
1	F	9	ASN	C-N-CA	9.53	130.70	120.12
1	J	155	VAL	CA-C-N	8.51	128.45	120.03
1	J	155	VAL	C-N-CA	8.51	128.45	120.03
1	D	52	LYS	N-CA-C	-8.42	98.00	110.46
1	E	297	THR	CA-C-N	8.35	128.55	119.87
1	E	297	THR	C-N-CA	8.35	128.55	119.87
1	H	9	ASN	CA-C-N	8.11	129.13	120.12
1	H	9	ASN	C-N-CA	8.11	129.13	120.12
1	H	382	VAL	N-CA-C	-8.01	102.73	110.42
1	J	52	LYS	N-CA-C	-7.85	98.32	110.17
1	C	254	VAL	N-CA-C	7.62	118.85	107.80
1	F	52	LYS	N-CA-C	-7.57	98.35	109.63
1	A	297	THR	CA-C-N	7.53	127.20	119.82
1	A	297	THR	C-N-CA	7.53	127.20	119.82
1	H	2	ARG	NE-CZ-NH2	7.49	125.94	119.20
1	L	52	LYS	N-CA-C	-7.44	98.26	109.94
1	D	9	ASN	CA-C-N	7.41	128.35	120.12
1	D	9	ASN	C-N-CA	7.41	128.35	120.12
1	C	2	ARG	N-CA-C	7.36	122.26	113.28
1	K	2	ARG	N-CA-C	7.36	122.26	113.28
1	D	141	GLY	N-CA-C	-7.33	102.69	112.14
1	B	52	LYS	N-CA-C	-7.30	99.15	110.17
1	B	141	GLY	N-CA-C	-7.19	102.86	112.14
1	H	351	ARG	CA-C-N	-7.10	112.32	119.92
1	H	351	ARG	C-N-CA	-7.10	112.32	119.92
1	H	52	LYS	N-CA-C	-7.09	99.97	110.46
1	K	254	VAL	N-CA-C	7.05	118.02	107.80
1	I	270	TYR	N-CA-CB	-6.98	102.02	110.53
1	G	297	THR	CA-C-N	6.93	127.08	119.87
1	G	297	THR	C-N-CA	6.93	127.08	119.87
1	G	105	ILE	N-CA-C	6.91	115.51	107.77

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L	9	ASN	CA-C-N	6.86	127.73	120.12
1	L	9	ASN	C-N-CA	6.86	127.73	120.12
1	B	51	LYS	N-CA-C	-6.85	103.89	111.36
1	A	254	VAL	N-CA-C	6.85	117.73	107.80
1	E	139	GLY	N-CA-C	-6.84	99.51	111.62
1	B	312	PRO	CA-C-N	-6.79	116.23	123.08
1	B	312	PRO	C-N-CA	-6.79	116.23	123.08
1	H	209	ASN	N-CA-C	-6.79	104.13	112.88
1	A	354	ASP	N-CA-C	-6.76	104.61	112.92
1	H	90	GLY	N-CA-C	6.66	118.99	110.20
1	J	9	ASN	CA-C-N	6.64	127.49	120.12
1	J	9	ASN	C-N-CA	6.64	127.49	120.12
1	G	254	VAL	N-CA-C	6.63	117.41	107.80
1	H	155	VAL	CA-C-N	6.57	126.54	120.03
1	H	155	VAL	C-N-CA	6.57	126.54	120.03
1	D	146	VAL	N-CA-C	6.55	116.71	110.42
1	F	393	THR	N-CA-C	-6.51	104.11	111.07
1	K	139	GLY	N-CA-C	-6.46	100.18	111.62
1	G	2	ARG	N-CA-C	6.46	121.24	113.23
1	D	258	VAL	N-CA-C	6.42	117.12	107.80
1	K	297	THR	CA-C-N	6.41	126.10	119.82
1	K	297	THR	C-N-CA	6.41	126.10	119.82
1	E	138	ILE	CA-C-N	-6.41	115.35	121.57
1	E	138	ILE	C-N-CA	-6.41	115.35	121.57
1	J	312	PRO	CA-C-N	-6.36	116.66	123.08
1	J	312	PRO	C-N-CA	-6.36	116.66	123.08
1	F	51	LYS	N-CA-C	-6.33	104.46	111.36
1	L	312	PRO	CA-C-N	-6.31	116.70	123.08
1	L	312	PRO	C-N-CA	-6.31	116.70	123.08
1	D	300	ARG	CA-C-N	-6.31	113.45	120.14
1	D	300	ARG	C-N-CA	-6.31	113.45	120.14
1	H	105	ILE	N-CA-C	6.30	114.83	107.77
1	A	11	LEU	N-CA-C	-6.28	105.44	113.23
1	D	251	GLY	N-CA-C	-6.21	107.14	115.21
1	A	141	GLY	N-CA-C	6.18	120.35	110.90
1	L	351	ARG	N-CA-C	-6.16	102.31	110.07
1	F	336	LYS	N-CA-C	-6.14	105.63	112.57
1	B	351	ARG	CA-C-N	-6.11	113.39	119.92
1	B	351	ARG	C-N-CA	-6.11	113.39	119.92
1	A	281	GLY	N-CA-C	6.08	123.24	114.67
1	D	155	VAL	CA-C-N	6.06	126.03	120.03
1	D	155	VAL	C-N-CA	6.06	126.03	120.03

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	312	PRO	CA-C-N	-6.05	116.97	123.08
1	H	312	PRO	C-N-CA	-6.05	116.97	123.08
1	D	264	ILE	N-CA-C	6.04	116.78	110.62
1	F	264	ILE	N-CA-C	6.03	116.77	110.62
1	F	351	ARG	CA-C-N	-6.03	113.36	120.11
1	F	351	ARG	C-N-CA	-6.03	113.36	120.11
1	C	297	THR	CA-C-N	6.02	126.13	119.87
1	C	297	THR	C-N-CA	6.02	126.13	119.87
1	D	351	ARG	CA-C-N	-6.02	113.48	119.92
1	D	351	ARG	C-N-CA	-6.02	113.48	119.92
1	E	337	ASP	N-CA-C	-6.00	104.48	112.94
1	F	338	VAL	N-CA-C	6.00	118.08	108.85
1	J	264	ILE	N-CA-C	5.98	116.72	110.62
1	E	16	LYS	N-CA-C	-5.98	106.04	113.15
1	C	264	ILE	N-CA-C	5.95	116.69	110.62
1	K	249	ALA	CA-C-N	-5.92	113.86	119.85
1	K	249	ALA	C-N-CA	-5.92	113.86	119.85
1	H	29	HIS	CA-C-N	5.91	127.23	119.84
1	H	29	HIS	C-N-CA	5.91	127.23	119.84
1	A	280	HIS	CA-C-N	-5.91	113.29	122.86
1	A	280	HIS	C-N-CA	-5.91	113.29	122.86
1	F	379	SER	N-CA-C	-5.91	104.97	111.82
1	I	142	SER	N-CA-C	5.90	118.82	109.50
1	H	255	ARG	N-CA-C	-5.88	99.61	109.07
1	H	2	ARG	CD-NE-CZ	5.86	132.60	124.40
1	B	374	THR	N-CA-C	-5.85	105.47	113.30
1	D	319	LEU	N-CA-C	5.84	117.73	111.36
1	G	300	ARG	N-CA-C	-5.82	101.50	110.58
1	C	105	ILE	N-CA-C	5.80	114.27	107.77
1	J	336	LYS	N-CA-C	-5.80	106.02	112.57
1	G	398	GLU	N-CA-C	-5.80	104.87	112.41
1	B	334	GLU	N-CA-C	-5.77	106.88	114.04
1	D	90	GLY	N-CA-C	5.76	117.66	110.29
1	G	331	ILE	N-CA-C	-5.74	104.33	110.36
1	H	300	ARG	CA-C-N	-5.73	114.07	120.14
1	H	300	ARG	C-N-CA	-5.73	114.07	120.14
1	L	63	LYS	N-CA-C	5.73	116.61	108.54
1	F	312	PRO	CA-C-N	-5.72	117.09	123.16
1	F	312	PRO	C-N-CA	-5.72	117.09	123.16
1	F	257	TYR	CA-CB-CG	5.72	124.19	113.90
1	I	227	VAL	N-CA-C	-5.68	104.09	112.04
1	D	357	LEU	N-CA-C	-5.67	105.17	111.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	K	338	VAL	N-CA-C	5.65	117.55	108.85
1	G	249	ALA	CA-C-N	-5.64	114.15	119.85
1	G	249	ALA	C-N-CA	-5.64	114.15	119.85
1	G	142	SER	N-CA-C	5.63	118.74	109.72
1	H	319	LEU	N-CA-C	5.63	117.50	111.36
1	C	355	ASP	CA-C-N	5.62	125.93	119.92
1	C	355	ASP	C-N-CA	5.62	125.93	119.92
1	K	227	VAL	N-CA-C	-5.61	104.18	112.04
1	K	138	ILE	CA-C-N	-5.60	116.14	121.57
1	K	138	ILE	C-N-CA	-5.60	116.14	121.57
1	L	314	ASN	CA-C-N	-5.60	114.97	122.42
1	L	314	ASN	C-N-CA	-5.60	114.97	122.42
1	D	297	THR	CA-C-N	5.60	125.69	119.87
1	D	297	THR	C-N-CA	5.60	125.69	119.87
1	K	390	ASN	N-CA-C	-5.59	106.43	113.20
1	E	11	LEU	N-CA-C	-5.58	106.31	113.23
1	J	51	LYS	N-CA-C	-5.58	105.28	111.36
1	L	58	HIS	N-CA-C	5.56	118.39	110.10
1	J	53	TYR	N-CA-C	5.56	122.64	110.80
1	E	354	ASP	N-CA-C	-5.54	106.10	112.92
1	I	132	VAL	N-CA-C	5.52	116.07	108.12
1	E	143	SER	N-CA-C	5.51	117.22	108.79
1	F	279	GLU	N-CA-C	-5.47	105.86	112.54
1	J	163	PHE	N-CA-CB	-5.46	101.35	110.80
1	L	297	THR	CA-C-N	5.46	125.55	119.87
1	L	297	THR	C-N-CA	5.46	125.55	119.87
1	H	258	VAL	N-CA-C	5.45	115.71	107.80
1	J	240	GLN	N-CA-C	-5.45	105.42	111.36
1	H	336	LYS	N-CA-C	-5.41	106.45	112.57
1	I	254	VAL	N-CA-C	5.41	115.65	107.80
1	J	319	LEU	N-CA-C	5.38	118.30	111.69
1	A	249	ALA	CA-C-N	-5.37	114.42	119.85
1	A	249	ALA	C-N-CA	-5.37	114.42	119.85
1	B	9	ASN	CA-C-N	5.37	126.08	120.12
1	B	9	ASN	C-N-CA	5.37	126.08	120.12
1	E	211	VAL	N-CA-C	5.37	115.58	107.75
1	F	300	ARG	CA-C-N	-5.35	114.47	120.14
1	F	300	ARG	C-N-CA	-5.35	114.47	120.14
1	I	249	ALA	CA-C-N	-5.35	114.44	119.85
1	I	249	ALA	C-N-CA	-5.35	114.44	119.85
1	F	334	GLU	N-CA-C	-5.34	107.41	114.04
1	H	51	LYS	N-CA-C	-5.29	105.60	111.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	215	ILE	N-CA-C	5.29	115.51	108.11
1	L	345	VAL	N-CA-C	5.28	115.59	107.77
1	D	227	VAL	N-CA-C	-5.27	104.66	112.04
1	F	138	ILE	CA-C-N	5.25	125.65	121.61
1	F	138	ILE	C-N-CA	5.25	125.65	121.61
1	G	11	LEU	N-CA-C	-5.24	106.73	113.23
1	L	4	ILE	N-CA-C	5.19	116.06	108.53
1	B	256	VAL	N-CA-C	5.17	115.41	108.17
1	D	66	VAL	N-CA-C	5.17	115.21	107.51
1	D	163	PHE	N-CA-CB	-5.16	101.66	111.00
1	I	217	MET	CA-C-N	-5.15	111.71	122.07
1	I	217	MET	C-N-CA	-5.15	111.71	122.07
1	I	143	SER	N-CA-C	5.15	116.12	108.86
1	J	227	VAL	N-CA-C	-5.14	104.84	112.04
1	L	288	ARG	N-CA-C	-5.14	106.98	114.12
1	J	351	ARG	CA-C-N	-5.13	114.43	119.92
1	J	351	ARG	C-N-CA	-5.13	114.43	119.92
1	H	8	LEU	N-CA-C	-5.13	100.94	109.46
1	D	312	PRO	CA-C-N	-5.12	117.91	123.08
1	D	312	PRO	C-N-CA	-5.12	117.91	123.08
1	I	350	GLY	N-CA-C	-5.12	108.55	115.21
1	D	142	SER	N-CA-C	-5.12	103.28	110.50
1	C	300	ARG	N-CA-C	-5.12	102.60	110.58
1	A	26	GLY	N-CA-C	5.10	119.56	112.57
1	E	338	VAL	N-CA-C	5.09	115.82	108.48
1	A	227	VAL	N-CA-C	-5.09	104.92	112.04
1	A	139	GLY	N-CA-C	-5.07	102.23	111.15
1	H	266	LYS	N-CA-C	-5.07	106.94	113.02
1	L	105	ILE	N-CA-C	5.07	113.44	107.77
1	L	27	LEU	N-CA-C	5.07	119.73	111.37
1	C	280	HIS	CA-C-N	-5.04	114.69	122.86
1	C	280	HIS	C-N-CA	-5.04	114.69	122.86
1	E	202	VAL	N-CA-C	5.04	115.83	108.58
1	H	27	LEU	N-CA-C	5.03	119.66	111.37
1	B	297	THR	CA-C-N	5.01	125.08	119.87
1	B	297	THR	C-N-CA	5.01	125.08	119.87
1	J	300	ARG	CA-C-N	-5.01	114.83	120.14
1	J	300	ARG	C-N-CA	-5.01	114.83	120.14
1	C	142	SER	N-CA-C	5.00	116.98	109.23

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	1	MET	Peptide
1	G	1	MET	Peptide
1	K	1	MET	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	3043	0	3119	153	0
1	B	3046	0	3119	150	0
1	C	3043	0	3119	124	0
1	D	3046	0	3119	168	0
1	E	3043	0	3119	153	0
1	F	3046	0	3119	157	0
1	G	3043	0	3119	102	0
1	H	3046	0	3119	150	0
1	I	3043	0	3119	106	0
1	J	3046	0	3119	127	0
1	K	3043	0	3119	118	0
1	L	3046	0	3119	149	0
2	A	10	0	0	3	0
2	B	5	0	0	0	0
2	C	5	0	0	1	0
2	D	10	0	0	1	0
2	E	10	0	0	0	0
2	F	5	0	0	1	0
2	G	10	0	0	2	0
2	H	5	0	0	0	0
2	I	10	0	0	0	0
2	J	5	0	0	1	0
2	K	10	0	0	0	0
2	L	5	0	0	2	0
3	A	111	0	0	44	0
3	B	96	0	0	36	0
3	C	84	0	0	24	0
3	D	82	0	0	37	0
3	E	123	0	0	33	0
3	F	110	0	0	35	0
3	G	45	0	0	9	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	H	74	0	0	29	0
3	I	31	0	0	6	0
3	J	43	0	0	15	0
3	K	72	0	0	22	0
3	L	71	0	0	28	0
All	All	37566	0	37428	1542	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 21.

