



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 08:49 PM UTC

PDB ID : 3MOP / pdb_00003mop
Title : The ternary Death Domain complex of MyD88, IRAK4, and IRAK2
Authors : Lin, S.-C.; Lo, Y.-C.; Wu, H.
Deposited on : 2010-04-23
Resolution : 3.40 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
Xtrriage (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

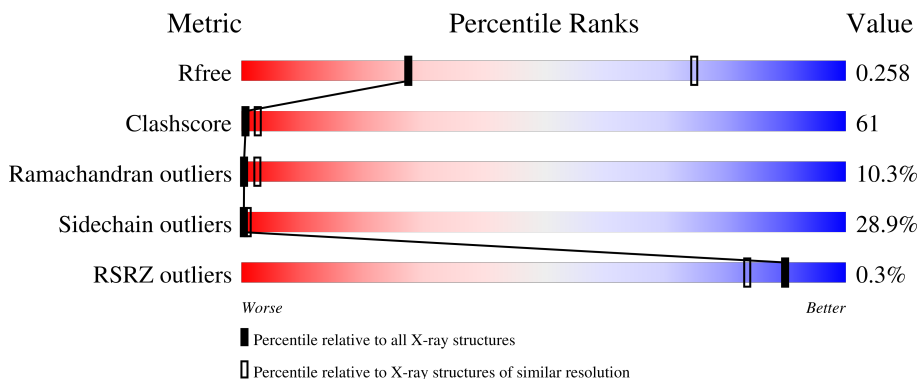
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.40 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.



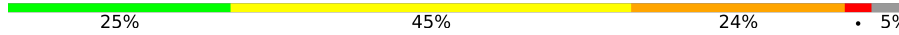
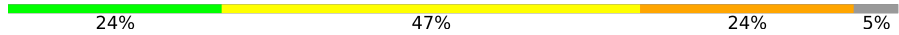
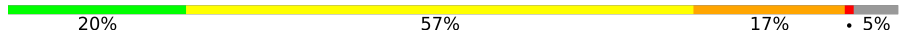
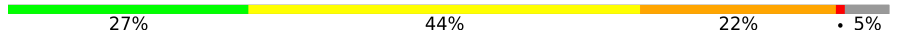
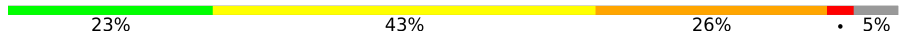
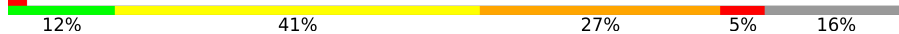

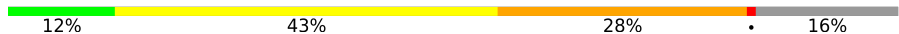
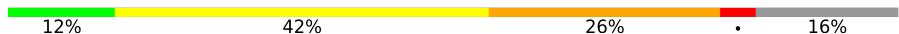
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1001 (3.44-3.36)
Clashscore	190562	1022 (3.44-3.36)
Ramachandran outliers	187476	1012 (3.44-3.36)
Sidechain outliers	187428	1012 (3.44-3.36)
RSRZ outliers	180081	1001 (3.44-3.36)

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	110	 % 31% 50% 14% • 5%
1	B	110	 29% 46% 19% • 5%
1	C	110	 20% 44% 29% • 5%
1	D	110	 25% 39% 30% • 5%
1	E	110	 % 15% 52% 25% • 5%

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Mol	Chain	Length	Quality of chain
1	F	110	
2	G	113	
2	H	113	
2	I	113	
2	J	113	
3	K	111	
3	L	111	
3	M	111	
3	N	111	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 11534 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 Myeloid differentiation primary response protein MyD88.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	105	833	525	148	156	4	0	0	0
1	B	105	833	525	148	156	4	0	0	0
1	C	105	833	525	148	156	4	0	0	0
1	D	105	833	525	148	156	4	0	0	0
1	E	105	833	525	148	156	4	0	0	0
1	F	105	833	525	148	156	4	0	0	0

There are 72 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	19	MET	-	expression tag	UNP Q99836
A	118	ALA	-	expression tag	UNP Q99836
A	119	ALA	-	expression tag	UNP Q99836
A	120	ALA	-	expression tag	UNP Q99836
A	121	LEU	-	expression tag	UNP Q99836
A	122	GLU	-	expression tag	UNP Q99836
A	123	HIS	-	expression tag	UNP Q99836
A	124	HIS	-	expression tag	UNP Q99836
A	125	HIS	-	expression tag	UNP Q99836
A	126	HIS	-	expression tag	UNP Q99836
A	127	HIS	-	expression tag	UNP Q99836
A	128	HIS	-	expression tag	UNP Q99836
B	19	MET	-	expression tag	UNP Q99836
B	118	ALA	-	expression tag	UNP Q99836
B	119	ALA	-	expression tag	UNP Q99836
B	120	ALA	-	expression tag	UNP Q99836
B	121	LEU	-	expression tag	UNP Q99836

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Chain	Residue	Modelled	Actual	Comment	Reference
B	122	GLU	-	expression tag	UNP Q99836
B	123	HIS	-	expression tag	UNP Q99836
B	124	HIS	-	expression tag	UNP Q99836
B	125	HIS	-	expression tag	UNP Q99836
B	126	HIS	-	expression tag	UNP Q99836
B	127	HIS	-	expression tag	UNP Q99836
B	128	HIS	-	expression tag	UNP Q99836
C	19	MET	-	expression tag	UNP Q99836
C	118	ALA	-	expression tag	UNP Q99836
C	119	ALA	-	expression tag	UNP Q99836
C	120	ALA	-	expression tag	UNP Q99836
C	121	LEU	-	expression tag	UNP Q99836
C	122	GLU	-	expression tag	UNP Q99836
C	123	HIS	-	expression tag	UNP Q99836
C	124	HIS	-	expression tag	UNP Q99836
C	125	HIS	-	expression tag	UNP Q99836
C	126	HIS	-	expression tag	UNP Q99836
C	127	HIS	-	expression tag	UNP Q99836
C	128	HIS	-	expression tag	UNP Q99836
D	19	MET	-	expression tag	UNP Q99836
D	118	ALA	-	expression tag	UNP Q99836
D	119	ALA	-	expression tag	UNP Q99836
D	120	ALA	-	expression tag	UNP Q99836
D	121	LEU	-	expression tag	UNP Q99836
D	122	GLU	-	expression tag	UNP Q99836
D	123	HIS	-	expression tag	UNP Q99836
D	124	HIS	-	expression tag	UNP Q99836
D	125	HIS	-	expression tag	UNP Q99836
D	126	HIS	-	expression tag	UNP Q99836
D	127	HIS	-	expression tag	UNP Q99836
D	128	HIS	-	expression tag	UNP Q99836
E	19	MET	-	expression tag	UNP Q99836
E	118	ALA	-	expression tag	UNP Q99836
E	119	ALA	-	expression tag	UNP Q99836
E	120	ALA	-	expression tag	UNP Q99836
E	121	LEU	-	expression tag	UNP Q99836
E	122	GLU	-	expression tag	UNP Q99836
E	123	HIS	-	expression tag	UNP Q99836
E	124	HIS	-	expression tag	UNP Q99836
E	125	HIS	-	expression tag	UNP Q99836
E	126	HIS	-	expression tag	UNP Q99836
E	127	HIS	-	expression tag	UNP Q99836