All (1542) 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:I:214:THR:HG21	1:J:161:THR:HG21	1.43	1.00
1:I:217:MET:HE1	1:I:237:VAL:HG11	1.46	0.97
1:B:310:LYS:HB2	1:B:318:LYS:HD2	1.45	0.96
1:G:214:THR:HG21	1:H:161:THR:HG21	1.48	0.95
1:C:217:MET:HE1	1:C:237:VAL:HG11	1.51	0.93
1:C:140:LYS:NZ	3:C:657:HOH:O	2.00	0.93
1:B:90:GLY:O	1:B:91:ARG:NH1	2.04	0.90
1:J:14:ILE:HD11	1:J:171:THR:HG21	1.55	0.89
1:K:318:LYS:HE2	1:K:396:ILE:HD13	1.55	0.88
1:K:71:ALA:H	1:K:280:HIS:HD2	1.21	0.87
1:A:214:THR:HG21	1:B:161:THR:HG21	1.56	0.87
1:D:90:GLY:O	1:D:91:ARG:NH1	2.07	0.87
1:A:26:GLY:HA2	3:A:640:HOH:O	1.74	0.87
1:J:90:GLY:O	1:J:91:ARG:NH1	2.08	0.86
1:L:158:SER:OG	1:L:348:GLN:O	1.94	0.84
1:L:90:GLY:O	1:L:91:ARG:NH1	2.11	0.84
1:F:14:ILE:HD11	1:F:171:THR:HG21	1.59	0.83
1:I:205:LEU:HD23	1:J:357:LEU:HD22	1.61	0.83
1:A:71:ALA:H	1:A:280:HIS:HD2	1.27	0.83
1:D:60:ASN:OD1	3:D:604:HOH:O	1.97	0.82
1:C:14:ILE:HD11	1:C:171:THR:HG21	1.61	0.82
1:K:189:PHE:HB3	3:K:670:HOH:O	1.78	0.82
1:K:144:ASP:OD1	3:K:640:HOH:O	1.97	0.81
1:E:68:GLY:O	1:F:91:ARG:NH2	2.12	0.81
1:C:201:LYS:NZ	1:D:160:ASP:OD2	2.14	0.81
1:H:245:ALA:HB1	1:H:254:VAL:HG21	1.62	0.81
1:F:158:SER:OG	1:F:348:GLN:O	1.98	0.80
1:A:364:ILE:O	3:A:650:HOH:O	1.98	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:211:VAL:HB	1:B:254:VAL:HG12	1.63	0.80
1:L:60:ASN:OD1	3:L:603:HOH:O	1.99	0.80
1:D:335:VAL:HG21	1:D:378:ILE:HG22	1.64	0.80
1:D:372:ASP:N	3:D:638:HOH:O	2.13	0.80
1:K:50:LEU:HD13	1:K:56:ILE:HG13	1.64	0.79
1:E:214:THR:HG21	1:F:161:THR:HG21	1.63	0.79
1:K:234:LYS:NZ	1:K:262:ASP:OD2	2.14	0.79
1:F:90:GLY:O	1:F:91:ARG:NH1	2.15	0.79
1:A:117:GLU:OE1	1:A:120:ARG:NH1	2.15	0.79
1:J:158:SER:OG	1:J:348:GLN:O	2.00	0.78
1:A:14:ILE:HD11	1:A:171:THR:HG21	1.64	0.78
1:L:52:LYS:O	1:L:54:GLY:N	2.17	0.78
1:I:197:GLY:N	1:I:218:ALA:O	2.14	0.78
1:B:335:VAL:HG21	1:B:378:ILE:HG22	1.66	0.78
1:F:217:MET:HE1	1:F:237:VAL:HG11	1.65	0.78
1:G:117:GLU:OE1	1:G:120:ARG:NH1	2.16	0.78
1:K:214:THR:HG21	1:L:161:THR:HG21	1.65	0.78
1:G:335:VAL:HG21	1:G:378:ILE:HG22	1.66	0.77
1:D:14:ILE:HD11	1:D:171:THR:HG21	1.65	0.77
1:H:326:LEU:HB3	1:H:385:MET:HE1	1.66	0.77
1:G:60:ASN:ND2	3:G:626:HOH:O	2.17	0.77
1:L:14:ILE:HD11	1:L:171:THR:HG21	1.65	0.77
1:J:217:MET:HE1	1:J:237:VAL:HG11	1.67	0.77
1:B:14:ILE:HD11	1:B:171:THR:HG21	1.65	0.77
1:I:273:VAL:O	3:I:615:HOH:O	2.03	0.77
1:K:205:LEU:HD23	1:L:357:LEU:HD22	1.66	0.77
1:K:398:GLU:N	1:K:398:GLU:OE1	2.18	0.77
1:L:213:LEU:HB3	1:L:256:VAL:HG12	1.67	0.77
1:J:326:LEU:HB3	1:J:385:MET:HE1	1.66	0.77
1:L:40:GLU:O	3:L:667:HOH:O	2.02	0.76
1:J:245:ALA:HB1	1:J:254:VAL:HG21	1.67	0.76
1:H:304:LEU:O	3:H:669:HOH:O	2.02	0.76
1:C:214:THR:HG21	1:D:161:THR:HG21	1.68	0.76
1:L:290:ASN:ND2	1:L:404:PHE:O	2.14	0.76
1:J:398:GLU:OE1	1:J:398:GLU:N	2.19	0.75
1:D:50:LEU:HD13	1:D:56:ILE:HG13	1.68	0.75
1:F:321:ASN:ND2	3:F:656:HOH:O	2.19	0.75
1:K:117:GLU:OE1	1:K:120:ARG:NH1	2.18	0.75
1:G:21:LEU:HD13	1:G:299:MET:HE1	1.69	0.75
1:K:228:ASN:HB3	3:K:606:HOH:O	1.87	0.75
1:H:213:LEU:HB3	1:H:256:VAL:HG12	1.67	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:71:ALA:H	1:C:280:HIS:HD2	1.32	0.75
1:H:30:PRO:HG3	3:H:664:HOH:O	1.86	0.75
1:H:178:TYR:HB2	1:H:295:LEU:HD21	1.68	0.75
1:J:53:TYR:CE1	1:J:96:VAL:HG21	2.22	0.75
1:J:60:ASN:ND2	3:J:621:HOH:O	2.19	0.74
1:G:217:MET:HE1	1:G:237:VAL:HG11	1.69	0.74
1:H:335:VAL:HG21	1:H:378:ILE:HG22	1.68	0.74
1:A:335:VAL:HG21	1:A:378:ILE:HG22	1.67	0.74
1:A:367:ASP:O	3:A:649:HOH:O	2.04	0.74
1:B:155:VAL:HG23	3:B:641:HOH:O	1.86	0.74
1:A:21:LEU:HD13	1:A:299:MET:HE1	1.68	0.74
1:H:310:LYS:HB2	1:H:318:LYS:HD2	1.69	0.74
1:E:335:VAL:HG21	1:E:378:ILE:HG22	1.70	0.74
1:E:299:MET:HE3	1:F:344:GLN:HE22	1.51	0.74
1:L:9:ASN:O	3:L:663:HOH:O	2.06	0.73
1:A:60:ASN:ND2	1:A:62:ASP:OD1	2.21	0.73
1:C:197:GLY:N	1:C:218:ALA:O	2.15	0.73
1:I:351:ARG:NE	1:I:355:ASP:O	2.20	0.73
1:L:335:VAL:HG21	1:L:378:ILE:HG22	1.70	0.73
1:C:318:LYS:HE2	1:C:396:ILE:HD13	1.69	0.73
1:G:14:ILE:HD11	1:G:171:THR:HG21	1.71	0.73
1:A:300:ARG:NH2	3:A:652:HOH:O	2.22	0.73
2:A:502:PO4:O1	3:A:702:HOH:O	2.06	0.72
1:D:118:TRP:O	1:D:122:ASN:ND2	2.20	0.72
1:L:13:ASP:O	3:L:615:HOH:O	2.07	0.72
1:J:52:LYS:O	1:J:54:GLY:N	2.23	0.72
1:K:71:ALA:H	1:K:280:HIS:CD2	2.07	0.72
1:D:310:LYS:HB2	1:D:318:LYS:HD2	1.71	0.72
1:L:157:LEU:O	1:L:314:ASN:HB2	1.90	0.71
1:L:391:LYS:O	3:L:639:HOH:O	2.07	0.71
1:F:52:LYS:O	1:F:54:GLY:N	2.23	0.71
1:I:270:TYR:HD1	3:J:633:HOH:O	1.72	0.71
1:A:217:MET:HE1	1:A:237:VAL:HG11	1.73	0.71
1:B:398:GLU:OE1	1:B:398:GLU:N	2.23	0.71
1:E:71:ALA:H	1:E:280:HIS:HD2	1.38	0.71
1:L:310:LYS:NZ	2:L:501:PO4:O3	2.23	0.71
1:K:61:LEU:N	3:K:662:HOH:O	2.22	0.71
1:E:398:GLU:OE1	1:E:398:GLU:N	2.23	0.71
1:D:143:SER:OG	3:D:669:HOH:O	2.09	0.71
1:G:71:ALA:H	1:G:280:HIS:HD2	1.39	0.71
1:K:113:GLU:OE1	3:K:661:HOH:O	2.09	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:372:ASP:N	1:L:375:LYS:HZ3	1.89	0.71
1:H:273:VAL:O	3:H:628:HOH:O	2.08	0.70
1:C:374:THR:HG21	3:C:632:HOH:O	1.91	0.70
1:K:14:ILE:HD11	1:K:171:THR:HG21	1.73	0.70
1:L:211:VAL:HB	1:L:254:VAL:HG12	1.72	0.70
1:A:144:ASP:OD2	3:A:638:HOH:O	2.10	0.70
1:E:287:GLY:O	3:E:650:HOH:O	2.08	0.70
1:H:14:ILE:HD11	1:H:171:THR:HG21	1.74	0.70
1:G:205:LEU:HD23	1:H:357:LEU:HD22	1.73	0.70
1:F:318:LYS:HE2	1:F:396:ILE:HD13	1.74	0.70
1:B:158:SER:OG	1:B:348:GLN:O	2.10	0.70
1:C:21:LEU:HD13	1:C:299:MET:HE1	1.72	0.70
1:G:337:ASP:O	1:G:365:THR:OG1	2.09	0.70
1:B:321:ASN:ND2	3:B:632:HOH:O	2.23	0.70
1:E:211:VAL:HB	1:E:254:VAL:HG12	1.73	0.70
1:A:398:GLU:N	1:A:398:GLU:OE1	2.24	0.69
1:E:10:PRO:HA	1:E:364:ILE:HD11	1.74	0.69
1:F:398:GLU:N	1:F:398:GLU:OE1	2.24	0.69
1:H:290:ASN:ND2	1:H:404:PHE:O	2.24	0.69
1:A:238:ARG:HB2	1:A:256:VAL:HG23	1.74	0.69
1:C:255:ARG:NH1	3:C:623:HOH:O	2.25	0.69
1:E:204:GLY:O	3:E:721:HOH:O	2.09	0.69
1:F:211:VAL:HB	1:F:254:VAL:HG12	1.74	0.69
1:I:49:TYR:OH	1:I:93:THR:O	2.09	0.69
1:J:178:TYR:HB2	1:J:295:LEU:HD21	1.73	0.69
1:F:135:ASP:OD1	3:F:700:HOH:O	2.09	0.69
1:J:310:LYS:HE2	1:J:315:HIS:CE1	2.27	0.69
1:G:365:THR:HG23	1:G:367:ASP:HB2	1.72	0.69
1:H:351:ARG:NE	1:H:355:ASP:O	2.26	0.69
1:F:348:GLN:N	3:F:687:HOH:O	2.24	0.69
1:A:60:ASN:C	1:A:60:ASN:HD22	2.00	0.69
1:B:107:VAL:HA	3:B:691:HOH:O	1.92	0.69
1:D:310:LYS:HE2	1:D:315:HIS:CE1	2.27	0.69
1:F:310:LYS:HB2	1:F:318:LYS:HD2	1.73	0.69
1:H:306:ALA:O	1:H:310:LYS:NZ	2.20	0.69
1:I:326:LEU:HB3	1:I:385:MET:HE1	1.74	0.69
1:A:346:LEU:HD11	1:B:21:LEU:HD11	1.74	0.69
1:F:24:ARG:HH21	1:F:182:ARG:HH11	1.40	0.69
1:H:310:LYS:HE2	1:H:315:HIS:CE1	2.28	0.68
1:H:60:ASN:ND2	3:H:603:HOH:O	2.02	0.68
1:J:257:TYR:OH	3:J:601:HOH:O	2.08	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:502:PO4:O3	3:A:664:HOH:O	2.09	0.68
1:H:211:VAL:HB	1:H:254:VAL:HG12	1.75	0.68
1:C:323:LEU:HD21	1:C:386:LEU:HD11	1.74	0.68
1:G:118:TRP:O	1:G:122:ASN:ND2	2.24	0.68
1:F:385:MET:HB2	3:F:705:HOH:O	1.92	0.68
1:F:351:ARG:HD3	3:F:687:HOH:O	1.93	0.68
1:K:325:ASN:HB3	3:K:655:HOH:O	1.93	0.68
1:B:293:VAL:O	3:B:695:HOH:O	2.12	0.68
1:L:217:MET:HE1	1:L:237:VAL:HG11	1.76	0.68
1:L:326:LEU:HB3	1:L:385:MET:HE1	1.75	0.68
1:C:35:ASP:OD1	1:C:287:GLY:N	2.25	0.67
1:D:344:GLN:O	3:D:662:HOH:O	2.12	0.67
1:E:14:ILE:HD11	1:E:171:THR:HG21	1.76	0.67
1:F:245:ALA:HB1	1:F:254:VAL:HG21	1.76	0.67
1:F:293:VAL:O	3:F:701:HOH:O	2.13	0.67
1:L:398:GLU:OE1	1:L:398:GLU:N	2.28	0.67
1:K:10:PRO:HA	1:K:364:ILE:HD11	1.75	0.67
1:D:217:MET:HE1	1:D:237:VAL:HG11	1.76	0.67
1:E:91:ARG:HH21	1:F:281:GLY:C	2.02	0.67
1:E:318:LYS:HE2	1:E:396:ILE:HD13	1.75	0.67
1:H:398:GLU:N	1:H:398:GLU:OE1	2.26	0.67
1:C:144:ASP:OD2	3:C:645:HOH:O	2.13	0.67
1:K:35:ASP:OD2	3:K:629:HOH:O	2.11	0.67
1:E:224:ILE:HG13	3:E:712:HOH:O	1.93	0.67
1:J:113:GLU:OE1	3:J:626:HOH:O	2.12	0.67
1:A:71:ALA:H	1:A:280:HIS:CD2	2.10	0.67
1:I:398:GLU:N	1:I:398:GLU:OE1	2.27	0.67
1:K:217:MET:HE1	1:K:237:VAL:HG11	1.75	0.67
1:L:311:ASN:HB3	1:L:315:HIS:HB2	1.77	0.67
1:C:207:ARG:HE	1:D:5:ASN:HD22	1.42	0.67
2:C:501:PO4:O2	3:C:637:HOH:O	2.11	0.67
1:C:335:VAL:HG21	1:C:378:ILE:HG22	1.77	0.67
1:E:312:PRO:O	1:E:393:THR:OG1	2.06	0.67
1:E:217:MET:HE1	1:E:237:VAL:HG11	1.76	0.67
1:F:311:ASN:ND2	1:F:314:ASN:OD1	2.28	0.67
1:G:398:GLU:OE1	1:G:398:GLU:N	2.28	0.67
1:K:372:ASP:N	1:K:375:LYS:HZ2	1.93	0.66
1:C:398:GLU:N	1:C:398:GLU:OE1	2.28	0.66
1:E:219:THR:OG1	1:E:270:TYR:O	2.12	0.66
1:F:335:VAL:HG21	1:F:378:ILE:HG22	1.78	0.66
1:C:71:ALA:H	1:C:280:HIS:CD2	2.14	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:68:GLY:O	1:L:91:ARG:NH2	2.28	0.66
1:L:245:ALA:HB1	1:L:254:VAL:HG21	1.76	0.66
1:D:53:TYR:CE1	1:D:96:VAL:HG21	2.30	0.66
1:K:323:LEU:HB2	1:K:392:LEU:HD21	1.76	0.66
1:F:310:LYS:HE2	1:F:315:HIS:CE1	2.31	0.66
1:J:213:LEU:HB3	1:J:256:VAL:HG12	1.78	0.66
1:K:372:ASP:N	3:K:648:HOH:O	2.29	0.66
1:B:13:ASP:O	3:B:616:HOH:O	2.13	0.66
1:F:326:LEU:HB3	1:F:385:MET:HE1	1.77	0.66
1:I:263:LYS:HB2	1:I:268:ILE:HD11	1.78	0.66
1:J:162:SER:OG	1:J:316:VAL:HG12	1.96	0.66
1:F:311:ASN:O	3:F:663:HOH:O	2.14	0.65
1:I:50:LEU:HD13	1:I:56:ILE:HG13	1.78	0.65
1:I:120:ARG:NH2	1:I:129:GLU:OE2	2.29	0.65
1:L:35:ASP:OD2	3:L:625:HOH:O	2.14	0.65
1:L:310:LYS:HB2	1:L:318:LYS:HD2	1.78	0.65
1:K:207:ARG:HE	1:L:5:ASN:HD22	1.45	0.65
1:B:385:MET:HB2	3:B:640:HOH:O	1.95	0.65
1:C:367:ASP:O	3:C:613:HOH:O	2.14	0.65
1:D:22:VAL:HG11	1:D:174:GLU:HA	1.77	0.65
1:G:17:LEU:O	1:G:206:ARG:NH2	2.22	0.65
1:C:223:LEU:HB2	3:C:646:HOH:O	1.95	0.65
1:K:373:GLU:N	3:K:648:HOH:O	2.30	0.65
1:E:24:ARG:NH1	3:E:681:HOH:O	2.29	0.65
1:C:293:VAL:O	3:C:676:HOH:O	2.15	0.65
1:K:120:ARG:NH2	3:K:663:HOH:O	2.30	0.65
1:B:352:PRO:HB3	3:B:641:HOH:O	1.97	0.65
1:G:2:ARG:O	1:G:3:ASN:HB2	1.95	0.65
1:B:245:ALA:HB1	1:B:254:VAL:HG21	1.79	0.64
1:J:161:THR:HG22	1:J:349:ILE:H	1.61	0.64
1:B:382:VAL:O	3:B:640:HOH:O	2.13	0.64
1:B:326:LEU:HB3	1:B:385:MET:HE1	1.80	0.64
1:A:178:TYR:HB2	1:A:295:LEU:HD21	1.79	0.64
1:E:21:LEU:HD13	1:E:299:MET:HE1	1.78	0.64
1:F:70:GLN:HB2	1:F:82:GLN:HB3	1.79	0.64
1:I:2:ARG:HH12	1:I:387:SER:HB3	1.62	0.64
1:L:35:ASP:OD1	1:L:287:GLY:N	2.30	0.64
1:E:120:ARG:NH2	1:E:129:GLU:OE2	2.31	0.64
1:G:273:VAL:O	3:G:615:HOH:O	2.15	0.64
1:F:213:LEU:HB3	1:F:256:VAL:HG12	1.78	0.64
1:G:2:ARG:HH12	1:G:387:SER:HB3	1.62	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:G:501:PO4:O3	3:G:635:HOH:O	2.15	0.64
1:G:318:LYS:HE2	1:G:396:ILE:HD13	1.78	0.64
1:G:29:HIS:NE2	2:G:501:PO4:O1	2.30	0.63
1:H:378:ILE:HG13	1:H:379:SER:N	2.13	0.63
1:J:253:ASN:OD1	1:J:255:ARG:NH2	2.31	0.63
1:A:68:GLY:O	1:B:91:ARG:NH2	2.30	0.63
1:D:10:PRO:HA	1:D:364:ILE:HD11	1.80	0.63
1:L:305:GLU:HB3	3:L:666:HOH:O	1.98	0.63
1:A:207:ARG:HE	1:B:5:ASN:HD22	1.46	0.63
1:D:161:THR:HG22	1:D:349:ILE:H	1.63	0.63
1:C:163:PHE:O	3:C:668:HOH:O	2.15	0.63
1:F:277:SER:O	1:F:282:ASP:HB2	1.98	0.63
1:H:53:TYR:CE1	1:H:96:VAL:HG21	2.34	0.63
1:H:162:SER:OG	1:H:316:VAL:HG12	1.99	0.63
1:A:253:ASN:HB2	3:A:671:HOH:O	1.99	0.63
1:C:80:ILE:HG13	1:C:131:HIS:CD2	2.34	0.63
1:C:69:GLY:O	1:C:70:GLN:HG3	1.99	0.63
1:E:205:LEU:HD23	1:F:357:LEU:HD22	1.79	0.63
1:E:329:ASN:O	1:E:333:GLN:HG3	2.00	0.62
1:F:161:THR:HG22	1:F:349:ILE:H	1.63	0.62
1:K:327:ILE:HG12	3:K:666:HOH:O	1.99	0.62
1:D:86:ILE:HG23	3:D:651:HOH:O	2.00	0.62
1:D:245:ALA:HB1	1:D:254:VAL:HG21	1.81	0.62
1:K:19:VAL:HG21	1:L:360:ASN:ND2	2.15	0.62
1:G:288:ARG:NH2	1:H:62:ASP:O	2.32	0.62
1:D:35:ASP:OD1	1:D:287:GLY:N	2.26	0.62
1:J:277:SER:O	1:J:282:ASP:HB2	2.00	0.62
1:A:205:LEU:HD23	1:B:357:LEU:HD22	1.82	0.62
1:E:378:ILE:HG13	1:E:379:SER:N	2.15	0.62
1:B:157:LEU:O	1:B:314:ASN:HB2	2.00	0.62
1:D:351:ARG:NE	1:D:355:ASP:O	2.30	0.62
1:H:51:LYS:NZ	3:H:649:HOH:O	2.31	0.62
1:A:238:ARG:HB2	1:A:256:VAL:CG2	2.30	0.62
1:C:378:ILE:HG13	1:C:379:SER:N	2.15	0.62
1:D:222:GLU:HB2	1:D:223:LEU:HD12	1.81	0.62
1:F:118:TRP:O	1:F:122:ASN:ND2	2.27	0.62
1:J:60:ASN:ND2	3:J:639:HOH:O	2.32	0.62
1:K:186:SER:N	3:K:670:HOH:O	2.33	0.62
1:H:274:THR:O	3:H:616:HOH:O	2.16	0.62
1:K:22:VAL:HG11	1:K:174:GLU:HA	1.81	0.62
1:D:365:THR:HG23	1:D:367:ASP:HB2	1.82	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:155:VAL:N	3:J:634:HOH:O	2.31	0.61
1:K:253:ASN:ND2	3:K:601:HOH:O	2.21	0.61
1:L:178:TYR:HB2	1:L:295:LEU:HD21	1.82	0.61
1:L:373:GLU:HG2	1:L:374:THR:H	1.64	0.61
1:F:346:LEU:HG	3:F:651:HOH:O	2.00	0.61
1:J:63:LYS:NZ	3:J:643:HOH:O	2.33	0.61
1:E:91:ARG:N	3:E:706:HOH:O	2.32	0.61
1:K:335:VAL:HG21	1:K:378:ILE:HG22	1.81	0.61
1:I:288:ARG:NH2	1:J:62:ASP:O	2.32	0.61
1:J:22:VAL:HG11	1:J:174:GLU:HA	1.83	0.61
1:K:21:LEU:HD13	1:K:299:MET:HE1	1.81	0.61
1:G:373:GLU:HG2	1:G:374:THR:H	1.66	0.61
1:L:277:SER:O	1:L:282:ASP:HB2	2.01	0.61
1:D:311:ASN:HB3	1:D:315:HIS:HB2	1.83	0.61
1:E:199:ASP:HB2	1:E:276:THR:HA	1.81	0.61
1:L:40:GLU:HG3	3:L:667:HOH:O	2.01	0.61
1:E:207:ARG:HE	1:F:5:ASN:HD22	1.47	0.61
1:K:2:ARG:O	1:K:3:ASN:HB2	1.99	0.61
1:A:19:VAL:HG21	1:B:360:ASN:ND2	2.15	0.61
1:E:69:GLY:O	1:E:70:GLN:HG3	2.00	0.61
1:F:311:ASN:HB3	1:F:315:HIS:HB2	1.82	0.61
1:L:354:ASP:HA	1:L:389:PHE:HE2	1.65	0.61
1:B:310:LYS:HE2	1:B:315:HIS:CE1	2.36	0.60
1:D:398:GLU:N	1:D:398:GLU:OE1	2.34	0.60
1:D:211:VAL:HB	1:D:254:VAL:HG12	1.84	0.60
1:L:300:ARG:NH1	3:L:636:HOH:O	2.34	0.60
1:H:217:MET:HE1	1:H:237:VAL:HG11	1.84	0.60
1:F:230:TYR:OH	1:F:262:ASP:OD2	2.09	0.60
1:J:238:ARG:HB2	1:J:256:VAL:HG23	1.83	0.60
1:K:181:GLU:OE1	1:K:182:ARG:HD2	2.02	0.60
1:A:120:ARG:NH2	1:A:129:GLU:OE2	2.33	0.60
1:L:118:TRP:O	1:L:122:ASN:ND2	2.33	0.60
1:B:110:ILE:HB	3:B:691:HOH:O	2.01	0.60
1:G:91:ARG:HH21	1:H:281:GLY:C	2.10	0.60
1:K:23:GLU:OE2	3:K:660:HOH:O	2.17	0.60
1:A:273:VAL:O	3:A:683:HOH:O	2.17	0.60
1:F:224:ILE:HG21	1:F:271:LEU:HD21	1.83	0.60