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Chain	Residue	Modelled	Actual	Comment	Reference
E	128	HIS	-	expression tag	UNP Q99836
F	19	MET	-	expression tag	UNP Q99836
F	118	ALA	-	expression tag	UNP Q99836
F	119	ALA	-	expression tag	UNP Q99836
F	120	ALA	-	expression tag	UNP Q99836
F	121	LEU	-	expression tag	UNP Q99836
F	122	GLU	-	expression tag	UNP Q99836
F	123	HIS	-	expression tag	UNP Q99836
F	124	HIS	-	expression tag	UNP Q99836
F	125	HIS	-	expression tag	UNP Q99836
F	126	HIS	-	expression tag	UNP Q99836
F	127	HIS	-	expression tag	UNP Q99836
F	128	HIS	-	expression tag	UNP Q99836

- Molecule 2 is a protein called Interleukin-1 receptor-associated kinase 4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	G	107	843	544	140	156	3	0	0	0
2	H	107	843	544	140	156	3	0	0	0
2	I	107	843	544	140	156	3	0	0	0
2	J	107	843	544	140	156	3	0	0	0

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
G	2	MET	-	expression tag	UNP Q9NWZ3
G	3	GLY	-	expression tag	UNP Q9NWZ3
G	107	LEU	-	expression tag	UNP Q9NWZ3
G	108	GLU	-	expression tag	UNP Q9NWZ3
G	109	HIS	-	expression tag	UNP Q9NWZ3
G	110	HIS	-	expression tag	UNP Q9NWZ3
G	111	HIS	-	expression tag	UNP Q9NWZ3
G	112	HIS	-	expression tag	UNP Q9NWZ3
G	113	HIS	-	expression tag	UNP Q9NWZ3
G	114	HIS	-	expression tag	UNP Q9NWZ3
H	2	MET	-	expression tag	UNP Q9NWZ3
H	3	GLY	-	expression tag	UNP Q9NWZ3
H	107	LEU	-	expression tag	UNP Q9NWZ3

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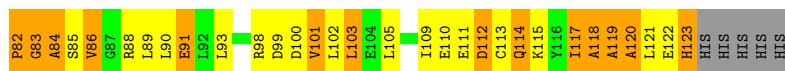
Chain	Residue	Modelled	Actual	Comment	Reference
H	108	GLU	-	expression tag	UNP Q9NWZ3
H	109	HIS	-	expression tag	UNP Q9NWZ3
H	110	HIS	-	expression tag	UNP Q9NWZ3
H	111	HIS	-	expression tag	UNP Q9NWZ3
H	112	HIS	-	expression tag	UNP Q9NWZ3
H	113	HIS	-	expression tag	UNP Q9NWZ3
H	114	HIS	-	expression tag	UNP Q9NWZ3
I	2	MET	-	expression tag	UNP Q9NWZ3
I	3	GLY	-	expression tag	UNP Q9NWZ3
I	107	LEU	-	expression tag	UNP Q9NWZ3
I	108	GLU	-	expression tag	UNP Q9NWZ3
I	109	HIS	-	expression tag	UNP Q9NWZ3
I	110	HIS	-	expression tag	UNP Q9NWZ3
I	111	HIS	-	expression tag	UNP Q9NWZ3
I	112	HIS	-	expression tag	UNP Q9NWZ3
I	113	HIS	-	expression tag	UNP Q9NWZ3
I	114	HIS	-	expression tag	UNP Q9NWZ3
J	2	MET	-	expression tag	UNP Q9NWZ3
J	3	GLY	-	expression tag	UNP Q9NWZ3
J	107	LEU	-	expression tag	UNP Q9NWZ3
J	108	GLU	-	expression tag	UNP Q9NWZ3
J	109	HIS	-	expression tag	UNP Q9NWZ3
J	110	HIS	-	expression tag	UNP Q9NWZ3
J	111	HIS	-	expression tag	UNP Q9NWZ3
J	112	HIS	-	expression tag	UNP Q9NWZ3
J	113	HIS	-	expression tag	UNP Q9NWZ3
J	114	HIS	-	expression tag	UNP Q9NWZ3

- Molecule 3 is a protein called Interleukin-1 receptor-associated kinase-like 2.

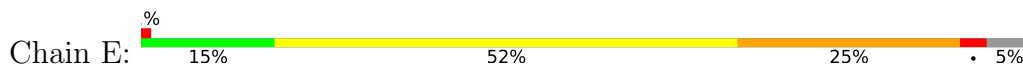
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	K	93	Total	C	N	O	S	0	0	0
			791	516	131	137	7			
3	L	93	Total	C	N	O	S	0	0	0
			791	516	131	137	7			
3	M	93	Total	C	N	O	S	0	0	0
			791	516	131	137	7			
3	N	93	Total	C	N	O	S	0	0	0
			791	516	131	137	7			

There are 4 discrepancies between the modelled and reference sequences:

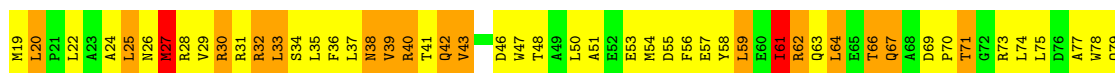
Chain	Residue	Modelled	Actual	Comment	Reference
K	50	TRP	ARG	engineered mutation	UNP O43187
L	50	TRP	ARG	engineered mutation	UNP O43187
M	50	TRP	ARG	engineered mutation	UNP O43187
N	50	TRP	ARG	engineered mutation	UNP O43187



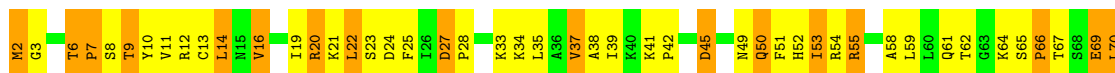
• Molecule 1: Myeloid differentiation primary response protein MyD88



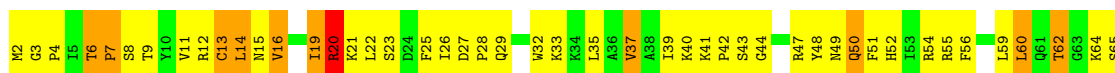
• Molecule 1: Myeloid differentiation primary response protein MyD88



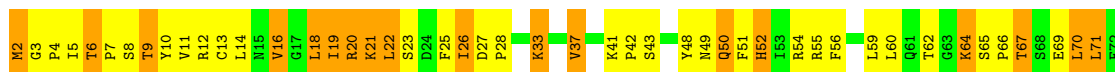
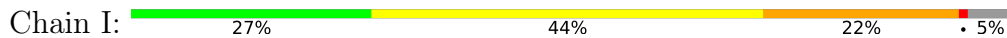
• Molecule 2: Interleukin-1 receptor-associated kinase 4



• Molecule 2: Interleukin-1 receptor-associated kinase 4



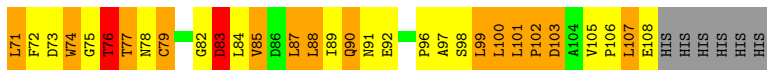
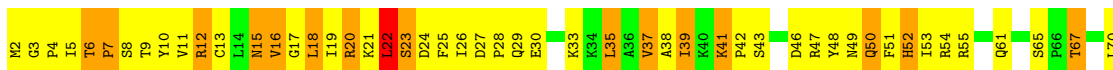
• Molecule 2: Interleukin-1 receptor-associated kinase 4





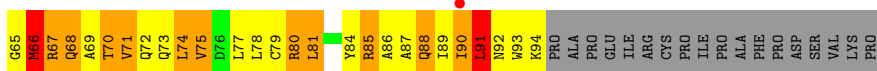
- Molecule 2: Interleukin-1 receptor-associated kinase 4

Chain J: 23% 43% 26% 5%



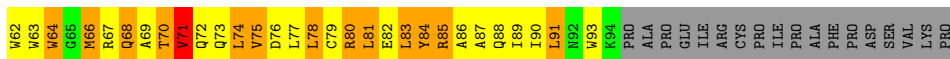
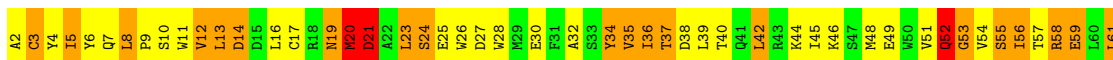
- Molecule 3: Interleukin-1 receptor-associated kinase-like 2