1:H:6:VAL:HG11	1:H:378:ILE:HD11	1.84	0.60
1:L:310:LYS:HE2	1:L:315:HIS:CE1	2.37	0.60
1:B:76:LYS:HG3	1:B:225:GLU:HA	1.84	0.60
1:I:348:GLN:NE2	1:J:212:ASP:OD2	2.34	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:19:VAL:HG21	1:L:360:ASN:HD21	1.67	0.60
1:C:232:ASN:HB2	3:C:604:HOH:O	2.02	0.60
1:E:57:LEU:HD22	1:E:93:THR:HG21	1.83	0.60
1:F:339:LYS:HB3	1:F:364:ILE:HG22	1.83	0.60
1:B:70:GLN:HB2	1:B:82:GLN:HB3	1.84	0.59
1:J:378:ILE:HG13	1:J:379:SER:N	2.14	0.59
1:B:277:SER:O	1:B:282:ASP:HB2	2.02	0.59
1:C:50:LEU:HD13	1:C:56:ILE:HG13	1.83	0.59
1:F:109:THR:N	3:F:704:HOH:O	2.34	0.59
1:A:402:THR:HG22	3:A:645:HOH:O	2.02	0.59
1:E:71:ALA:H	1:E:280:HIS:CD2	2.18	0.59
1:F:204:GLY:O	3:F:703:HOH:O	2.17	0.59
1:K:323:LEU:HD21	1:K:386:LEU:HD11	1.82	0.59
1:C:288:ARG:HG3	3:C:671:HOH:O	2.01	0.59
1:D:2:ARG:N	3:D:612:HOH:O	2.35	0.59
1:K:199:ASP:HB2	1:K:276:THR:HA	1.84	0.59
1:F:378:ILE:HG13	1:F:379:SER:N	2.16	0.59
1:G:234:LYS:NZ	1:G:262:ASP:OD2	2.33	0.59
1:H:9:ASN:OD1	1:H:11:LEU:HD13	2.03	0.59
1:L:53:TYR:CE1	1:L:96:VAL:HG21	2.38	0.59
1:B:178:TYR:HB2	1:B:295:LEU:HD21	1.84	0.59
1:E:218:ALA:HA	1:E:270:TYR:CD2	2.38	0.59
1:A:299:MET:HE3	1:B:344:GLN:OE1	2.03	0.59
1:B:217:MET:HE1	1:B:237:VAL:HG11	1.84	0.59
1:F:53:TYR:CE1	1:F:96:VAL:HG21	2.37	0.59
1:A:195:GLU:N	1:A:195:GLU:OE1	2.34	0.59
1:B:112:ILE:HG21	3:C:638:HOH:O	2.01	0.59
1:D:160:ASP:HB3	3:D:640:HOH:O	2.02	0.59
1:H:218:ALA:HA	1:H:270:TYR:HD2	1.68	0.59
1:I:117:GLU:OE1	1:I:120:ARG:NH1	2.34	0.59
1:B:218:ALA:HA	1:B:270:TYR:HD2	1.66	0.59
1:C:19:VAL:HG21	1:D:360:ASN:ND2	2.18	0.59
1:D:277:SER:O	1:D:282:ASP:HB2	2.03	0.59
1:E:117:GLU:OE1	1:E:120:ARG:NH1	2.35	0.59
1:F:338:VAL:HB	3:F:695:HOH:O	2.02	0.59
1:G:348:GLN:NE2	1:H:212:ASP:OD2	2.36	0.59
1:I:354:ASP:HA	1:I:389:PHE:HE2	1.66	0.59
1:A:349:ILE:HD12	3:A:706:HOH:O	2.01	0.58
1:B:21:LEU:HD13	1:B:299:MET:HE1	1.84	0.58
1:C:220:ILE:HB	3:C:646:HOH:O	2.03	0.58
1:I:207:ARG:HE	1:J:5:ASN:HD22	1.50	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:311:ASN:ND2	1:L:314:ASN:OD1	2.36	0.58
1:C:218:ALA:HA	1:C:270:TYR:CE2	2.38	0.58
1:E:9:ASN:O	1:E:12:SER:HB3	2.04	0.58
1:E:398:GLU:HB2	3:E:715:HOH:O	2.04	0.58
1:I:35:ASP:OD1	1:I:287:GLY:N	2.30	0.58
1:I:318:LYS:HE2	1:I:396:ILE:HD13	1.84	0.58
1:B:22:VAL:HG11	1:B:174:GLU:HA	1.84	0.58
1:G:326:LEU:HB3	1:G:385:MET:HE1	1.85	0.58
1:A:24:ARG:NE	3:A:640:HOH:O	2.36	0.58
1:A:52:LYS:NZ	3:A:698:HOH:O	2.36	0.58
1:B:52:LYS:O	1:B:54:GLY:N	2.35	0.58
1:D:178:TYR:HB2	1:D:295:LEU:HD21	1.84	0.58
1:G:50:LEU:HD13	1:G:56:ILE:HG13	1.84	0.58
1:H:308:ALA:HA	3:H:631:HOH:O	2.02	0.58
1:H:237:VAL:O	1:H:240:GLN:HB2	2.04	0.58
1:K:211:VAL:HB	1:K:254:VAL:HG12	1.84	0.58
1:K:238:ARG:HB2	1:K:256:VAL:CG2	2.34	0.58
1:A:199:ASP:HB2	1:A:276:THR:HA	1.85	0.58
1:B:365:THR:C	1:B:367:ASP:H	2.12	0.58
1:D:117:GLU:OE1	1:D:120:ARG:NH1	2.36	0.58
1:A:228:ASN:CG	1:E:264:ILE:HG21	2.29	0.58
1:E:40:GLU:HB2	3:E:718:HOH:O	2.03	0.58
1:J:373:GLU:HG2	1:J:374:THR:H	1.69	0.57
1:A:118:TRP:O	1:A:122:ASN:ND2	2.29	0.57
1:A:196:VAL:HB	3:A:701:HOH:O	2.03	0.57
1:A:222:GLU:HB2	1:A:223:LEU:HD12	1.86	0.57
1:A:268:ILE:C	1:A:269:LEU:HD12	2.29	0.57
1:C:158:SER:O	1:C:349:ILE:HA	2.04	0.57
1:C:288:ARG:NH2	1:D:62:ASP:O	2.36	0.57
1:H:5:ASN:C	1:H:5:ASN:HD22	2.12	0.57
1:A:320:TYR:OH	3:A:679:HOH:O	2.18	0.57
1:E:386:LEU:HA	3:E:677:HOH:O	2.05	0.57
1:J:312:PRO:HB2	3:J:628:HOH:O	2.04	0.57
1:A:14:ILE:N	3:A:648:HOH:O	2.26	0.57
1:A:243:ASP:HB3	3:A:668:HOH:O	2.05	0.57
1:B:294:GLY:HA3	3:B:695:HOH:O	2.03	0.57
1:B:400:LYS:HE3	3:B:620:HOH:O	2.05	0.57
1:A:315:HIS:HB3	3:A:695:HOH:O	2.03	0.57
1:B:161:THR:HG22	1:B:349:ILE:H	1.69	0.57
1:F:22:VAL:HG11	1:F:174:GLU:HA	1.86	0.57
1:I:218:ALA:HA	1:I:270:TYR:HE2	1.67	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:218:ALA:HA	1:L:270:TYR:HD2	1.70	0.57
1:A:9:ASN:O	1:A:12:SER:HB3	2.03	0.57
1:B:344:GLN:HA	3:B:650:HOH:O	2.02	0.57
1:D:36:ALA:HB3	3:D:646:HOH:O	2.04	0.57
1:J:238:ARG:HB2	1:J:256:VAL:CG2	2.34	0.57
1:J:181:GLU:OE1	1:J:182:ARG:HD2	2.05	0.57
1:D:37:VAL:HG13	3:D:646:HOH:O	2.05	0.56
1:I:19:VAL:HG21	1:J:360:ASN:ND2	2.19	0.56
1:B:108:GLY:O	1:B:112:ILE:HG12	2.04	0.56
1:F:162:SER:OG	1:F:316:VAL:HG12	2.05	0.56
1:D:157:LEU:O	1:D:314:ASN:HB2	2.05	0.56
1:E:22:VAL:HG11	1:E:174:GLU:HA	1.86	0.56
1:F:163:PHE:CD1	1:F:163:PHE:C	2.83	0.56
1:F:222:GLU:HG3	3:F:635:HOH:O	2.04	0.56
1:F:331:ILE:HG23	3:F:666:HOH:O	2.05	0.56
1:G:323:LEU:HD12	1:G:385:MET:HE2	1.88	0.56
1:I:25:LYS:NZ	2:J:501:PO4:O2	2.38	0.56
1:L:95:GLU:HB2	1:L:140:LYS:NZ	2.20	0.56
1:C:195:GLU:HA	3:C:646:HOH:O	2.05	0.56
1:D:8:LEU:HD11	1:D:365:THR:HG22	1.88	0.56
1:D:65:LEU:O	3:D:651:HOH:O	2.16	0.56
1:E:351:ARG:NE	1:E:355:ASP:O	2.37	0.56
1:I:197:GLY:HA3	1:I:218:ALA:HB3	1.87	0.56
1:B:199:ASP:HB2	1:B:276:THR:HA	1.87	0.56
1:C:351:ARG:NH2	3:C:602:HOH:O	2.38	0.56
1:H:316:VAL:HG13	1:H:320:TYR:CD2	2.41	0.56
1:K:91:ARG:HH21	1:L:281:GLY:C	2.13	0.56
1:A:288:ARG:NE	3:A:690:HOH:O	2.39	0.56
1:D:167:PHE:CZ	1:D:293:VAL:HG11	2.40	0.56
1:H:103:ASP:O	3:H:658:HOH:O	2.18	0.56
1:H:39:GLU:OE1	1:H:318:LYS:NZ	2.33	0.56
1:I:130:ARG:HB2	3:I:630:HOH:O	2.05	0.56
1:I:299:MET:HE3	1:J:344:GLN:OE1	2.06	0.56
1:J:9:ASN:OD1	1:J:11:LEU:HD13	2.06	0.56
1:C:11:LEU:HD23	1:D:11:LEU:HG	1.88	0.56
1:C:354:ASP:C	1:C:356:PRO:HD3	2.31	0.56
1:E:213:LEU:HD23	1:E:256:VAL:HG12	1.88	0.56
1:F:294:GLY:HA3	3:F:701:HOH:O	2.06	0.56
1:K:80:ILE:HG13	1:K:131:HIS:CD2	2.41	0.56
1:K:381:ILE:HG22	3:K:666:HOH:O	2.06	0.56
1:C:25:LYS:HD2	1:C:289:GLY:HA2	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:199:ASP:HB2	1:C:276:THR:HA	1.87	0.56
1:C:203:MET:HE3	1:D:346:LEU:HD11	1.88	0.56
1:C:374:THR:HA	3:C:644:HOH:O	2.05	0.56
1:D:2:ARG:NH2	1:D:387:SER:HB3	2.21	0.56
1:H:118:TRP:O	1:H:122:ASN:ND2	2.30	0.56
1:J:120:ARG:NH2	3:J:629:HOH:O	2.10	0.56
1:A:24:ARG:CZ	3:A:640:HOH:O	2.54	0.55
1:C:219:THR:OG1	1:C:270:TYR:O	2.22	0.55
1:C:348:GLN:NE2	1:D:212:ASP:OD2	2.35	0.55
1:E:212:ASP:OD2	1:F:357:LEU:HD21	2.06	0.55
1:I:323:LEU:HB2	1:I:392:LEU:HD21	1.86	0.55
1:A:70:GLN:OE1	1:B:140:LYS:N	2.24	0.55
1:E:19:VAL:HG21	1:F:360:ASN:ND2	2.21	0.55
1:E:374:THR:HA	3:E:678:HOH:O	2.06	0.55
1:G:35:ASP:OD1	1:G:287:GLY:N	2.32	0.55
1:K:326:LEU:HB3	1:K:385:MET:HE1	1.88	0.55
1:B:373:GLU:HG2	1:B:374:THR:H	1.71	0.55
1:H:310:LYS:HA	3:H:633:HOH:O	2.06	0.55
1:L:316:VAL:HG13	1:L:320:TYR:CD2	2.41	0.55
1:D:52:LYS:O	1:D:54:GLY:N	2.40	0.55
1:I:335:VAL:HG21	1:I:378:ILE:HG22	1.87	0.55
1:K:219:THR:OG1	1:K:270:TYR:O	2.24	0.55
1:A:144:ASP:HB3	3:A:661:HOH:O	2.05	0.55
1:C:37:VAL:HB	1:C:115:VAL:HG22	1.88	0.55
1:E:6:VAL:HB	1:E:375:LYS:HE3	1.88	0.55
1:G:346:LEU:HD11	1:H:21:LEU:HD11	1.89	0.55
1:I:91:ARG:HH21	1:J:281:GLY:C	2.14	0.55
1:A:373:GLU:HG2	1:A:374:THR:H	1.71	0.55
1:B:9:ASN:OD1	1:B:11:LEU:HD13	2.07	0.55
1:C:310:LYS:HE2	1:C:315:HIS:NE2	2.22	0.55
1:E:257:TYR:HB3	1:E:260:THR:CG2	2.37	0.55
1:F:6:VAL:CG1	1:F:378:ILE:HD11	2.37	0.55
1:F:53:TYR:O	1:F:55:VAL:HG22	2.07	0.55
1:F:316:VAL:HG21	3:F:706:HOH:O	2.07	0.55
1:I:80:ILE:HG13	1:I:131:HIS:CD2	2.41	0.55
1:A:10:PRO:HA	1:A:364:ILE:HD11	1.87	0.55
1:A:74:ARG:HD3	3:A:609:HOH:O	2.06	0.55
1:B:379:SER:HB2	3:B:687:HOH:O	2.06	0.55
1:C:178:TYR:HB2	1:C:295:LEU:HD21	1.89	0.55
1:D:326:LEU:HB3	1:D:385:MET:HE1	1.88	0.55
1:H:140:LYS:O	3:H:637:HOH:O	2.18	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:310:LYS:HB2	1:J:318:LYS:HD2	1.89	0.55
1:L:324:ALA:N	3:L:657:HOH:O	2.39	0.55
1:B:213:LEU:HB3	1:B:256:VAL:HG12	1.89	0.55
1:I:304:LEU:HD12	1:J:296:ILE:HG23	1.88	0.55
1:D:103:ASP:OD1	1:D:103:ASP:N	2.39	0.55
1:D:181:GLU:OE2	1:D:185:ASN:ND2	2.39	0.55
1:D:301:PRO:HA	3:D:636:HOH:O	2.07	0.55
1:E:138:ILE:HG22	1:E:139:GLY:O	2.07	0.55
1:F:354:ASP:HA	1:F:389:PHE:HE2	1.72	0.55
1:H:209:ASN:O	1:H:253:ASN:N	2.30	0.55
1:I:14:ILE:HD11	1:I:171:THR:HG21	1.88	0.55
1:K:238:ARG:HB2	1:K:256:VAL:HG23	1.89	0.55
1:L:181:GLU:OE1	1:L:182:ARG:HD2	2.07	0.55
1:C:288:ARG:NH1	2:D:501:PO4:O1	2.38	0.54
1:D:373:GLU:HG2	1:D:374:THR:H	1.72	0.54
1:G:257:TYR:HB3	1:G:260:THR:CG2	2.37	0.54
1:D:344:GLN:HG3	3:D:619:HOH:O	2.07	0.54
1:E:268:ILE:C	1:E:269:LEU:HD12	2.32	0.54
1:F:83:PRO:O	3:F:614:HOH:O	2.19	0.54
1:F:157:LEU:O	1:F:314:ASN:HB2	2.07	0.54
1:G:218:ALA:HA	1:G:270:TYR:HD2	1.72	0.54
1:L:238:ARG:HG3	3:L:640:HOH:O	2.07	0.54
1:L:270:TYR:OH	1:L:277:SER:HB3	2.07	0.54
1:L:378:ILE:HG13	1:L:379:SER:N	2.21	0.54
1:H:48:TYR:CZ	1:H:52:LYS:HD2	2.42	0.54
1:J:270:TYR:OH	1:J:277:SER:HB3	2.06	0.54
1:E:338:VAL:HB	3:E:664:HOH:O	2.07	0.54
1:H:169:PRO:HD2	1:H:340:PHE:HD1	1.72	0.54
1:K:46:SER:OG	1:K:59:HIS:NE2	2.40	0.54
1:A:35:ASP:OD1	1:A:287:GLY:N	2.32	0.54
1:H:218:ALA:HA	1:H:270:TYR:CD2	2.42	0.54
1:A:6:VAL:CG1	1:A:378:ILE:HD11	2.36	0.54
1:B:328:ALA:HB2	1:B:343:VAL:HG23	1.90	0.54
1:G:71:ALA:H	1:G:280:HIS:CD2	2.22	0.54
1:H:315:HIS:N	3:H:640:HOH:O	2.35	0.54
1:I:184:LEU:O	1:I:190:LYS:NZ	2.41	0.54
1:B:217:MET:HE1	1:B:237:VAL:HG21	1.88	0.54
1:F:218:ALA:HA	1:F:270:TYR:HD2	1.73	0.54
1:F:382:VAL:O	3:F:705:HOH:O	2.18	0.54
1:I:378:ILE:HG13	1:I:379:SER:N	2.22	0.54
1:C:326:LEU:HB3	1:C:385:MET:HE1	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:70:GLN:HB2	1:D:82:GLN:HB3	1.88	0.54
1:E:50:LEU:N	3:E:674:HOH:O	2.41	0.54
1:F:219:THR:OG1	1:F:270:TYR:O	2.26	0.54
1:G:19:VAL:HG21	1:H:360:ASN:ND2	2.22	0.54
1:I:346:LEU:CD1	1:J:21:LEU:HD11	2.38	0.54
1:L:179:GLU:O	3:L:658:HOH:O	2.17	0.54
1:B:378:ILE:HG13	1:B:379:SER:N	2.22	0.54
1:G:210:GLU:HG2	1:G:253:ASN:HB3	1.89	0.54
1:H:158:SER:HB3	1:H:348:GLN:O	2.08	0.54
1:J:335:VAL:HG21	1:J:378:ILE:HG22	1.88	0.54
1:L:345:VAL:HG21	3:L:657:HOH:O	2.08	0.54
1:B:2:ARG:NH2	1:B:387:SER:HB3	2.23	0.54
1:B:45:LEU:HD13	3:B:691:HOH:O	2.08	0.54
1:C:68:GLY:O	1:D:91:ARG:NH2	2.40	0.54
1:D:2:ARG:HG3	1:D:3:ASN:H	1.73	0.54
1:D:373:GLU:N	3:D:638:HOH:O	2.38	0.54
1:I:289:GLY:N	3:I:625:HOH:O	2.40	0.54
1:C:70:GLN:HG2	1:C:280:HIS:CD2	2.44	0.53
1:C:294:GLY:HA3	3:C:676:HOH:O	2.07	0.53
1:D:158:SER:HB3	1:D:348:GLN:O	2.08	0.53
1:F:195:GLU:OE1	1:F:195:GLU:N	2.39	0.53
1:G:312:PRO:O	1:G:393:THR:OG1	2.16	0.53
1:H:222:GLU:HG3	3:H:628:HOH:O	2.07	0.53
1:A:329:ASN:O	1:A:333:GLN:HG3	2.08	0.53
1:B:6:VAL:CG1	1:B:378:ILE:HD11	2.38	0.53
1:F:2:ARG:N	3:F:682:HOH:O	2.40	0.53
1:J:147:GLY:O	3:J:631:HOH:O	2.18	0.53
1:J:354:ASP:HA	1:J:389:PHE:HE2	1.74	0.53
1:B:207:ARG:HD3	3:B:602:HOH:O	2.08	0.53
1:C:70:GLN:OE1	1:D:140:LYS:N	2.34	0.53
1:D:311:ASN:N	3:D:604:HOH:O	2.22	0.53
1:D:318:LYS:HE2	1:D:396:ILE:HD13	1.90	0.53
1:I:103:ASP:OD1	1:I:103:ASP:N	2.36	0.53
1:L:268:ILE:C	1:L:269:LEU:HD12	2.34	0.53
1:L:316:VAL:HB	3:L:656:HOH:O	2.08	0.53
1:A:357:LEU:O	1:A:357:LEU:HD23	2.08	0.53
1:B:165:VAL:HG22	3:B:650:HOH:O	2.08	0.53
1:C:205:LEU:HD23	1:D:357:LEU:HD22	1.90	0.53
1:C:273:VAL:HB	3:C:682:HOH:O	2.08	0.53
1:E:299:MET:HE3	1:F:344:GLN:NE2	2.22	0.53
1:J:74:ARG:HD3	3:J:603:HOH:O	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:318:LYS:HE2	1:J:396:ILE:HD13	1.90	0.53
1:A:228:ASN:ND2	1:E:264:ILE:HG12	2.24	0.53
1:A:337:ASP:O	3:A:707:HOH:O	2.18	0.53
1:H:21:LEU:HD13	1:H:299:MET:HE1	1.90	0.53
1:E:256:VAL:HG11	3:E:661:HOH:O	2.09	0.53
1:F:302:MET:N	3:F:641:HOH:O	2.34	0.53
1:K:195:GLU:OE1	1:K:195:GLU:N	2.39	0.53
1:L:218:ALA:HA	1:L:270:TYR:CD2	2.44	0.53
1:H:277:SER:O	1:H:282:ASP:HB2	2.08	0.53
1:K:22:VAL:HB	1:K:177:VAL:HG21	1.90	0.53
1:K:70:GLN:HA	1:K:280:HIS:CD2	2.44	0.53
1:F:9:ASN:OD1	1:F:11:LEU:HD13	2.08	0.53
1:I:373:GLU:HG2	1:I:374:THR:H	1.74	0.53
1:K:378:ILE:HG13	1:K:379:SER:N	2.23	0.53
1:L:300:ARG:HG2	3:L:636:HOH:O	2.08	0.53
1:I:207:ARG:NE	1:J:5:ASN:HD22	2.06	0.53
1:J:199:ASP:HB2	1:J:276:THR:HA	1.91	0.53
1:L:14:ILE:HB	1:L:17:LEU:HD12	1.91	0.53
1:L:144:ASP:OD1	1:L:145:LEU:HD12	2.09	0.53
1:L:161:THR:HG22	1:L:349:ILE:H	1.73	0.53
1:F:238:ARG:HB2	1:F:256:VAL:CG2	2.38	0.53
1:K:196:VAL:HB	3:K:642:HOH:O	2.09	0.53
1:A:263:LYS:HB2	1:A:268:ILE:CG1	2.39	0.52
1:I:329:ASN:O	1:I:333:GLN:HG3	2.09	0.52
1:L:230:TYR:OH	1:L:262:ASP:OD2	2.22	0.52
1:A:218:ALA:HA	1:A:270:TYR:HD2	1.74	0.52
1:C:31:ASP:HB3	3:C:628:HOH:O	2.09	0.52
1:D:169:PRO:HD2	1:D:340:PHE:HD1	1.74	0.52
1:D:375:LYS:HG3	3:D:638:HOH:O	2.08	0.52
1:K:299:MET:HE3	1:L:344:GLN:OE1	2.09	0.52
1:L:282:ASP:N	3:L:632:HOH:O	2.41	0.52
1:A:271:LEU:HG	3:A:655:HOH:O	2.09	0.52
1:A:340:PHE:N	3:A:650:HOH:O	2.42	0.52
1:B:53:TYR:CE1	1:B:96:VAL:HG21	2.43	0.52
1:C:120:ARG:NH2	1:C:129:GLU:OE2	2.42	0.52
1:E:222:GLU:HG3	3:E:663:HOH:O	2.09	0.52
1:F:328:ALA:HB2	1:F:343:VAL:HG23	1.90	0.52
1:G:323:LEU:HD21	1:G:386:LEU:HD11	1.91	0.52
1:J:211:VAL:HB	1:J:254:VAL:HG12	1.91	0.52
1:K:118:TRP:O	1:K:122:ASN:ND2	2.37	0.52
1:B:22:VAL:HB	1:B:177:VAL:HG21	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:207:ARG:NE	1:D:5:ASN:HD22	2.07	0.52
1:F:6:VAL:HG11	1:F:378:ILE:HD11	1.92	0.52
1:F:299:MET:HB2	3:F:648:HOH:O	2.09	0.52
1:L:9:ASN:OD1	1:L:11:LEU:HD13	2.09	0.52
1:G:201:LYS:NZ	1:H:160:ASP:OD2	2.43	0.52
1:G:346:LEU:HD23	1:G:347:GLY:N	2.25	0.52
1:H:140:LYS:N	3:H:637:HOH:O	2.42	0.52
1:J:10:PRO:HA	1:J:364:ILE:HD11	1.91	0.52
1:J:120:ARG:NH2	1:J:129:GLU:OE2	2.43	0.52
1:J:217:MET:O	1:J:270:TYR:HD2	1.93	0.52
1:J:365:THR:C	1:J:367:ASP:H	2.18	0.52
1:D:162:SER:OG	1:D:316:VAL:HG12	2.09	0.52
1:D:316:VAL:HG13	1:D:320:TYR:CD2	2.44	0.52
1:I:290:ASN:HB3	1:I:294:GLY:C	2.35	0.52
1:J:14:ILE:HB	1:J:17:LEU:HD12	1.91	0.52
1:D:222:GLU:HG3	3:D:634:HOH:O	2.10	0.52
1:G:222:GLU:HG3	3:G:615:HOH:O	2.09	0.52
1:I:238:ARG:HB2	1:I:256:VAL:HG23	1.92	0.52
1:E:136:TYR:HD2	1:E:138:ILE:HG12	1.75	0.52
1:E:296:ILE:HG21	3:F:658:HOH:O	2.10	0.52
1:F:218:ALA:HA	1:F:270:TYR:CD2	2.45	0.52
1:K:103:ASP:OD1	1:K:103:ASP:N	2.34	0.52