Chain K: 2% 12% 41% 27% 5% 16%



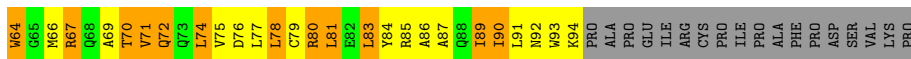
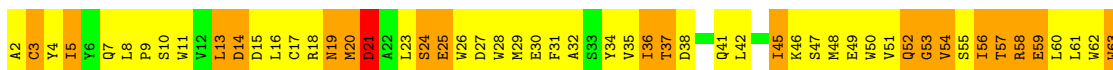
- Molecule 3: Interleukin-1 receptor-associated kinase-like 2

Chain L: 13% 38% 30% 16%



- Molecule 3: Interleukin-1 receptor-associated kinase-like 2

Chain M: 12% 43% 28% 16%



- Molecule 3: Interleukin-1 receptor-associated kinase-like 2

Chain N: 12% 42% 26% 16%



M62	M63	M64	G65	M66	R67	Q68	A69	T70	V71	Q72	Q73	L74	V75	D76	L77	L78	C79	R80	L81	E82	L83	Y84	R85	A86	A87	Q88	I89	I90	L91	I92	M93	R94	PRO	ALA	PRO	PRO	GLU	GLU	ILE	ILE	ARG	CYS	PRO	PRO	ILE	ILE	PRO	ALA	PHE	PRO	ASP	ASP	SER	VAL	VAL	LYS	LYS	PRO
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4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	101.71Å 307.12Å 187.62Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.96 – 3.40 19.96 – 3.40	Depositor EDS
% Data completeness (in resolution range)	82.3 (19.96-3.40) 81.8 (19.96-3.40)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.49 (at 3.31Å)	Xtrriage
Refinement program	PHENIX 1.6_289	Depositor
R, R_{free}	0.216 , 0.261 0.211 , 0.258	Depositor DCC
R_{free} test set	2000 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	110.9	Xtrriage
Anisotropy	0.396	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.25 , 153.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	11534	wwPDB-VP
Average B, all atoms (Å ²)	179.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.47% 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 i

5.1 Standard geometry i

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	A	0.56	0/845	0.89	1/1144 (0.1%)
1	B	0.74	0/845	1.09	3/1144 (0.3%)
1	C	0.67	0/845	1.09	3/1144 (0.3%)
1	D	0.82	0/845	1.26	7/1144 (0.6%)
1	E	0.97	1/845 (0.1%)	1.43	12/1144 (1.0%)
1	F	0.74	0/845	1.14	3/1144 (0.3%)
2	G	0.71	0/863	1.11	4/1174 (0.3%)
2	H	0.80	0/863	1.12	1/1174 (0.1%)
2	I	0.77	0/863	1.10	1/1174 (0.1%)
2	J	0.58	0/863	1.05	2/1174 (0.2%)
3	K	0.61	0/812	0.95	3/1106 (0.3%)
3	L	0.65	0/812	1.01	4/1106 (0.4%)
3	M	0.57	0/812	0.94	1/1106 (0.1%)
3	N	0.54	0/812	0.95	3/1106 (0.3%)
All	All	0.71	1/11770 (0.0%)	1.09	48/15984 (0.3%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	45	ALA	CA-CB	-7.27	1.43	1.53

All (48) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	E	69	ASP	CA-C-N	-13.72	105.78	119.64
1	E	69	ASP	C-N-CA	-13.72	105.78	119.64
1	F	64	LEU	N-CA-C	-10.54	102.60	114.62
1	D	69	ASP	CA-C-N	-9.48	109.14	119.19
1	D	69	ASP	C-N-CA	-9.48	109.14	119.19
1	E	20	LEU	CA-C-N	8.63	130.22	120.98
1	E	20	LEU	C-N-CA	8.63	130.22	120.98
1	D	86	VAL	N-CA-C	-7.68	102.96	110.72
1	C	64	LEU	N-CA-C	-7.47	105.33	114.75

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	86	VAL	N-CA-C	-6.91	103.74	110.72
1	E	49	ALA	N-CA-C	-6.38	104.42	111.82
1	D	29	VAL	N-CA-C	6.37	116.53	110.42
1	E	26	ASN	N-CA-C	6.32	118.38	108.96
1	B	69	ASP	CA-C-N	6.29	125.91	119.56
1	B	69	ASP	C-N-CA	6.29	125.91	119.56
3	N	71	VAL	N-CA-C	-6.22	107.43	113.53
3	L	8	LEU	CA-C-N	6.15	126.51	119.93
3	L	8	LEU	C-N-CA	6.15	126.51	119.93
1	C	104	GLU	N-CA-C	-6.08	105.62	113.16
3	L	71	VAL	N-CA-C	-6.05	107.60	113.53
1	E	70	PRO	CA-C-N	-5.93	112.33	120.28
1	E	70	PRO	C-N-CA	-5.93	112.33	120.28
3	K	71	VAL	N-CA-C	-5.93	107.72	113.53
1	C	36	PHE	N-CA-C	-5.91	105.07	112.93
2	I	73	ASP	N-CA-C	5.89	118.49	111.71
1	E	116	TYR	N-CA-C	-5.88	103.67	111.96
3	L	64	TRP	N-CA-C	-5.86	105.97	113.23
3	M	64	TRP	N-CA-C	-5.83	104.89	112.23
1	D	81	ARG	CA-C-N	5.63	126.88	119.84
1	D	81	ARG	C-N-CA	5.63	126.88	119.84
2	H	74	TRP	N-CA-C	-5.63	106.00	112.92
1	B	115	LYS	N-CA-C	-5.62	106.59	113.50
1	E	61	ILE	CB-CA-C	-5.62	104.56	112.14
1	F	39	VAL	CB-CA-C	-5.51	104.25	111.25
1	F	61	ILE	CB-CA-C	-5.46	104.69	112.22
1	D	26	ASN	N-CA-C	5.42	117.04	108.96
3	N	50	TRP	CB-CA-C	5.36	118.42	111.50
3	K	91	LEU	N-CA-C	-5.35	106.76	113.72
3	K	64	TRP	N-CA-C	-5.35	106.26	112.89
2	G	65	SER	CA-C-N	-5.34	114.53	120.45
2	G	65	SER	C-N-CA	-5.34	114.53	120.45
1	A	45	ALA	N-CA-C	5.32	118.11	107.62
2	G	33	LYS	N-CA-C	-5.29	105.52	111.28
2	J	74	TRP	N-CA-C	-5.22	106.75	113.02
2	G	66	PRO	N-CA-C	-5.22	107.83	114.35
1	E	90	LEU	N-CA-C	5.21	116.77	111.14
3	N	91	LEU	N-CA-C	-5.12	107.00	114.12
2	J	22	LEU	N-CA-C	-5.10	107.03	113.20

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	833	0	844	91	0
1	B	833	0	844	96	0
1	C	833	0	844	112	0
1	D	833	0	844	115	0
1	E	833	0	844	127	0
1	F	833	0	844	119	0
2	G	843	0	849	103	0
2	H	843	0	849	119	0
2	I	843	0	849	106	0
2	J	843	0	849	116	0
3	K	791	0	784	109	0
3	L	791	0	784	120	0
3	M	791	0	784	111	0
3	N	791	0	784	118	0
All	All	11534	0	11596	1417	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 61.