1:L:181:GLU:OE2	1:L:185:ASN:ND2	2.38	0.52
1:L:222:GLU:HB2	1:L:223:LEU:HD12	1.92	0.52
1:C:2:ARG:O	1:C:3:ASN:HB2	2.10	0.52
1:G:288:ARG:HG3	3:G:644:HOH:O	2.10	0.52
1:K:346:LEU:HD23	1:K:347:GLY:N	2.25	0.52
1:C:68:GLY:HA2	1:C:278:ALA:O	2.10	0.51
1:C:86:ILE:N	3:C:642:HOH:O	2.23	0.51
1:E:326:LEU:HB3	1:E:385:MET:HE1	1.93	0.51
1:G:329:ASN:O	1:G:333:GLN:HG3	2.08	0.51
1:J:24:ARG:HH21	1:J:182:ARG:HH11	1.56	0.51
1:K:75:PHE:CZ	1:K:227:VAL:HG22	2.45	0.51
1:K:304:LEU:HD12	1:L:296:ILE:HG12	1.92	0.51
1:A:2:ARG:HH12	1:A:387:SER:HB3	1.75	0.51
1:A:323:LEU:HB2	1:A:392:LEU:HD21	1.92	0.51
1:E:316:VAL:HG21	3:E:639:HOH:O	2.10	0.51
1:E:337:ASP:O	1:E:365:THR:OG1	2.19	0.51
1:F:199:ASP:HB2	1:F:276:THR:HA	1.90	0.51
1:I:323:LEU:HD21	1:I:386:LEU:HD11	1.91	0.51
1:L:15:GLU:HG3	1:L:171:THR:HB	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:238:ARG:NH2	3:A:653:HOH:O	2.29	0.51
1:D:9:ASN:OD1	1:D:11:LEU:HD13	2.10	0.51
1:D:237:VAL:O	1:D:240:GLN:HB2	2.11	0.51
1:F:382:VAL:O	1:F:386:LEU:HD13	2.10	0.51
1:H:81:ILE:HD13	3:H:627:HOH:O	2.11	0.51
1:K:217:MET:HE1	1:K:237:VAL:HG21	1.91	0.51
1:L:14:ILE:CD1	1:L:171:THR:HG21	2.40	0.51
1:D:56:ILE:HD11	1:D:397:LEU:HD11	1.92	0.51
1:F:127:ASP:OD1	1:F:130:ARG:HG3	2.10	0.51
1:F:238:ARG:HB2	1:F:256:VAL:HG23	1.92	0.51
1:I:270:TYR:O	1:I:270:TYR:CD2	2.63	0.51
1:A:29:HIS:NE2	2:A:501:PO4:O2	2.42	0.51
1:C:373:GLU:HG2	1:C:374:THR:H	1.76	0.51
1:J:291:ARG:HG3	1:J:302:MET:HG2	1.92	0.51
1:L:2:ARG:HG3	1:L:3:ASN:H	1.74	0.51
1:B:161:THR:HG23	1:B:349:ILE:HG22	1.92	0.51
1:E:8:LEU:HD11	1:E:365:THR:HG22	1.92	0.51
1:I:70:GLN:HG2	1:I:280:HIS:CE1	2.45	0.51
1:I:222:GLU:HG3	3:I:615:HOH:O	2.09	0.51
1:B:195:GLU:OE1	1:B:195:GLU:N	2.40	0.51
1:B:218:ALA:HA	1:B:270:TYR:CD2	2.45	0.51
1:E:342:GLN:HG3	3:E:625:HOH:O	2.11	0.51
1:E:354:ASP:HB2	1:E:386:LEU:O	2.11	0.51
1:L:34:ALA:C	1:L:286:THR:HG22	2.36	0.51
1:B:20:GLU:HA	1:B:299:MET:SD	2.50	0.51
1:D:320:TYR:HE2	1:D:347:GLY:HA3	1.75	0.51
1:H:6:VAL:CG1	1:H:378:ILE:HD11	2.40	0.51
1:A:300:ARG:HB3	3:A:689:HOH:O	2.10	0.51
1:A:378:ILE:HG13	1:A:379:SER:N	2.25	0.51
1:D:207:ARG:HD3	3:D:601:HOH:O	2.11	0.51
1:E:25:LYS:HD2	1:E:289:GLY:HA2	1.92	0.51
1:G:373:GLU:HG2	1:G:374:THR:N	2.26	0.51
1:H:182:ARG:NH1	3:H:605:HOH:O	2.43	0.51
1:A:289:GLY:N	3:A:704:HOH:O	2.43	0.51
1:B:60:ASN:ND2	3:B:661:HOH:O	2.35	0.51
1:C:217:MET:HE3	1:C:258:VAL:HG13	1.93	0.51
1:D:339:LYS:HG3	1:D:366:TYR:CE1	2.46	0.51
1:E:138:ILE:HD12	3:E:706:HOH:O	2.11	0.51
1:H:373:GLU:HG2	1:H:374:THR:H	1.76	0.51
1:K:19:VAL:HG11	1:L:360:ASN:HD22	1.74	0.51
1:D:268:ILE:C	1:D:269:LEU:HD12	2.36	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:21:LEU:HD13	1:I:299:MET:HE1	1.92	0.50
1:I:219:THR:OG1	1:I:270:TYR:O	2.29	0.50
1:K:62:ASP:OD2	1:K:310:LYS:NZ	2.35	0.50
1:L:12:SER:HB3	3:L:663:HOH:O	2.11	0.50
1:D:124:ARG:NE	1:D:198:GLU:OE2	2.40	0.50
1:F:197:GLY:N	1:F:218:ALA:O	2.34	0.50
1:H:365:THR:C	1:H:367:ASP:H	2.19	0.50
1:L:318:LYS:HE2	1:L:396:ILE:HD13	1.92	0.50
1:A:50:LEU:HD13	1:A:56:ILE:HG13	1.93	0.50
1:A:199:ASP:OD2	1:A:277:SER:OG	2.21	0.50
3:A:694:HOH:O	1:B:90:GLY:N	2.40	0.50
1:H:268:ILE:C	1:H:269:LEU:HD12	2.36	0.50
1:A:14:ILE:CD1	1:A:171:THR:HG21	2.40	0.50
1:B:6:VAL:HG11	1:B:378:ILE:HD11	1.92	0.50
1:E:53:TYR:CE1	1:E:96:VAL:HG21	2.47	0.50
1:G:22:VAL:HB	1:G:177:VAL:HG21	1.94	0.50
1:H:52:LYS:O	1:H:54:GLY:N	2.42	0.50
1:L:162:SER:OG	1:L:316:VAL:HG12	2.11	0.50
1:A:69:GLY:O	1:A:70:GLN:HG3	2.12	0.50
1:A:348:GLN:NE2	1:B:212:ASP:OD2	2.35	0.50
1:F:163:PHE:HB2	1:F:346:LEU:HD13	1.93	0.50
1:J:354:ASP:HB2	1:J:386:LEU:O	2.12	0.50
1:B:160:ASP:HB3	3:B:693:HOH:O	2.10	0.50
1:D:217:MET:HE1	1:D:237:VAL:HG21	1.94	0.50
1:E:16:LYS:HD3	3:E:620:HOH:O	2.12	0.50
1:E:76:LYS:HG3	1:E:225:GLU:HA	1.94	0.50
1:H:289:GLY:N	3:H:626:HOH:O	2.43	0.50
1:K:218:ALA:HA	1:K:270:TYR:CD2	2.46	0.50
1:L:163:PHE:HB3	1:L:346:LEU:HD13	1.94	0.50
1:E:36:ALA:O	3:E:718:HOH:O	2.20	0.50
1:H:354:ASP:HA	1:H:389:PHE:HE2	1.77	0.50
1:K:76:LYS:HG3	1:K:225:GLU:HA	1.94	0.50
1:B:14:ILE:CD1	1:B:171:THR:HG21	2.38	0.50
1:D:60:ASN:HA	3:D:604:HOH:O	2.11	0.50
1:F:323:LEU:O	1:F:326:LEU:N	2.45	0.50
1:G:199:ASP:HB2	1:G:276:THR:HA	1.94	0.50
1:I:217:MET:HE3	1:I:258:VAL:HG13	1.93	0.50
1:J:354:ASP:HA	1:J:389:PHE:CE2	2.47	0.50
1:K:22:VAL:HG21	1:K:174:GLU:HG2	1.93	0.50
1:C:320:TYR:OH	3:C:630:HOH:O	2.18	0.50
1:E:55:VAL:O	3:E:674:HOH:O	2.17	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:80:ILE:HG13	1:F:131:HIS:CD2	2.47	0.50
1:F:373:GLU:HG2	1:F:374:THR:H	1.76	0.50
1:G:218:ALA:HA	1:G:270:TYR:CD2	2.47	0.50
1:H:50:LEU:HD13	1:H:56:ILE:HG13	1.93	0.50
1:H:103:ASP:OD1	1:H:103:ASP:N	2.45	0.50
1:J:2:ARG:HG3	1:J:3:ASN:H	1.77	0.50
1:A:269:LEU:HB3	3:A:655:HOH:O	2.12	0.49
1:C:238:ARG:HB2	1:C:256:VAL:CG2	2.42	0.49
1:D:90:GLY:N	3:D:644:HOH:O	2.38	0.49
1:E:103:ASP:OD1	1:E:103:ASP:N	2.42	0.49
1:I:10:PRO:HA	1:I:364:ILE:HD11	1.94	0.49
1:J:157:LEU:O	1:J:314:ASN:HB2	2.12	0.49
1:J:290:ASN:ND2	1:J:404:PHE:O	2.30	0.49
1:K:273:VAL:H	1:K:279:GLU:CD	2.19	0.49
1:C:60:ASN:ND2	1:C:62:ASP:OD1	2.43	0.49
1:E:353:ILE:HD12	1:E:389:PHE:CE1	2.46	0.49
1:F:51:LYS:HE2	3:F:696:HOH:O	2.12	0.49
1:F:316:VAL:HG13	1:F:320:TYR:CD2	2.47	0.49
1:G:210:GLU:HA	1:G:253:ASN:HB3	1.93	0.49
1:K:348:GLN:NE2	1:L:212:ASP:OD2	2.45	0.49
1:A:24:ARG:NH2	3:A:685:HOH:O	2.34	0.49
1:C:257:TYR:HB3	1:C:260:THR:CG2	2.42	0.49
1:G:348:GLN:HB2	1:G:357:LEU:HD12	1.94	0.49
1:H:311:ASN:N	3:H:603:HOH:O	2.29	0.49
1:I:195:GLU:OE1	1:I:195:GLU:N	2.43	0.49
1:J:268:ILE:C	1:J:269:LEU:HD12	2.38	0.49
1:A:181:GLU:OE2	1:A:185:ASN:ND2	2.45	0.49
1:A:257:TYR:HD2	1:B:348:GLN:HE22	1.58	0.49
1:A:326:LEU:HB3	1:A:385:MET:HE1	1.93	0.49
1:C:130:ARG:HD3	3:C:651:HOH:O	2.11	0.49
1:D:181:GLU:OE1	1:D:182:ARG:HD2	2.11	0.49
1:F:24:ARG:HH21	1:F:182:ARG:NH1	2.08	0.49
1:L:238:ARG:HB2	1:L:256:VAL:HG23	1.94	0.49
1:A:158:SER:O	1:A:349:ILE:HA	2.13	0.49
1:A:373:GLU:HG2	1:A:374:THR:N	2.27	0.49
1:B:356:PRO:HG2	1:B:386:LEU:HD21	1.95	0.49
1:D:310:LYS:HB2	1:D:318:LYS:HB2	1.93	0.49
1:E:323:LEU:HB2	1:E:392:LEU:HD21	1.94	0.49
1:E:335:VAL:O	1:E:338:VAL:HG22	2.12	0.49
1:F:14:ILE:CD1	1:F:171:THR:HG21	2.39	0.49
1:L:365:THR:C	1:L:367:ASP:H	2.21	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:57:LEU:HD22	1:A:93:THR:HG21	1.95	0.49
1:B:218:ALA:HB2	1:B:270:TYR:HE2	1.78	0.49
1:D:6:VAL:CG1	1:D:378:ILE:HD11	2.43	0.49
1:D:61:LEU:O	3:D:641:HOH:O	2.20	0.49
1:E:25:LYS:HD3	1:E:29:HIS:CE1	2.48	0.49
1:E:273:VAL:N	1:E:279:GLU:OE2	2.31	0.49
1:I:335:VAL:O	1:I:338:VAL:HG22	2.13	0.49
1:D:143:SER:O	1:D:146:VAL:HG12	2.13	0.49
1:E:158:SER:O	1:E:349:ILE:HA	2.12	0.49
1:L:310:LYS:HB2	1:L:318:LYS:HB2	1.95	0.49
1:A:351:ARG:NE	1:A:355:ASP:O	2.45	0.49
1:B:291:ARG:HG3	1:B:302:MET:HG2	1.94	0.49
1:B:354:ASP:HA	1:B:389:PHE:HE2	1.78	0.49
1:C:103:ASP:OD1	1:C:103:ASP:N	2.41	0.49
1:K:365:THR:HG23	1:K:367:ASP:HB2	1.95	0.49
1:B:75:PHE:CZ	1:B:227:VAL:HG22	2.48	0.49
1:C:10:PRO:HA	1:C:364:ILE:HD11	1.95	0.49
1:F:52:LYS:NZ	3:F:697:HOH:O	2.27	0.49
1:J:237:VAL:O	1:J:240:GLN:HB2	2.13	0.49
1:K:167:PHE:CZ	1:K:293:VAL:HG21	2.48	0.49
1:A:228:ASN:HD21	1:E:264:ILE:HG12	1.78	0.48
1:B:162:SER:OG	1:B:316:VAL:HG12	2.13	0.48
1:C:346:LEU:HD23	1:C:347:GLY:N	2.28	0.48
1:E:274:THR:N	1:E:279:GLU:OE1	2.41	0.48
1:I:118:TRP:O	1:I:122:ASN:ND2	2.32	0.48
1:I:310:LYS:O	1:I:318:LYS:NZ	2.41	0.48
1:K:190:LYS:HD3	3:K:642:HOH:O	2.12	0.48
1:K:213:LEU:HD23	1:K:256:VAL:HG12	1.94	0.48
1:A:213:LEU:HB3	1:A:256:VAL:HG12	1.95	0.48
1:B:165:VAL:HG23	1:B:303:SER:HB2	1.95	0.48
1:C:218:ALA:HA	1:C:270:TYR:CD2	2.47	0.48
1:D:320:TYR:HE2	1:D:347:GLY:CA	2.26	0.48
1:D:328:ALA:HB2	1:D:343:VAL:HG23	1.95	0.48
1:E:169:PRO:O	3:E:673:HOH:O	2.20	0.48
1:H:70:GLN:HB2	1:H:82:GLN:HB3	1.94	0.48
1:H:120:ARG:NH2	3:H:634:HOH:O	2.25	0.48
1:A:338:VAL:HB	3:A:650:HOH:O	2.13	0.48
1:E:2:ARG:HH12	1:E:387:SER:HB3	1.78	0.48
1:F:2:ARG:HG3	1:F:3:ASN:H	1.78	0.48
1:D:27:LEU:O	3:D:678:HOH:O	2.20	0.48
1:D:249:ALA:O	1:D:252:TYR:HB2	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:207:ARG:NH2	1:F:5:ASN:HB2	2.27	0.48
1:H:2:ARG:NH1	1:H:387:SER:HB3	2.28	0.48
1:K:223:LEU:HD12	1:K:223:LEU:N	2.28	0.48
1:L:195:GLU:OE1	1:L:195:GLU:N	2.39	0.48
1:E:206:ARG:HD2	1:E:252:TYR:CZ	2.48	0.48
1:F:253:ASN:OD1	1:F:255:ARG:NH2	2.45	0.48
1:H:39:GLU:HA	1:H:309:GLY:HA2	1.95	0.48
1:I:243:ASP:HB3	3:I:604:HOH:O	2.13	0.48
1:J:48:TYR:OH	3:J:625:HOH:O	2.16	0.48
1:L:354:ASP:HA	1:L:389:PHE:CE2	2.46	0.48
1:A:190:LYS:HD3	3:A:701:HOH:O	2.12	0.48
1:C:138:ILE:HG22	1:C:139:GLY:O	2.14	0.48
1:E:68:GLY:HA2	1:E:278:ALA:O	2.13	0.48
1:L:70:GLN:HB2	1:L:82:GLN:HB3	1.96	0.48
1:E:63:LYS:HE2	1:F:283:ASP:HB2	1.95	0.48
1:C:268:ILE:C	1:C:269:LEU:HD12	2.38	0.48
1:E:242:LEU:HG	3:E:661:HOH:O	2.13	0.48
1:G:6:VAL:CG1	1:G:378:ILE:HD11	2.44	0.48
1:I:20:GLU:HA	1:I:299:MET:SD	2.54	0.48
1:I:193:LEU:HD23	1:I:193:LEU:HA	1.70	0.48
1:I:373:GLU:HG2	1:I:374:THR:N	2.29	0.48
1:L:15:GLU:CG	1:L:171:THR:HB	2.44	0.48
1:L:171:THR:HG23	1:L:300:ARG:HH21	1.78	0.48
1:L:329:ASN:O	1:L:333:GLN:HG3	2.14	0.48
1:E:280:HIS:CE1	1:F:140:LYS:O	2.67	0.48
1:F:103:ASP:OD1	1:F:103:ASP:N	2.44	0.48
1:H:157:LEU:O	1:H:314:ASN:HB2	2.14	0.48
1:H:291:ARG:HG3	1:H:302:MET:HG2	1.96	0.48
1:I:280:HIS:CE1	1:J:140:LYS:O	2.67	0.48
1:A:71:ALA:N	1:A:280:HIS:HD2	2.05	0.48
1:B:7:GLN:O	1:B:363:VAL:HG12	2.13	0.48
1:D:13:ASP:HA	3:D:676:HOH:O	2.13	0.48
1:D:230:TYR:CD2	1:D:269:LEU:HD23	2.49	0.48
1:D:301:PRO:HD2	3:D:635:HOH:O	2.12	0.48
1:E:238:ARG:NH2	3:E:717:HOH:O	2.46	0.48
1:F:163:PHE:HE1	1:F:165:VAL:HG22	1.79	0.48
1:F:244:LEU:HA	1:F:247:LYS:HD3	1.94	0.48
1:F:310:LYS:HA	3:F:667:HOH:O	2.13	0.48
1:H:235:GLU:O	1:H:239:ASN:ND2	2.47	0.48
1:A:263:LYS:HB2	1:A:268:ILE:HG12	1.96	0.47
1:A:323:LEU:HD21	1:A:386:LEU:HD11	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:220:ILE:HB	3:D:677:HOH:O	2.14	0.47
1:F:15:GLU:HG3	1:F:171:THR:HB	1.96	0.47
1:F:167:PHE:CZ	1:F:293:VAL:HG11	2.48	0.47
1:G:158:SER:O	1:G:349:ILE:HA	2.13	0.47
1:I:310:LYS:HE2	1:I:315:HIS:NE2	2.29	0.47
1:K:116:LYS:HB3	3:K:663:HOH:O	2.14	0.47
1:B:58:HIS:ND1	3:B:676:HOH:O	2.35	0.47
1:B:342:GLN:O	3:B:612:HOH:O	2.20	0.47
1:C:329:ASN:O	1:C:333:GLN:HG3	2.15	0.47
1:D:323:LEU:HB2	1:D:392:LEU:HD21	1.95	0.47
1:G:96:VAL:HG11	3:G:616:HOH:O	2.14	0.47
1:K:365:THR:C	1:K:367:ASP:H	2.22	0.47
1:A:339:LYS:HG3	1:A:366:TYR:HE1	1.79	0.47
1:B:320:TYR:HE2	1:B:347:GLY:HA3	1.80	0.47
1:C:346:LEU:HD11	1:D:21:LEU:HD11	1.97	0.47
1:D:21:LEU:HD13	1:D:299:MET:HE1	1.96	0.47
1:D:77:GLY:O	1:D:221:SER:OG	2.31	0.47
1:E:242:LEU:N	3:E:661:HOH:O	2.47	0.47
1:E:301:PRO:O	3:E:713:HOH:O	2.20	0.47
1:K:160:ASP:OD1	1:K:161:THR:N	2.47	0.47
1:K:307:THR:OG1	1:K:403:LEU:HB3	2.14	0.47
1:L:382:VAL:O	1:L:386:LEU:HD13	2.14	0.47
1:C:195:GLU:OE1	1:C:195:GLU:N	2.46	0.47
1:D:142:SER:N	3:D:672:HOH:O	2.23	0.47
1:D:163:PHE:CZ	1:D:165:VAL:HG22	2.49	0.47
1:G:335:VAL:O	1:G:338:VAL:HG22	2.14	0.47
1:A:224:ILE:HD12	3:A:703:HOH:O	2.14	0.47
1:I:318:LYS:O	1:I:322:VAL:HG23	2.14	0.47
1:F:193:LEU:HA	1:F:193:LEU:HD23	1.71	0.47
3:F:704:HOH:O	1:G:83:PRO:HD3	2.14	0.47
1:L:6:VAL:CG1	1:L:378:ILE:HD11	2.45	0.47
1:L:224:ILE:HG21	1:L:271:LEU:HD21	1.97	0.47
1:L:321:ASN:ND2	3:L:623:HOH:O	2.48	0.47
1:A:255:ARG:HH21	1:E:246:SER:CB	2.27	0.47
1:B:111:ILE:O	1:B:114:SER:OG	2.21	0.47
1:B:296:ILE:HG22	1:B:298:PRO:HD3	1.95	0.47
1:D:378:ILE:HG13	1:D:379:SER:N	2.30	0.47
1:E:14:ILE:HB	1:E:17:LEU:HD12	1.96	0.47
1:E:230:TYR:HD1	3:E:712:HOH:O	1.98	0.47
1:G:238:ARG:HB2	1:G:256:VAL:CG2	2.44	0.47
1:H:108:GLY:O	1:H:112:ILE:HG12	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:127:ASP:OD1	1:J:130:ARG:HG3	2.14	0.47
1:A:6:VAL:HG11	1:A:378:ILE:HD11	1.97	0.47
1:A:217:MET:HE1	1:A:237:VAL:HG21	1.96	0.47
1:E:167:PHE:CZ	1:E:293:VAL:HG21	2.50	0.47
1:E:381:ILE:HD12	3:E:671:HOH:O	2.15	0.47
1:L:21:LEU:HD13	1:L:299:MET:HE1	1.96	0.47
1:L:305:GLU:HG2	2:L:501:PO4:O2	2.15	0.47
1:E:22:VAL:HG21	1:E:174:GLU:HG2	1.97	0.47
1:B:305:GLU:HB2	3:B:632:HOH:O	2.14	0.47
1:D:14:ILE:CD1	1:D:171:THR:HG21	2.38	0.47
1:G:178:TYR:HB2	1:G:295:LEU:HD21	1.96	0.47
1:H:5:ASN:HD22	1:H:6:VAL:N	2.13	0.47
1:K:14:ILE:CD1	1:K:171:THR:HG21	2.43	0.47
1:L:161:THR:HG23	1:L:349:ILE:HG22	1.97	0.47
1:A:304:LEU:HD12	1:B:296:ILE:HG12	1.97	0.46
1:D:161:THR:HG23	1:D:349:ILE:HG22	1.97	0.46
1:D:270:TYR:OH	1:D:277:SER:HB3	2.15	0.46
1:F:117:GLU:OE1	1:F:120:ARG:NH1	2.47	0.46
1:F:403:LEU:HD22	3:F:662:HOH:O	2.14	0.46
1:G:195:GLU:OE1	1:G:195:GLU:N	2.40	0.46
1:H:354:ASP:HB2	1:H:386:LEU:O	2.15	0.46
1:I:91:ARG:HG3	3:J:617:HOH:O	2.15	0.46
1:J:75:PHE:CZ	1:J:227:VAL:HG22	2.50	0.46
1:J:356:PRO:HG2	1:J:386:LEU:HD21	1.97	0.46
1:B:316:VAL:HG13	1:B:320:TYR:CD2	2.49	0.46
1:D:373:GLU:HG2	1:D:374:THR:N	2.30	0.46
1:E:213:LEU:HB3	1:E:256:VAL:HG12	1.97	0.46
1:E:395:LEU:HA	1:E:398:GLU:OE2	2.16	0.46
1:H:161:THR:HG22	1:H:349:ILE:H	1.79	0.46
1:H:206:ARG:HD2	1:H:252:TYR:CZ	2.50	0.46
1:L:218:ALA:HB2	1:L:270:TYR:HE2	1.81	0.46
1:L:373:GLU:OE2	3:L:630:HOH:O	2.20	0.46
1:A:70:GLN:HG2	1:A:280:HIS:CD2	2.49	0.46
1:D:89:ALA:HB1	3:D:644:HOH:O	2.15	0.46
1:F:181:GLU:OE1	1:F:182:ARG:NH1	2.49	0.46
1:G:346:LEU:CD1	1:H:21:LEU:HD11	2.45	0.46
1:H:335:VAL:O	1:H:338:VAL:HG22	2.15	0.46
1:J:118:TRP:O	1:J:122:ASN:ND2	2.37	0.46
1:K:217:MET:HE3	1:K:258:VAL:HG13	1.97	0.46
1:L:52:LYS:C	1:L:54:GLY:H	2.19	0.46
1:L:351:ARG:NE	1:L:355:ASP:O	2.43	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:339:LYS:N	3:A:650:HOH:O	2.48	0.46
1:C:222:GLU:HB2	1:C:223:LEU:HD12	1.97	0.46
1:E:91:ARG:HG3	3:E:680:HOH:O	2.15	0.46
1:H:392:LEU:O	1:H:396:ILE:HG13	2.15	0.46
1:J:373:GLU:HG2	1:J:374:THR:N	2.31	0.46
1:K:354:ASP:C	1:K:356:PRO:HD3	2.39	0.46
1:A:21:LEU:O	1:A:298:PRO:HD2	2.16	0.46
1:B:234:LYS:NZ	1:B:262:ASP:OD2	2.40	0.46
1:D:62:ASP:CG	1:D:310:LYS:HD3	2.41	0.46
1:G:2:ARG:NH1	1:G:387:SER:HB3	2.29	0.46
1:H:14:ILE:CD1	1:H:171:THR:HG21	2.43	0.46
1:H:379:SER:O	1:H:383:ASP:HB2	2.16	0.46
1:J:230:TYR:CD2	1:J:269:LEU:HD23	2.51	0.46
1:L:56:ILE:HD11	1:L:397:LEU:HD11	1.97	0.46
1:A:58:HIS:HD2	3:A:647:HOH:O	1.97	0.46
1:A:162:SER:OG	1:A:316:VAL:HG23	2.16	0.46
1:A:193:LEU:HA	1:A:193:LEU:HD23	1.74	0.46
1:A:218:ALA:HA	1:A:270:TYR:CD2	2.51	0.46
1:B:165:VAL:HG12	1:B:166:GLY:N	2.31	0.46
1:B:248:ILE:HG22	3:B:685:HOH:O	2.16	0.46
1:D:217:MET:O	1:D:270:TYR:HD2	1.98	0.46
1:F:338:VAL:HG11	3:F:666:HOH:O	2.14	0.46
1:G:222:GLU:HB2	1:G:223:LEU:HD12	1.96	0.46
1:J:70:GLN:HB2	1:J:82:GLN:HB3	1.97	0.46
1:H:95:GLU:HB2	1:H:140:LYS:NZ	2.30	0.46
1:I:53:TYR:O	1:I:55:VAL:HG22	2.15	0.46
1:J:316:VAL:HG13	1:J:320:TYR:CD2	2.49	0.46
1:J:351:ARG:NE	1:J:355:ASP:O	2.43	0.46
1:A:14:ILE:HB	1:A:17:LEU:HD12	1.98	0.46
1:C:365:THR:HG23	1:C:367:ASP:HB2	1.98	0.46
1:E:19:VAL:HG21	1:F:360:ASN:HD21	1.81	0.46
1:E:49:TYR:OH	1:E:93:THR:O	2.23	0.46
1:E:307:THR:OG1	1:E:403:LEU:HB3	2.16	0.46
1:F:181:GLU:OE2	1:F:185:ASN:ND2	2.38	0.46
1:F:291:ARG:HG3	1:F:302:MET:HG2	1.97	0.46
1:H:340:PHE:O	1:H:364:ILE:N	2.48	0.46
1:J:365:THR:C	1:J:367:ASP:N	2.74	0.46
1:C:340:PHE:HB2	1:C:364:ILE:HB	1.98	0.46
1:F:205:LEU:O	1:F:211:VAL:HA	2.16	0.46
1:G:60:ASN:C	1:G:60:ASN:OD1	2.58	0.46
1:J:235:GLU:O	1:J:239:ASN:ND2	2.47	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:223:LEU:HD13	3:K:669:HOH:O	2.15	0.46
1:L:114:SER:HB2	3:L:667:HOH:O	2.15	0.46
1:A:58:HIS:CE1	3:A:646:HOH:O	2.68	0.46
1:B:301:PRO:HD2	3:B:671:HOH:O	2.15	0.46
1:H:217:MET:HE1	1:H:237:VAL:HG21	1.97	0.46
1:H:403:LEU:HD13	3:H:631:HOH:O	2.16	0.46
1:L:373:GLU:HG2	1:L:374:THR:N	2.29	0.46
1:E:70:GLN:HA	1:E:280:HIS:CD2	2.50	0.45
1:E:223:LEU:N	1:E:223:LEU:HD12	2.31	0.45
1:E:310:LYS:HE2	1:E:315:HIS:NE2	2.30	0.45
1:F:365:THR:C	1:F:367:ASP:H	2.24	0.45
1:I:20:GLU:OE1	1:I:206:ARG:NE	2.45	0.45
1:I:68:GLY:HA2	1:I:278:ALA:O	2.16	0.45
1:I:217:MET:HE1	1:I:237:VAL:CG1	2.33	0.45
1:L:328:ALA:HB2	1:L:343:VAL:HG23	1.98	0.45
1:B:116:LYS:HB3	3:B:654:HOH:O	2.16	0.45
1:C:238:ARG:HB2	1:C:256:VAL:HG23	1.99	0.45
1:C:304:LEU:HD23	1:C:304:LEU:HA	1.69	0.45
1:D:143:SER:HA	1:D:146:VAL:HG12	1.99	0.45
1:D:291:ARG:HG3	1:D:302:MET:HG2	1.96	0.45
1:E:181:GLU:OE1	1:E:182:ARG:HD2	2.16	0.45
1:E:238:ARG:HB2	1:E:256:VAL:CG2	2.46	0.45
1:F:310:LYS:HB2	1:F:318:LYS:HB2	1.98	0.45
1:G:14:ILE:HB	1:G:17:LEU:HD12	1.98	0.45
1:J:14:ILE:CD1	1:J:171:THR:HG21	2.36	0.45
1:A:17:LEU:O	1:A:206:ARG:NH2	2.39	0.45
1:E:8:LEU:HD12	1:E:363:VAL:HG13	1.97	0.45
1:E:346:LEU:HA	1:E:346:LEU:HD23	1.71	0.45
1:F:61:LEU:HD23	1:F:90:GLY:HA3	1.99	0.45
1:J:339:LYS:HB3	1:J:364:ILE:HG22	1.97	0.45
1:K:8:LEU:HD11	1:K:365:THR:HG22	1.99	0.45
1:L:199:ASP:HB2	1:L:276:THR:HA	1.97	0.45
1:L:212:ASP:OD1	1:L:255:ARG:HD2	2.17	0.45
1:L:296:ILE:HG22	1:L:298:PRO:HD3	1.97	0.45
1:B:323:LEU:HB2	1:B:392:LEU:HD21	1.98	0.45
1:D:15:GLU:HG3	1:D:171:THR:HB	1.98	0.45
1:D:345:VAL:HG22	3:D:662:HOH:O	2.15	0.45
1:E:291:ARG:HA	1:E:291:ARG:NE	2.32	0.45
1:G:238:ARG:HB2	1:G:256:VAL:HG23	1.99	0.45
1:G:304:LEU:HD12	1:H:296:ILE:HG23	1.98	0.45
1:I:21:LEU:O	1:I:298:PRO:HD2	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:63:LYS:HE2	1:J:283:ASP:HB2	1.98	0.45
1:I:283:ASP:OD2	1:J:90:GLY:HA2	2.16	0.45
1:J:56:ILE:HD11	1:J:397:LEU:HD11	1.99	0.45
1:K:4:ILE:HG21	1:K:379:SER:HB2	1.99	0.45
1:K:49:TYR:OH	1:K:93:THR:O	2.22	0.45
1:A:60:ASN:ND2	1:A:60:ASN:C	2.68	0.45
1:A:288:ARG:NH2	1:B:62:ASP:O	2.49	0.45
1:B:70:GLN:NE2	3:C:657:HOH:O	2.50	0.45
1:C:230:TYR:OH	1:C:262:ASP:OD2	2.25	0.45
1:D:120:ARG:NH2	1:D:129:GLU:OE2	2.50	0.45
1:D:356:PRO:HG2	1:D:386:LEU:HD21	1.97	0.45
1:K:20:GLU:HG2	1:K:173:LEU:HD23	1.99	0.45
1:K:319:LEU:HD23	1:K:319:LEU:HA	1.80	0.45
1:L:305:GLU:HB2	3:L:623:HOH:O	2.15	0.45
1:B:220:ILE:HB	1:B:223:LEU:HD13	1.98	0.45
1:B:360:ASN:OD1	1:B:361:VAL:N	2.50	0.45
1:E:207:ARG:NE	1:F:5:ASN:HD22	2.14	0.45
1:E:316:VAL:HG23	1:E:317:GLY:N	2.31	0.45
1:F:357:LEU:HB2	3:F:651:HOH:O	2.16	0.45
1:L:238:ARG:HB2	1:L:256:VAL:CG2	2.46	0.45
1:B:10:PRO:HA	1:B:364:ILE:HD11	1.98	0.45
1:C:346:LEU:CD1	1:D:21:LEU:HD11	2.47	0.45
1:E:210:GLU:HG2	1:E:253:ASN:HB3	1.98	0.45
1:G:68:GLY:HA2	1:G:278:ALA:O	2.17	0.45
1:H:112:ILE:HD11	1:H:136:TYR:HD1	1.82	0.45
1:H:120:ARG:NH2	1:H:129:GLU:OE2	2.49	0.45
1:H:207:ARG:O	1:H:210:GLU:HB2	2.17	0.45
1:H:322:VAL:O	1:H:326:LEU:HD13	2.17	0.45
1:J:164:GLY:HA2	1:J:321:ASN:ND2	2.32	0.45
1:A:160:ASP:HA	3:A:706:HOH:O	2.16	0.45
1:A:201:LYS:HE3	3:B:636:HOH:O	2.16	0.45
1:B:320:TYR:HE2	1:B:347:GLY:CA	2.30	0.45
1:C:197:GLY:HA3	1:C:218:ALA:HB3	1.99	0.45
1:H:35:ASP:HA	1:H:286:THR:HG22	1.99	0.45
1:I:268:ILE:C	1:I:269:LEU:HD12	2.42	0.45
1:B:157:LEU:N	1:B:157:LEU:HD12	2.32	0.45
1:B:344:GLN:HG3	3:B:612:HOH:O	2.17	0.45
1:B:398:GLU:HB2	3:B:620:HOH:O	2.17	0.45
1:E:70:GLN:HG2	1:E:280:HIS:CD2	2.52	0.45
1:G:64:THR:HB	1:G:286:THR:HG23	1.99	0.45
1:H:195:GLU:OE1	1:H:195:GLU:N	2.45	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:112:ILE:HD11	1:B:136:TYR:HD1	1.82	0.45
1:C:290:ASN:HB2	1:C:404:PHE:O	2.17	0.45
1:G:281:GLY:CA	1:H:91:ARG:HE	2.30	0.45
1:H:181:GLU:OE1	1:H:182:ARG:HD2	2.17	0.45
1:A:298:PRO:O	1:B:301:PRO:HB3	2.17	0.44
1:B:104:GLN:HA	3:B:659:HOH:O	2.17	0.44
1:C:14:ILE:CD1	1:C:171:THR:HG21	2.39	0.44
1:F:329:ASN:O	1:F:333:GLN:HG3	2.17	0.44
1:G:14:ILE:CD1	1:G:171:THR:HG21	2.45	0.44
1:G:70:GLN:HB2	1:G:82:GLN:HB3	1.99	0.44
1:H:22:VAL:HG11	1:H:174:GLU:HA	1.99	0.44
1:H:399:GLY:HA2	3:H:674:HOH:O	2.16	0.44
1:I:190:LYS:HE3	1:I:196:VAL:O	2.17	0.44
1:J:169:PRO:HD2	1:J:340:PHE:HD1	1.82	0.44
1:J:321:ASN:ND2	3:J:619:HOH:O	2.50	0.44
1:B:351:ARG:NE	1:B:355:ASP:O	2.43	0.44
1:E:25:LYS:NZ	2:F:501:PO4:O2	2.46	0.44
1:H:38:ALA:HB1	1:H:61:LEU:O	2.17	0.44
1:I:374:THR:HG21	3:I:629:HOH:O	2.16	0.44
1:K:199:ASP:OD2	1:K:277:SER:OG	2.33	0.44
1:L:117:GLU:OE1	1:L:120:ARG:NH1	2.49	0.44
1:B:167:PHE:HB3	1:B:342:GLN:HB2	1.98	0.44
1:C:60:ASN:ND2	1:C:60:ASN:O	2.45	0.44
1:C:283:ASP:O	1:D:63:LYS:HE2	2.17	0.44
1:D:161:THR:CG2	1:D:349:ILE:H	2.30	0.44
1:I:15:GLU:HG3	1:I:171:THR:HB	2.00	0.44
1:I:217:MET:CE	1:I:237:VAL:HG11	2.34	0.44
1:J:6:VAL:CG1	1:J:378:ILE:HD11	2.48	0.44
1:J:162:SER:HB2	1:J:305:GLU:OE1	2.17	0.44
1:K:207:ARG:NH2	1:L:5:ASN:HB2	2.32	0.44
1:L:224:ILE:HG23	3:L:662:HOH:O	2.16	0.44
1:A:339:LYS:HG3	1:A:366:TYR:CE1	2.52	0.44
1:B:311:ASN:HB3	1:B:315:HIS:HB2	1.98	0.44
1:C:60:ASN:HD22	1:C:60:ASN:C	2.26	0.44
1:E:330:LYS:HD3	1:E:381:ILE:HG23	1.98	0.44
1:J:249:ALA:O	1:J:252:TYR:HB2	2.17	0.44
1:K:157:LEU:HD21	1:K:350:GLY:HA2	1.98	0.44
1:A:231:ILE:HD11	1:A:264:ILE:HD11	1.99	0.44
1:D:95:GLU:HG2	1:D:104:GLN:OE1	2.17	0.44
1:D:342:GLN:O	3:D:619:HOH:O	2.21	0.44
1:F:335:VAL:O	1:F:338:VAL:HG22	2.16	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:365:THR:O	1:J:367:ASP:N	2.39	0.44
1:D:199:ASP:HB2	1:D:276:THR:HA	1.98	0.44
1:E:83:PRO:O	3:E:622:HOH:O	2.21	0.44
1:F:93:THR:HA	3:F:684:HOH:O	2.17	0.44
1:F:178:TYR:HB2	1:F:295:LEU:HD21	1.99	0.44
1:A:22:VAL:HG11	1:A:174:GLU:HA	2.00	0.44
1:A:67:VAL:HG13	3:A:694:HOH:O	2.18	0.44
1:C:273:VAL:H	1:C:279:GLU:CD	2.26	0.44
1:F:165:VAL:HG12	1:F:166:GLY:N	2.33	0.44
1:F:181:GLU:OE1	1:F:182:ARG:HD2	2.17	0.44
1:F:268:ILE:C	1:F:269:LEU:HD12	2.42	0.44
1:G:103:ASP:OD1	1:G:103:ASP:N	2.39	0.44
1:H:69:GLY:HA3	3:H:659:HOH:O	2.18	0.44
1:H:91:ARG:HG2	1:H:141:GLY:HA2	2.00	0.44
1:H:126:LEU:HD13	1:H:274:THR:HB	1.99	0.44
1:A:335:VAL:O	1:A:338:VAL:HG22	2.18	0.44
1:B:136:TYR:HD2	1:B:138:ILE:HG12	1.82	0.44
1:C:35:ASP:CG	1:C:287:GLY:H	2.21	0.44
1:I:43:ARG:NH1	1:I:398:GLU:HA	2.33	0.44
1:I:69:GLY:O	1:I:70:GLN:HG3	2.17	0.44
1:A:5:ASN:N	3:A:663:HOH:O	2.50	0.44
1:C:60:ASN:ND2	1:C:60:ASN:C	2.76	0.44
1:D:91:ARG:HG2	1:D:141:GLY:HA2	2.00	0.44
1:G:318:LYS:O	1:G:322:VAL:HG23	2.18	0.44
1:G:351:ARG:HD3	1:G:355:ASP:O	2.17	0.44
1:H:14:ILE:HB	1:H:17:LEU:HD12	2.00	0.44
1:H:161:THR:HG22	1:H:348:GLN:HA	1.99	0.44
1:H:206:ARG:HD2	1:H:252:TYR:OH	2.18	0.44
1:I:270:TYR:O	1:I:270:TYR:HD2	2.01	0.44
1:L:103:ASP:OD1	1:L:103:ASP:N	2.46	0.44
1:L:142:SER:HB3	1:L:145:LEU:HD13	2.00	0.44
1:A:63:LYS:HE2	1:B:283:ASP:HB2	2.00	0.43
1:B:38:ALA:HB1	1:B:61:LEU:O	2.18	0.43
1:G:385:MET:HE2	1:G:385:MET:HB3	1.56	0.43
1:H:222:GLU:HB2	1:H:223:LEU:HD12	2.00	0.43
1:I:18:GLN:HB3	1:I:19:VAL:HG23	2.00	0.43
1:L:53:TYR:CZ	1:L:96:VAL:HG21	2.53	0.43
1:B:84:ILE:HA	3:B:696:HOH:O	2.17	0.43
1:B:110:ILE:HD12	3:B:691:HOH:O	2.17	0.43
1:B:325:ASN:HB2	3:B:627:HOH:O	2.17	0.43
1:D:195:GLU:HA	3:D:677:HOH:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:245:ALA:CB	1:D:254:VAL:HG21	2.47	0.43
1:F:217:MET:HE1	1:F:237:VAL:HG21	2.00	0.43
1:G:25:LYS:HD2	1:G:289:GLY:HA2	2.00	0.43
1:H:44:LYS:HE3	3:H:636:HOH:O	2.18	0.43
1:H:76:LYS:HG3	1:H:225:GLU:HA	1.99	0.43
1:J:7:GLN:O	1:J:363:VAL:HG12	2.19	0.43
1:L:395:LEU:HA	1:L:398:GLU:OE2	2.18	0.43
1:B:307:THR:OG1	1:B:403:LEU:HB3	2.17	0.43
1:D:50:LEU:HD13	1:D:56:ILE:CG1	2.46	0.43
1:D:300:ARG:HG3	3:D:654:HOH:O	2.18	0.43
1:E:35:ASP:OD1	1:E:287:GLY:N	2.32	0.43
1:G:323:LEU:HB2	1:G:392:LEU:HD21	1.99	0.43
1:H:311:ASN:HB2	3:H:603:HOH:O	2.18	0.43
1:I:348:GLN:HB2	1:I:357:LEU:HD12	2.00	0.43
1:J:360:ASN:OD1	1:J:361:VAL:N	2.51	0.43
1:K:257:TYR:HB3	1:K:260:THR:CG2	2.48	0.43
1:B:373:GLU:HG2	1:B:374:THR:N	2.34	0.43
1:C:67:VAL:HG13	3:D:644:HOH:O	2.18	0.43
1:D:217:MET:O	1:D:270:TYR:CD2	2.71	0.43
1:E:315:HIS:O	1:E:319:LEU:HG	2.18	0.43
1:F:212:ASP:OD1	1:F:255:ARG:HD2	2.18	0.43
1:F:312:PRO:HB2	3:F:657:HOH:O	2.18	0.43
1:G:203:MET:HE3	1:H:346:LEU:HD11	2.00	0.43
1:I:357:LEU:O	1:I:357:LEU:HD23	2.18	0.43
1:J:241:ILE:O	1:J:245:ALA:N	2.43	0.43
1:A:11:LEU:HD23	1:B:11:LEU:HG	2.01	0.43
1:A:296:ILE:HG12	1:B:304:LEU:HD12	2.00	0.43
1:B:193:LEU:HB3	1:B:196:VAL:HG23	2.00	0.43
1:B:290:ASN:ND2	1:B:404:PHE:O	2.35	0.43
1:C:19:VAL:HG21	1:D:360:ASN:HD21	1.80	0.43
1:E:56:ILE:HD12	1:E:313:VAL:HG23	2.00	0.43
1:F:22:VAL:HB	1:F:177:VAL:HG21	2.00	0.43
1:F:120:ARG:NH2	1:F:129:GLU:OE2	2.52	0.43
1:G:47:LEU:HD23	1:G:47:LEU:HA	1.74	0.43
1:H:29:HIS:HA	1:H:30:PRO:HD2	1.82	0.43
1:I:163:PHE:CZ	1:I:165:VAL:HG22	2.53	0.43
1:I:391:LYS:HD2	1:I:391:LYS:HA	1.78	0.43
1:J:165:VAL:HG23	1:J:303:SER:HB2	2.00	0.43
1:C:205:LEU:HD12	1:C:206:ARG:N	2.33	0.43
1:C:210:GLU:HG2	1:C:253:ASN:HB3	2.00	0.43
1:E:160:ASP:OD1	1:E:161:THR:N	2.50	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:34:ALA:C	1:F:286:THR:HG22	2.43	0.43
1:F:304:LEU:HB2	3:F:658:HOH:O	2.18	0.43
1:J:340:PHE:O	1:J:364:ILE:N	2.49	0.43
1:K:375:LYS:HG3	3:K:648:HOH:O	2.18	0.43
1:A:281:GLY:H	1:B:91:ARG:HE	1.65	0.43
1:C:299:MET:HE3	1:D:344:GLN:HE22	1.83	0.43
1:D:6:VAL:HG11	1:D:378:ILE:HD11	1.99	0.43
1:F:25:LYS:HD2	1:F:289:GLY:HA2	2.01	0.43
1:G:10:PRO:HA	1:G:364:ILE:HD11	2.01	0.43
1:G:48:TYR:OH	3:G:617:HOH:O	2.14	0.43
1:H:81:ILE:HD12	1:H:81:ILE:HG23	1.80	0.43
1:I:319:LEU:HD23	1:I:319:LEU:HA	1.88	0.43
1:J:23:GLU:HB3	1:J:296:ILE:HB	2.01	0.43
1:J:167:PHE:CZ	1:J:293:VAL:HG11	2.54	0.43
1:J:213:LEU:HD23	1:J:256:VAL:HG12	2.01	0.43
1:K:60:ASN:O	1:K:90:GLY:HA3	2.19	0.43
1:L:120:ARG:NH2	1:L:129:GLU:OE2	2.52	0.43
1:A:70:GLN:HA	1:A:280:HIS:CD2	2.53	0.43
1:A:207:ARG:NH2	1:B:5:ASN:HB2	2.34	0.43
1:C:373:GLU:HG2	1:C:374:THR:N	2.32	0.43
1:E:47:LEU:HA	1:E:47:LEU:HD23	1.78	0.43
1:G:120:ARG:NH2	1:G:129:GLU:OE2	2.52	0.43
1:H:104:GLN:HA	3:H:658:HOH:O	2.18	0.43
1:J:338:VAL:HA	1:J:365:THR:HG22	2.01	0.43
1:K:216:ALA:HB2	3:L:635:HOH:O	2.17	0.43
1:L:312:PRO:O	1:L:393:THR:OG1	2.28	0.43
1:A:80:ILE:HG13	1:A:131:HIS:CD2	2.54	0.43
1:B:15:GLU:HG3	1:B:171:THR:HB	2.01	0.43
1:C:283:ASP:OD2	1:D:90:GLY:HA2	2.19	0.43
1:E:75:PHE:CZ	1:E:227:VAL:HG22	2.53	0.43
1:E:257:TYR:CD2	1:F:348:GLN:NE2	2.85	0.43
1:E:271:LEU:HD23	3:E:712:HOH:O	2.17	0.43
1:E:315:HIS:NE2	3:E:676:HOH:O	2.19	0.43
1:G:310:LYS:HE2	1:G:315:HIS:NE2	2.33	0.43
1:H:163:PHE:CE1	1:H:165:VAL:HG22	2.54	0.43
1:I:20:GLU:HG2	1:I:173:LEU:HD23	2.00	0.43
1:K:227:VAL:HG12	1:K:264:ILE:HD11	2.01	0.43
1:K:329:ASN:O	1:K:333:GLN:HG3	2.19	0.43
1:L:230:TYR:CD2	1:L:269:LEU:HD23	2.54	0.43
1:L:318:LYS:O	1:L:322:VAL:HG23	2.19	0.43
1:A:58:HIS:CD2	3:A:647:HOH:O	2.72	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:91:ARG:HH21	1:B:282:ASP:N	2.16	0.43
1:B:53:TYR:CZ	1:B:96:VAL:HG21	2.54	0.43
1:E:245:ALA:HB1	1:E:254:VAL:HG21	2.01	0.43
1:F:360:ASN:OD1	1:F:361:VAL:N	2.52	0.43
1:K:14:ILE:HB	1:K:17:LEU:HD12	2.01	0.43
1:K:238:ARG:HB2	1:K:256:VAL:HG21	2.01	0.43
1:K:268:ILE:C	1:K:269:LEU:HD12	2.44	0.43
1:L:91:ARG:HG2	1:L:141:GLY:HA2	2.00	0.43
1:C:158:SER:C	1:C:160:ASP:H	2.27	0.42
1:D:65:LEU:C	3:D:651:HOH:O	2.60	0.42
1:G:304:LEU:HD23	1:G:304:LEU:HA	1.69	0.42
1:H:357:LEU:HB2	3:H:666:HOH:O	2.19	0.42
1:H:374:THR:HA	3:H:644:HOH:O	2.17	0.42
1:K:136:TYR:HD2	1:K:138:ILE:HG12	1.84	0.42
1:K:163:PHE:HE1	1:K:344:GLN:HB3	1.84	0.42
1:L:51:LYS:NZ	3:L:651:HOH:O	2.49	0.42
1:L:266:LYS:HA	1:L:266:LYS:HD3	1.81	0.42
1:D:346:LEU:HA	1:D:346:LEU:HD12	1.66	0.42
1:F:6:VAL:HB	1:F:375:LYS:HE3	2.00	0.42
1:G:22:VAL:HG21	1:G:174:GLU:HG2	2.01	0.42
1:I:112:ILE:HD12	1:I:134:VAL:HB	2.00	0.42
1:L:273:VAL:H	1:L:279:GLU:CD	2.27	0.42
1:L:294:GLY:HA3	3:L:642:HOH:O	2.17	0.42
1:A:125:TYR:HB2	1:A:274:THR:O	2.19	0.42
1:D:7:GLN:O	1:D:363:VAL:HG12	2.18	0.42
1:D:27:LEU:HA	1:D:32:TYR:CE1	2.55	0.42
1:E:19:VAL:HG11	1:F:360:ASN:HD22	1.85	0.42
1:F:23:GLU:HB3	1:F:296:ILE:HB	2.01	0.42
1:F:230:TYR:CD2	1:F:269:LEU:HD23	2.54	0.42
1:K:342:GLN:HG3	3:K:638:HOH:O	2.19	0.42