All (1417) 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)
3:N:13:LEU:HD21	3:N:61:LEU:HD23	1.24	1.16
1:B:114:GLN:HA	1:B:117:ILE:HG22	1.23	1.12
3:L:20:MET:HB3	3:L:57:THR:HG21	1.29	1.11
2:G:100:LEU:HD22	2:G:101:LEU:HD22	1.22	1.10
2:H:2:MET:HG3	2:H:3:GLY:H	1.09	1.10
2:G:34:LYS:HD2	2:G:93:PHE:HZ	1.18	1.07
3:M:55:SER:HB2	3:M:58:ARG:HB2	1.07	1.07
1:C:22:LEU:HD22	1:C:22:LEU:H	1.20	1.06
1:F:30:ARG:HG3	1:F:30:ARG:HH11	1.16	1.06
3:K:55:SER:HB3	3:K:58:ARG:HB2	1.35	1.05
3:N:45:ILE:O	3:N:56:ILE:HD13	1.57	1.04
1:D:40:ARG:NH1	1:D:40:ARG:HB2	1.74	1.03
2:J:11:VAL:HG23	2:J:79:CYS:O	1.56	1.02

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:46:ASP:OD1	1:D:48:THR:HG23	1.56	1.02
1:E:37:LEU:HB2	1:E:71:THR:HG21	1.41	1.02
2:H:7:PRO:HG3	2:H:101:LEU:HD13	1.41	1.02
2:H:2:MET:CG	2:H:3:GLY:H	1.73	1.01
1:E:20:LEU:HD12	1:E:20:LEU:H	1.24	1.01
1:B:81:ARG:HB3	1:B:82:PRO:HD2	1.40	1.00
3:K:16:LEU:HB2	3:K:89:ILE:HD11	1.45	0.99
2:I:50:GLN:HE21	3:L:58:ARG:HH22	1.01	0.97
1:F:85:SER:HB3	1:F:88:ARG:HB2	1.45	0.95
3:M:13:LEU:HD21	3:M:61:LEU:HD13	1.48	0.95
3:M:64:TRP:HZ3	3:M:74:LEU:HD23	1.29	0.95
2:G:34:LYS:HD2	2:G:93:PHE:CZ	2.02	0.94
2:J:84:LEU:HD23	2:J:100:LEU:HD11	1.49	0.94
1:D:33:LEU:HD13	1:D:105:LEU:HD12	1.50	0.94
3:M:20:MET:CB	3:M:57:THR:HG21	1.97	0.94
3:M:9:PRO:HG2	3:M:93:TRP:CH2	2.03	0.93
1:F:43:VAL:HG11	2:J:48:TYR:CE1	2.03	0.93
1:C:85:SER:HB3	1:C:88:ARG:HB2	1.51	0.93
1:E:38:ASN:HD21	1:E:71:THR:H	1.15	0.92
1:F:67:GLN:NE2	1:F:67:GLN:H	1.67	0.92
1:D:32:ARG:HH11	1:D:32:ARG:HG2	1.30	0.92
3:N:37:THR:HG22	3:N:38:ASP:H	1.32	0.91
1:C:50:LEU:HD11	1:C:54:MET:HE3	1.49	0.91
2:H:95:ALA:HA	3:L:68:GLN:HE22	1.35	0.91
3:L:55:SER:HB2	3:L:58:ARG:HB2	1.52	0.90
2:J:41:LYS:HE3	2:J:41:LYS:HA	1.53	0.90
2:J:100:LEU:HB3	2:J:101:LEU:HD23	1.52	0.90
3:N:55:SER:CB	3:N:58:ARG:HB2	2.01	0.90
3:N:20:MET:O	3:N:28:TRP:HD1	1.54	0.89
2:G:41:LYS:HB3	2:G:42:PRO:HD2	1.54	0.89
1:B:58:TYR:CE1	1:E:30:ARG:CZ	2.56	0.89