1:L:49:TYR:HB3	1:L:55:VAL:O	2.19	0.42
1:A:207:ARG:NE	1:B:5:ASN:HD22	2.15	0.42
1:B:163:PHE:CZ	1:B:165:VAL:HG22	2.55	0.42
1:C:76:LYS:HG3	1:C:225:GLU:HA	2.01	0.42
1:C:281:GLY:H	1:D:91:ARG:HE	1.66	0.42
1:C:357:LEU:HD23	1:C:357:LEU:O	2.20	0.42
1:D:27:LEU:HG	3:D:678:HOH:O	2.18	0.42
1:D:127:ASP:OD1	1:D:130:ARG:HG3	2.20	0.42
1:G:29:HIS:HA	1:G:30:PRO:HD2	1.91	0.42
1:H:25:LYS:HD2	1:H:289:GLY:HA2	2.00	0.42
1:H:311:ASN:ND2	1:H:314:ASN:OD1	2.53	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:235:GLU:O	1:I:239:ASN:ND2	2.52	0.42
1:J:263:LYS:HB2	1:J:268:ILE:HG13	2.01	0.42
1:K:205:LEU:HD12	1:K:206:ARG:N	2.35	0.42
1:L:354:ASP:HB2	1:L:386:LEU:O	2.19	0.42
1:A:12:SER:O	3:A:666:HOH:O	2.21	0.42
1:A:143:SER:HA	3:A:697:HOH:O	2.18	0.42
1:A:222:GLU:HG3	3:A:683:HOH:O	2.18	0.42
1:F:106:PRO:O	1:F:110:ILE:HG13	2.19	0.42
1:F:222:GLU:HB2	1:F:223:LEU:HD12	2.02	0.42
1:G:291:ARG:NH1	1:G:304:LEU:HA	2.35	0.42
1:I:346:LEU:HD11	1:J:21:LEU:HD11	2.01	0.42
1:J:351:ARG:NH2	1:J:355:ASP:OD2	2.52	0.42
1:K:190:LYS:NZ	3:K:642:HOH:O	2.52	0.42
1:B:164:GLY:O	3:B:650:HOH:O	2.22	0.42
1:C:340:PHE:O	1:C:364:ILE:N	2.52	0.42
1:F:161:THR:CG2	1:F:349:ILE:H	2.30	0.42
1:J:103:ASP:OD1	1:J:103:ASP:N	2.41	0.42
1:A:231:ILE:HD11	1:A:264:ILE:HG12	2.02	0.42
1:C:22:VAL:HG11	1:C:174:GLU:HA	2.01	0.42
1:E:290:ASN:HB2	1:E:404:PHE:O	2.20	0.42
1:E:391:LYS:HD2	1:E:394:GLU:HB3	2.02	0.42
1:F:2:ARG:HG3	3:F:702:HOH:O	2.20	0.42
1:F:276:THR:C	1:F:278:ALA:H	2.28	0.42
1:H:164:GLY:HA2	1:H:321:ASN:ND2	2.35	0.42
1:I:270:TYR:CD1	3:J:633:HOH:O	2.56	0.42
1:I:285:MET:HE1	1:J:64:THR:O	2.20	0.42
1:J:357:LEU:HD23	1:J:357:LEU:O	2.20	0.42
1:J:382:VAL:O	1:J:386:LEU:HD13	2.20	0.42
1:A:231:ILE:HD11	1:A:264:ILE:CD1	2.50	0.42
1:B:365:THR:C	1:B:367:ASP:N	2.78	0.42
1:C:52:LYS:HB3	1:C:52:LYS:HE3	1.86	0.42
1:C:53:TYR:O	1:C:55:VAL:HG22	2.20	0.42
1:C:207:ARG:HD2	1:D:358:ILE:HD12	2.02	0.42
1:D:64:THR:HB	1:D:286:THR:HG23	2.02	0.42
1:E:38:ALA:HB2	1:E:64:THR:OG1	2.20	0.42
1:E:385:MET:HE2	1:E:385:MET:HB3	1.58	0.42
1:F:161:THR:HG23	1:F:349:ILE:HG22	2.01	0.42
1:I:106:PRO:O	1:I:110:ILE:HG13	2.20	0.42
1:I:200:ILE:HA	1:I:216:ALA:O	2.19	0.42
1:J:171:THR:HG23	1:J:300:ARG:HH21	1.84	0.42
1:J:212:ASP:OD1	1:J:255:ARG:HD2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:224:ILE:HG21	1:J:271:LEU:HD21	2.02	0.42
1:J:385:MET:HE2	1:J:385:MET:HB3	1.81	0.42
1:A:53:TYR:O	1:A:55:VAL:HG22	2.20	0.42
1:B:268:ILE:C	1:B:269:LEU:HD12	2.44	0.42
1:D:178:TYR:HE2	3:D:647:HOH:O	2.02	0.42
1:D:296:ILE:HG22	1:D:298:PRO:HD3	2.00	0.42
1:E:7:GLN:O	1:E:363:VAL:HG12	2.20	0.42
1:F:249:ALA:O	1:F:252:TYR:HB2	2.20	0.42
1:G:160:ASP:OD1	1:G:161:THR:N	2.52	0.42
1:H:165:VAL:HG12	1:H:166:GLY:N	2.35	0.42
1:I:257:TYR:HB3	1:I:260:THR:CG2	2.50	0.42
1:J:25:LYS:HD2	1:J:289:GLY:HA2	2.01	0.42
1:L:142:SER:HB3	1:L:145:LEU:CD1	2.50	0.42
1:A:158:SER:HA	1:A:316:VAL:HG11	2.02	0.42
1:A:206:ARG:HD2	1:A:252:TYR:CZ	2.55	0.42
1:B:375:LYS:HE2	1:B:375:LYS:HB3	1.76	0.42
1:D:39:GLU:HA	1:D:309:GLY:HA2	2.02	0.42
1:D:63:LYS:N	3:D:641:HOH:O	2.52	0.42
1:D:64:THR:HB	1:D:286:THR:CG2	2.50	0.42
1:D:218:ALA:HA	1:D:270:TYR:HD2	1.84	0.42
1:D:318:LYS:O	1:D:322:VAL:HG23	2.20	0.42
1:E:130:ARG:NE	3:E:699:HOH:O	2.53	0.42
1:E:203:MET:HE3	1:F:346:LEU:HD11	2.02	0.42
1:G:9:ASN:O	1:G:12:SER:HB3	2.20	0.42
1:G:20:GLU:C	1:G:21:LEU:HD12	2.45	0.42
1:G:281:GLY:H	1:H:91:ARG:HE	1.68	0.42
1:G:378:ILE:HG13	1:G:379:SER:N	2.34	0.42
1:H:22:VAL:HB	1:H:177:VAL:HG21	2.02	0.42
1:H:23:GLU:HB3	1:H:296:ILE:HB	2.02	0.42
1:L:22:VAL:HB	1:L:177:VAL:HG21	2.00	0.42
1:A:224:ILE:HG21	1:A:271:LEU:HD21	2.01	0.41
1:A:356:PRO:HG2	1:A:386:LEU:HD21	2.02	0.41
1:C:126:LEU:HD12	1:C:126:LEU:HA	1.85	0.41
1:D:8:LEU:HD12	1:D:363:VAL:HG13	2.00	0.41
1:D:354:ASP:HA	1:D:389:PHE:HE2	1.84	0.41
1:E:50:LEU:HD13	1:E:56:ILE:HG13	2.02	0.41
1:F:230:TYR:CG	1:F:269:LEU:HD23	2.54	0.41
1:F:305:GLU:HB2	3:F:656:HOH:O	2.19	0.41
1:G:62:ASP:OD2	1:G:310:LYS:NZ	2.43	0.41
1:G:70:GLN:HA	1:G:280:HIS:CD2	2.55	0.41
1:G:299:MET:HE3	1:H:344:GLN:OE1	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:11:LEU:HD23	1:J:11:LEU:HG	2.01	0.41
1:I:19:VAL:HG11	1:J:360:ASN:HD22	1.85	0.41
1:I:207:ARG:HE	1:J:5:ASN:ND2	2.16	0.41
1:J:38:ALA:HB1	1:J:61:LEU:O	2.20	0.41
1:K:291:ARG:HA	1:K:291:ARG:NE	2.35	0.41
1:A:273:VAL:H	1:A:279:GLU:CD	2.28	0.41
1:B:80:ILE:HG13	1:B:131:HIS:CD2	2.55	0.41
1:C:335:VAL:O	1:C:338:VAL:HG22	2.21	0.41
1:D:22:VAL:HB	1:D:177:VAL:HG21	2.02	0.41
1:D:76:LYS:HG3	1:D:225:GLU:HA	2.01	0.41
1:D:163:PHE:CE1	1:D:165:VAL:HG22	2.55	0.41
1:D:213:LEU:HB3	1:D:256:VAL:HG12	2.02	0.41
1:D:238:ARG:HB2	1:D:256:VAL:CG2	2.50	0.41
1:E:365:THR:C	1:E:367:ASP:H	2.28	0.41
1:F:15:GLU:CG	1:F:171:THR:HB	2.51	0.41
1:F:375:LYS:HE2	1:F:375:LYS:HB3	1.81	0.41
1:H:74:ARG:HD3	3:H:621:HOH:O	2.20	0.41
1:H:165:VAL:HG23	1:H:303:SER:HB2	2.02	0.41
1:H:223:LEU:HD12	1:H:223:LEU:N	2.35	0.41
1:H:270:TYR:OH	1:H:277:SER:HB3	2.20	0.41
1:H:354:ASP:HA	1:H:389:PHE:CE2	2.54	0.41
1:K:245:ALA:HB1	1:K:254:VAL:HG21	2.01	0.41
1:L:281:GLY:N	3:L:632:HOH:O	2.53	0.41
1:A:19:VAL:HG21	1:B:360:ASN:HD21	1.83	0.41
1:B:373:GLU:H	1:B:373:GLU:CD	2.25	0.41
1:C:222:GLU:HG3	3:C:683:HOH:O	2.20	0.41
1:D:165:VAL:HG12	1:D:166:GLY:N	2.35	0.41
1:D:357:LEU:O	1:D:357:LEU:HD23	2.20	0.41
1:E:91:ARG:HH21	1:F:282:ASP:N	2.18	0.41
1:H:71:ALA:HB2	1:H:80:ILE:HD13	2.02	0.41
1:H:199:ASP:HB2	1:H:276:THR:HA	2.01	0.41
1:I:85:TYR:CD2	1:J:137:LYS:HD2	2.55	0.41
1:J:90:GLY:O	1:J:91:ARG:HD2	2.20	0.41
1:J:222:GLU:HB2	1:J:223:LEU:HD12	2.02	0.41
1:K:57:LEU:HD22	1:K:93:THR:HG21	2.01	0.41
1:C:319:LEU:HD23	1:C:319:LEU:HA	1.78	0.41
1:D:165:VAL:HG23	1:D:303:SER:HB2	2.02	0.41
1:E:53:TYR:CZ	1:E:96:VAL:HG21	2.54	0.41
1:E:80:ILE:HG13	1:E:131:HIS:CD2	2.55	0.41
1:H:48:TYR:O	1:H:49:TYR:C	2.64	0.41
1:H:95:GLU:HG2	3:H:658:HOH:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:373:GLU:HG2	1:H:374:THR:N	2.34	0.41
1:J:218:ALA:HA	1:J:270:TYR:CD2	2.55	0.41
1:K:385:MET:HB3	1:K:385:MET:HE2	1.56	0.41
1:L:9:ASN:HA	1:L:10:PRO:HD3	1.77	0.41
1:L:183:HIS:O	1:L:186:SER:HB3	2.20	0.41
1:L:276:THR:C	1:L:278:ALA:H	2.28	0.41
1:B:120:ARG:NH2	3:B:654:HOH:O	2.28	0.41
1:C:217:MET:HE1	1:C:237:VAL:HG21	2.03	0.41
1:D:34:ALA:C	1:D:286:THR:HG22	2.45	0.41
1:D:49:TYR:O	1:D:52:LYS:O	2.39	0.41
1:F:268:ILE:HD13	1:F:268:ILE:HG21	1.84	0.41
1:H:170:LEU:HB3	1:H:174:GLU:HB2	2.01	0.41
1:H:274:THR:N	1:H:279:GLU:OE1	2.50	0.41
1:H:328:ALA:HB2	1:H:343:VAL:HG23	2.02	0.41
1:H:360:ASN:OD1	1:H:361:VAL:N	2.54	0.41
1:K:71:ALA:N	1:K:280:HIS:HD2	2.03	0.41
1:L:237:VAL:O	1:L:240:GLN:HB2	2.20	0.41
1:B:310:LYS:O	1:B:318:LYS:NZ	2.39	0.41
1:C:355:ASP:N	1:C:356:PRO:HD3	2.33	0.41
1:F:33:ILE:O	1:F:37:VAL:HG22	2.20	0.41
1:F:53:TYR:O	1:F:55:VAL:N	2.52	0.41
1:F:209:ASN:O	1:F:253:ASN:N	2.43	0.41
1:F:305:GLU:HG3	3:F:627:HOH:O	2.21	0.41
1:F:394:GLU:HG3	3:F:680:HOH:O	2.19	0.41
1:G:281:GLY:H	1:H:91:ARG:HH21	1.68	0.41
1:H:81:ILE:HD13	1:H:81:ILE:HA	1.70	0.41
1:H:234:LYS:NZ	1:H:259:ASN:HB2	2.35	0.41
1:H:365:THR:O	1:H:367:ASP:N	2.49	0.41
1:J:263:LYS:HB2	1:J:268:ILE:CG1	2.50	0.41
1:K:69:GLY:O	1:K:70:GLN:HG3	2.20	0.41
1:K:212:ASP:OD2	1:L:357:LEU:HD21	2.20	0.41
1:K:222:GLU:HB2	1:K:223:LEU:HD12	2.01	0.41
1:L:161:THR:HG22	1:L:348:GLN:HA	2.03	0.41
1:B:34:ALA:C	1:B:286:THR:HG22	2.44	0.41
1:B:105:ILE:HG23	1:B:105:ILE:HD12	1.71	0.41
1:D:29:HIS:HA	1:D:30:PRO:HD2	1.90	0.41
1:E:199:ASP:OD2	1:E:277:SER:OG	2.23	0.41
1:G:140:LYS:NZ	3:G:638:HOH:O	2.33	0.41
1:H:377:GLU:HB2	3:H:644:HOH:O	2.21	0.41
1:I:9:ASN:O	1:I:12:SER:HB3	2.21	0.41
1:I:201:LYS:NZ	1:J:160:ASP:OD2	2.53	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:238:ARG:HB2	1:I:256:VAL:CG2	2.51	0.41
1:J:310:LYS:HB2	1:J:318:LYS:HB2	2.03	0.41
1:J:329:ASN:O	1:J:333:GLN:HG3	2.20	0.41
1:K:138:ILE:HG22	1:K:139:GLY:O	2.20	0.41
1:K:210:GLU:HG2	1:K:253:ASN:HB3	2.03	0.41
1:L:80:ILE:HD12	1:L:80:ILE:HG23	1.82	0.41
1:L:184:LEU:HB3	1:L:200:ILE:HD13	2.01	0.41
1:L:363:VAL:HG22	1:L:364:ILE:N	2.36	0.41
1:A:15:GLU:HG3	1:A:171:THR:HB	2.01	0.41
1:A:103:ASP:OD1	1:A:103:ASP:N	2.44	0.41
1:D:53:TYR:CZ	1:D:96:VAL:HG21	2.55	0.41
1:E:318:LYS:O	1:E:322:VAL:HG23	2.21	0.41
3:E:650:HOH:O	1:F:288:ARG:HG2	2.19	0.41
1:F:356:PRO:HG2	1:F:386:LEU:HD21	2.03	0.41
1:H:310:LYS:HB2	1:H:318:LYS:HB2	2.02	0.41
1:I:197:GLY:HA3	1:I:218:ALA:CB	2.51	0.41
1:A:53:TYR:CZ	1:A:96:VAL:HG21	2.56	0.41
1:A:68:GLY:HA2	1:A:278:ALA:O	2.21	0.41
1:A:394:GLU:O	1:A:394:GLU:HG2	2.21	0.41
1:B:14:ILE:HB	1:B:17:LEU:HD12	2.02	0.41
1:B:162:SER:O	1:B:346:LEU:HD12	2.21	0.41
1:B:181:GLU:OE1	1:B:182:ARG:HD2	2.21	0.41
1:B:306:ALA:O	1:B:310:LYS:NZ	2.32	0.41
1:C:84:ILE:N	1:C:131:HIS:O	2.42	0.41
1:C:217:MET:HB3	1:C:217:MET:HE2	1.81	0.41
1:C:237:VAL:O	1:C:240:GLN:HB2	2.20	0.41
1:C:297:THR:HA	1:C:298:PRO:HD2	1.93	0.41
1:C:351:ARG:HD3	1:C:355:ASP:O	2.21	0.41
1:E:43:ARG:NH1	1:E:398:GLU:HA	2.36	0.41
1:E:393:THR:C	1:E:395:LEU:N	2.79	0.41
1:F:163:PHE:CE1	1:F:165:VAL:CG2	3.04	0.41
1:G:20:GLU:HA	1:G:299:MET:SD	2.61	0.41
1:G:71:ALA:HB2	1:G:80:ILE:HD13	2.02	0.41
1:G:96:VAL:HG21	3:G:616:HOH:O	2.20	0.41
1:H:193:LEU:HD23	1:H:193:LEU:HA	1.86	0.41
1:I:186:SER:HB3	1:I:189:PHE:HB3	2.02	0.41
1:I:210:GLU:HG2	1:I:253:ASN:HB3	2.01	0.41
1:I:353:ILE:HG13	1:I:389:PHE:CE2	2.56	0.41
1:J:53:TYR:CZ	1:J:96:VAL:HG21	2.56	0.41
1:K:22:VAL:HG21	1:K:174:GLU:CG	2.50	0.41
1:K:292:GLY:O	1:K:402:THR:OG1	2.31	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:8:LEU:HD12	1:L:363:VAL:HG13	2.03	0.41
1:L:29:HIS:HA	1:L:30:PRO:HD2	1.87	0.41
1:L:205:LEU:HD12	1:L:206:ARG:N	2.36	0.41
1:L:320:TYR:HE2	1:L:347:GLY:CA	2.34	0.41
1:A:91:ARG:HH21	1:B:281:GLY:C	2.29	0.41
1:A:257:TYR:HB3	1:A:260:THR:CG2	2.51	0.41
1:B:155:VAL:HA	1:B:156:PRO:HD3	1.87	0.41
1:B:238:ARG:HB2	1:B:256:VAL:CG2	2.51	0.41
1:D:311:ASN:ND2	1:D:314:ASN:OD1	2.54	0.41
1:D:325:ASN:HB2	3:D:630:HOH:O	2.19	0.41
1:E:169:PRO:HD2	1:E:340:PHE:HD1	1.85	0.41
1:E:259:ASN:HA	3:E:714:HOH:O	2.21	0.41
1:E:351:ARG:HG2	1:E:355:ASP:O	2.21	0.41
1:F:90:GLY:O	1:F:91:ARG:HD2	2.20	0.41
1:F:340:PHE:O	1:F:364:ILE:N	2.54	0.41
1:G:260:THR:HG21	1:H:348:GLN:NE2	2.36	0.41
1:K:304:LEU:HD23	1:K:304:LEU:HA	1.81	0.41
1:L:339:LYS:HG3	1:L:366:TYR:CE1	2.56	0.41
1:A:354:ASP:C	1:A:356:PRO:HD3	2.46	0.40
1:C:268:ILE:O	1:C:268:ILE:HG13	2.21	0.40
1:D:218:ALA:HA	1:D:270:TYR:CD2	2.56	0.40
1:D:290:ASN:HB3	1:D:294:GLY:C	2.46	0.40
1:E:264:ILE:HD12	1:E:264:ILE:HA	1.94	0.40
1:H:356:PRO:HG2	1:H:386:LEU:HD21	2.03	0.40
1:I:356:PRO:HG2	1:I:386:LEU:HD21	2.04	0.40
1:A:71:ALA:HB2	1:A:80:ILE:HD13	2.03	0.40
1:A:95:GLU:O	1:A:142:SER:HA	2.22	0.40
1:A:365:THR:C	1:A:367:ASP:H	2.29	0.40
1:B:127:ASP:OD1	1:B:130:ARG:HG3	2.21	0.40
1:C:285:MET:HE1	1:D:64:THR:C	2.46	0.40
1:E:217:MET:HE3	1:E:258:VAL:HG13	2.03	0.40
1:F:64:THR:HB	1:F:286:THR:HG23	2.04	0.40
1:J:217:MET:O	1:J:270:TYR:CD2	2.73	0.40
1:L:161:THR:CG2	1:L:349:ILE:H	2.34	0.40
1:L:323:LEU:HD21	1:L:386:LEU:HD12	2.04	0.40
1:L:385:MET:HE2	1:L:385:MET:HB3	1.74	0.40
1:A:60:ASN:O	1:A:90:GLY:HA3	2.22	0.40
1:A:280:HIS:CE1	1:B:140:LYS:O	2.75	0.40
1:B:3:ASN:HB2	3:B:642:HOH:O	2.22	0.40
1:D:195:GLU:OE1	1:D:195:GLU:N	2.47	0.40
1:D:319:LEU:HA	1:D:319:LEU:HD23	1.93	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:394:GLU:O	1:F:394:GLU:HG2	2.22	0.40
1:G:211:VAL:HB	1:G:254:VAL:HG12	2.03	0.40
1:H:329:ASN:O	1:H:333:GLN:HG3	2.20	0.40
1:I:222:GLU:HB2	1:I:223:LEU:HD12	2.03	0.40
1:J:46:SER:OG	1:J:59:HIS:CE1	2.74	0.40
1:K:326:LEU:CB	1:K:385:MET:HE1	2.50	0.40
1:L:111:ILE:O	1:L:114:SER:OG	2.32	0.40
1:L:224:ILE:HD12	3:L:662:HOH:O	2.21	0.40
1:A:167:PHE:CZ	1:A:293:VAL:HG21	2.56	0.40
1:B:142:SER:N	3:B:630:HOH:O	2.48	0.40
1:B:229:HIS:CE1	3:B:690:HOH:O	2.74	0.40
1:B:230:TYR:CD2	1:B:269:LEU:HD23	2.57	0.40
1:C:138:ILE:HD13	1:C:138:ILE:HA	1.85	0.40
1:C:205:LEU:HD12	1:C:206:ARG:H	1.87	0.40
1:D:310:LYS:O	1:D:318:LYS:NZ	2.47	0.40
1:D:322:VAL:O	1:D:326:LEU:HD13	2.22	0.40
1:E:206:ARG:NH1	1:E:208:GLY:O	2.54	0.40
1:E:326:LEU:CB	1:E:385:MET:HE1	2.50	0.40
1:F:217:MET:O	1:F:270:TYR:CD2	2.75	0.40
1:G:283:ASP:OD2	1:H:90:GLY:HA2	2.21	0.40
1:H:56:ILE:HD11	1:H:397:LEU:HD11	2.01	0.40
1:J:155:VAL:HA	1:J:156:PRO:HD3	1.82	0.40
1:L:21:LEU:O	1:L:298:PRO:HD2	2.22	0.40
1:C:280:HIS:CE1	1:D:142:SER:HA	2.56	0.40
1:D:14:ILE:HD13	3:D:654:HOH:O	2.21	0.40
1:E:192:LYS:HA	1:E:192:LYS:HD2	1.75	0.40
1:E:205:LEU:HD12	1:E:206:ARG:H	1.86	0.40
1:E:269:LEU:HD12	1:E:269:LEU:N	2.37	0.40
1:E:348:GLN:HB2	1:E:357:LEU:HD12	2.02	0.40
1:E:356:PRO:HG2	1:E:386:LEU:HD21	2.03	0.40
1:G:252:TYR:HD1	1:G:252:TYR:HA	1.77	0.40
1:G:268:ILE:C	1:G:269:LEU:HD12	2.46	0.40
1:G:354:ASP:C	1:G:356:PRO:HD3	2.47	0.40
1:H:106:PRO:O	1:H:110:ILE:HG13	2.22	0.40
1:K:167:PHE:CE2	1:K:293:VAL:HG21	2.56	0.40
1:K:288:ARG:NH2	1:L:62:ASP:O	2.48	0.40
1:L:155:VAL:N	3:L:647:HOH:O	2.53	0.40
1:L:306:ALA:O	1:L:310:LYS:NZ	2.48	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all 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	385/413 (93%)	371 (96%)	13 (3%)	1 (0%)	36	66
1	B	386/413 (94%)	372 (96%)	12 (3%)	2 (0%)	24	55
1	C	385/413 (93%)	370 (96%)	14 (4%)	1 (0%)	36	66
1	D	386/413 (94%)	372 (96%)	12 (3%)	2 (0%)	24	55
1	E	385/413 (93%)	372 (97%)	12 (3%)	1 (0%)	36	66
1	F	386/413 (94%)	369 (96%)	14 (4%)	3 (1%)	16	44
1	G	385/413 (93%)	370 (96%)	13 (3%)	2 (0%)	24	55
1	H	386/413 (94%)	371 (96%)	13 (3%)	2 (0%)	24	55
1	I	385/413 (93%)	370 (96%)	14 (4%)	1 (0%)	36	66
1	J	386/413 (94%)	371 (96%)	13 (3%)	2 (0%)	24	55
1	K	385/413 (93%)	369 (96%)	14 (4%)	2 (0%)	24	55
1	L	386/413 (94%)	372 (96%)	12 (3%)	2 (0%)	24	55
All	All	4626/4956 (93%)	4449 (96%)	156 (3%)	21 (0%)	24	55

All (21) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	53	TYR
1	D	53	TYR
1	F	53	TYR
1	G	3	ASN
1	H	53	TYR
1	J	53	TYR
1	L	53	TYR
1	F	54	GLY
1	A	366	TYR
1	D	366	TYR
1	H	366	TYR
1	K	3	ASN

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Mol	Chain	Res	Type
1	K	366	TYR
1	L	366	TYR
1	B	366	TYR
1	C	366	TYR
1	E	366	TYR
1	G	366	TYR
1	I	366	TYR
1	J	366	TYR
1	F	366	TYR

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	336/354 (95%)	329 (98%)	7 (2%)	47	79
1	B	336/354 (95%)	329 (98%)	7 (2%)	47	79
1	C	336/354 (95%)	328 (98%)	8 (2%)	43	77
1	D	336/354 (95%)	328 (98%)	8 (2%)	43	77
1	E	336/354 (95%)	330 (98%)	6 (2%)	51	82
1	F	336/354 (95%)	330 (98%)	6 (2%)	51	82
1	G	336/354 (95%)	327 (97%)	9 (3%)	39	74
1	H	336/354 (95%)	326 (97%)	10 (3%)	36	72
1	I	336/354 (95%)	329 (98%)	7 (2%)	47	79
1	J	336/354 (95%)	329 (98%)	7 (2%)	47	79
1	K	336/354 (95%)	327 (97%)	9 (3%)	39	74
1	L	336/354 (95%)	330 (98%)	6 (2%)	51	82
All	All	4032/4248 (95%)	3942 (98%)	90 (2%)	45	78

All (90) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET

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Mol	Chain	Res	Type
1	A	60	ASN
1	A	87	ILE
1	A	237	VAL
1	A	293	VAL
1	A	299	MET
1	A	365	THR
1	B	87	ILE
1	B	91	ARG
1	B	165	VAL
1	B	276	THR
1	B	293	VAL
1	B	299	MET
1	B	316	VAL
1	C	1	MET
1	C	2	ARG
1	C	60	ASN
1	C	87	ILE
1	C	293	VAL
1	C	299	MET
1	C	316	VAL
1	C	378	ILE
1	D	87	ILE
1	D	91	ARG
1	D	165	VAL
1	D	276	THR
1	D	293	VAL
1	D	299	MET
1	D	316	VAL
1	D	374	THR
1	E	60	ASN
1	E	87	ILE
1	E	276	THR
1	E	293	VAL
1	E	299	MET
1	E	378	ILE
1	F	87	ILE
1	F	91	ARG
1	F	163	PHE
1	F	299	MET
1	F	316	VAL
1	F	374	THR
1	G	1	MET

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Mol	Chain	Res	Type
1	G	2	ARG
1	G	3	ASN
1	G	4	ILE
1	G	60	ASN
1	G	87	ILE
1	G	293	VAL
1	G	299	MET
1	G	316	VAL
1	H	5	ASN
1	H	81	ILE
1	H	87	ILE
1	H	91	ARG
1	H	165	VAL
1	H	293	VAL
1	H	299	MET
1	H	316	VAL
1	H	374	THR
1	H	383	ASP
1	I	87	ILE
1	I	165	VAL
1	I	171	THR
1	I	293	VAL
1	I	299	MET
1	I	365	THR
1	I	378	ILE
1	J	87	ILE
1	J	91	ARG
1	J	165	VAL
1	J	293	VAL
1	J	299	MET
1	J	316	VAL
1	J	374	THR
1	K	1	MET
1	K	2	ARG
1	K	4	ILE
1	K	60	ASN
1	K	87	ILE
1	K	186	SER
1	K	293	VAL
1	K	299	MET
1	K	316	VAL
1	L	87	ILE

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Mol	Chain	Res	Type
1	L	91	ARG
1	L	293	VAL
1	L	299	MET
1	L	316	VAL
1	L	374	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (43) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	183	HIS
1	A	239	ASN
1	A	280	HIS
1	B	3	ASN
1	B	18	GLN
1	B	70	GLN
1	B	253	ASN
1	C	280	HIS
1	D	5	ASN
1	D	18	GLN
1	D	58	HIS
1	D	82	GLN
1	D	315	HIS
1	E	18	GLN
1	E	60	ASN
1	E	280	HIS
1	F	5	ASN
1	F	104	GLN
1	F	183	HIS
1	F	267	ASN
1	F	315	HIS
1	G	18	GLN
1	G	59	HIS
1	G	280	HIS
1	G	325	ASN
1	H	3	ASN
1	H	315	HIS
1	I	7	GLN
1	I	18	GLN
1	I	121	ASN
1	I	183	HIS
1	I	280	HIS
1	J	3	ASN

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Mol	Chain	Res	Type
1	J	5	ASN
1	J	18	GLN
1	J	104	GLN
1	J	315	HIS
1	K	121	ASN
1	K	253	ASN
1	K	280	HIS
1	L	59	HIS
1	L	239	ASN
1	L	315	HIS

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 ⓘ

18 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	PO4	A	502	-	4,4,4	1.08	0	6,6,6	0.49	0
2	PO4	K	501	-	4,4,4	1.16	0	6,6,6	0.53	0
2	PO4	K	502	-	4,4,4	1.02	0	6,6,6	0.49	0
2	PO4	E	502	-	4,4,4	1.09	0	6,6,6	0.54	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	PO4	G	502	-	4,4,4	1.03	0	6,6,6	0.54	0
2	PO4	D	501	-	4,4,4	1.09	0	6,6,6	0.83	0
2	PO4	I	501	-	4,4,4	1.12	0	6,6,6	0.44	0
2	PO4	G	501	-	4,4,4	1.15	0	6,6,6	0.81	0
2	PO4	D	502	-	4,4,4	1.01	0	6,6,6	0.47	0
2	PO4	H	501	-	4,4,4	1.05	0	6,6,6	0.62	0
2	PO4	L	501	-	4,4,4	1.27	0	6,6,6	0.62	0
2	PO4	C	501	-	4,4,4	1.13	0	6,6,6	0.68	0
2	PO4	E	501	-	4,4,4	1.12	0	6,6,6	0.60	0
2	PO4	B	501	-	4,4,4	1.06	0	6,6,6	0.50	0
2	PO4	F	501	-	4,4,4	1.14	0	6,6,6	0.61	0
2	PO4	I	502	-	4,4,4	1.03	0	6,6,6	0.51	0
2	PO4	J	501	-	4,4,4	1.05	0	6,6,6	0.52	0
2	PO4	A	501	-	4,4,4	1.21	0	6,6,6	0.31	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

8 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	502	PO4	2	0
2	D	501	PO4	1	0
2	G	501	PO4	2	0
2	L	501	PO4	2	0
2	C	501	PO4	1	0
2	F	501	PO4	1	0
2	J	501	PO4	1	0
2	A	501	PO4	1	0

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 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	391/413 (94%)	-0.13	5 (1%) 75 66	23, 40, 71, 132	0
1	B	392/413 (94%)	-0.10	5 (1%) 75 66	25, 46, 75, 116	0
1	C	391/413 (94%)	0.01	1 (0%) 90 86	32, 53, 83, 119	0
1	D	392/413 (94%)	0.03	5 (1%) 75 66	31, 53, 83, 126	0
1	E	391/413 (94%)	-0.19	2 (0%) 87 82	18, 35, 67, 126	0
1	F	392/413 (94%)	-0.11	3 (0%) 82 75	22, 42, 71, 120	0
1	G	391/413 (94%)	-0.04	5 (1%) 75 66	23, 49, 85, 125	0
1	H	392/413 (94%)	0.07	9 (2%) 61 51	26, 53, 90, 126	0
1	I	391/413 (94%)	0.12	4 (1%) 79 72	40, 60, 91, 128	0
1	J	392/413 (94%)	0.14	9 (2%) 61 51	35, 61, 95, 126	0
1	K	391/413 (94%)	0.05	1 (0%) 90 86	33, 54, 87, 125	0
1	L	392/413 (94%)	0.08	6 (1%) 72 63	32, 58, 94, 146	0
All	All	4698/4956 (94%)	-0.01	55 (1%) 76 68	18, 50, 87, 146	0

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	L	147	GLY	4.9
1	D	12	SER	4.5
1	A	0	ILE	4.3
1	L	12	SER	4.2
1	J	163	PHE	3.9
1	H	12	SER	3.8
1	E	0	ILE	3.7
1	D	10	PRO	3.6
1	I	0	ILE	3.6
1	H	2	ARG	3.6
1	K	0	ILE	3.4

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Mol	Chain	Res	Type	RSRZ
1	J	12	SER	3.3
1	D	147	GLY	3.3
1	H	10	PRO	3.2
1	F	12	SER	3.1
1	H	163	PHE	3.1
1	B	163	PHE	3.0
1	G	0	ILE	2.8
1	B	372	ASP	2.8
1	J	376	ASN	2.6
1	C	0	ILE	2.6
1	L	10	PRO	2.5
1	G	400	LYS	2.5
1	J	337	ASP	2.5
1	H	147	GLY	2.5
1	F	11	LEU	2.4
1	H	144	ASP	2.4
1	J	187	LYS	2.4
1	J	372	ASP	2.4
1	G	372	ASP	2.4
1	L	11	LEU	2.4
1	B	367	ASP	2.3
1	H	367	ASP	2.3
1	A	372	ASP	2.3
1	H	257	TYR	2.3
1	L	372	ASP	2.2
1	A	155	VAL	2.2
1	A	159	GLY	2.2
1	B	12	SER	2.2
1	J	11	LEU	2.2
1	A	309	GLY	2.2
1	I	218	ALA	2.2
1	H	372	ASP	2.1
1	G	373	GLU	2.1
1	F	372	ASP	2.1
1	G	159	GLY	2.1
1	J	164	GLY	2.1
1	B	336	LYS	2.1
1	D	163	PHE	2.1
1	E	372	ASP	2.0
1	I	217	MET	2.0
1	D	389	PHE	2.0
1	I	159	GLY	2.0

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Mol	Chain	Res	Type	RSRZ
1	L	359	ALA	2.0
1	J	257	TYR	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](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	PO4	A	502	5/5	0.61	0.22	146,149,158,159	0
2	PO4	E	502	5/5	0.63	0.23	137,138,146,154	0
2	PO4	D	502	5/5	0.64	0.42	211,313,334,380	0
2	PO4	I	502	5/5	0.64	0.20	180,183,195,202	0
2	PO4	G	502	5/5	0.68	0.17	101,104,132,148	0
2	PO4	K	502	5/5	0.72	0.16	117,119,177,181	0
2	PO4	G	501	5/5	0.74	0.18	95,109,127,148	0
2	PO4	D	501	5/5	0.83	0.11	84,84,88,93	0
2	PO4	I	501	5/5	0.83	0.14	98,98,105,115	0
2	PO4	H	501	5/5	0.85	0.12	21,79,121,136	0
2	PO4	K	501	5/5	0.86	0.19	117,135,139,142	0
2	PO4	C	501	5/5	0.86	0.12	70,77,83,85	0
2	PO4	B	501	5/5	0.87	0.11	72,75,81,84	0
2	PO4	E	501	5/5	0.87	0.17	108,122,130,136	0
2	PO4	A	501	5/5	0.88	0.11	62,77,90,99	0
2	PO4	J	501	5/5	0.88	0.15	114,118,154,157	0
2	PO4	L	501	5/5	0.90	0.12	98,103,107,117	0
2	PO4	F	501	5/5	0.92	0.14	103,112,123,124	0

6.5 Other polymers [i](#)

There are no such residues in this entry.