3:L:5:ILE:HD11	3:L:69:ALA:CB	2.03	0.89
3:L:20:MET:H	3:L:20:MET:CE	1.86	0.88
2:G:96:PRO:HD3	3:K:66:MET:CE	2.02	0.88
1:E:43:VAL:HG21	2:I:48:TYR:CE1	2.09	0.88
1:F:25:LEU:HD21	1:F:86:VAL:HG22	1.56	0.88
3:K:38:ASP:OD2	3:K:41:GLN:HG2	1.74	0.88
2:J:18:LEU:C	2:J:18:LEU:HD12	1.98	0.87
1:D:40:ARG:HB2	1:D:40:ARG:HH11	1.35	0.87
2:G:55:ARG:HG3	2:G:55:ARG:HH11	1.39	0.87
2:I:50:GLN:HE21	3:L:58:ARG:NH2	1.71	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:8:LEU:HD12	3:L:12:VAL:HG11	1.56	0.87
1:C:38:ASN:HD21	1:C:71:THR:H	1.20	0.87
1:E:81:ARG:HB3	1:E:82:PRO:HD2	1.55	0.87
2:J:25:PHE:HB3	2:J:96:PRO:HG3	1.54	0.87
2:J:50:GLN:HE21	3:M:58:ARG:NH2	1.72	0.87
3:N:55:SER:HB2	3:N:58:ARG:HB2	1.56	0.87
2:G:81:VAL:HG12	2:G:100:LEU:HD21	1.57	0.87
3:N:28:TRP:CZ2	3:N:56:ILE:HG13	2.11	0.86
2:J:7:PRO:HG3	2:J:101:LEU:HD13	1.57	0.86
1:F:36:PHE:O	1:F:39:VAL:HG22	1.75	0.86
3:M:55:SER:CB	3:M:58:ARG:HB2	2.00	0.86
2:H:2:MET:HG3	2:H:4:PRO:HD2	1.56	0.86
1:D:38:ASN:HD21	1:D:71:THR:H	1.23	0.86
3:K:70:THR:H	3:K:73:GLN:NE2	1.74	0.86
3:M:55:SER:HB2	3:M:58:ARG:CB	2.02	0.86
3:L:20:MET:CB	3:L:57:THR:HG21	2.06	0.85
2:G:95:ALA:HA	3:K:68:GLN:HE22	1.40	0.85
2:I:18:LEU:C	2:I:18:LEU:HD23	2.02	0.85
1:C:58:TYR:HB2	1:F:26:ASN:HB3	1.56	0.85
3:N:86:ALA:O	3:N:89:ILE:HG13	1.77	0.84
1:D:43:VAL:HG21	2:H:48:TYR:CE1	2.11	0.84
3:L:5:ILE:HD11	3:L:69:ALA:HB1	1.57	0.84
1:F:39:VAL:HG23	1:F:39:VAL:O	1.76	0.84
3:K:84:TYR:O	3:K:88:GLN:HG2	1.76	0.83
1:C:47:TRP:HB3	1:C:71:THR:HG22	1.59	0.83
2:H:2:MET:HG3	2:H:3:GLY:N	1.92	0.83
3:L:35:VAL:HG12	3:L:37:THR:OG1	1.79	0.83
1:A:122:GLU:HG3	1:A:123:HIS:H	1.43	0.83
1:D:93:LEU:HD13	1:D:101:VAL:HG11	1.59	0.83
3:L:16:LEU:HD12	3:L:90:ILE:CG2	2.08	0.83
3:L:84:TYR:O	3:L:88:GLN:HG3	1.79	0.82
3:L:64:TRP:HZ3	3:L:74:LEU:HD23	1.45	0.82
3:M:64:TRP:CZ3	3:M:74:LEU:HD23	2.15	0.82
2:J:50:GLN:HE21	3:M:58:ARG:HH22	1.24	0.82
2:I:11:VAL:HG11	2:I:79:CYS:SG	2.18	0.82
2:G:81:VAL:CG1	2:G:100:LEU:HD21	2.10	0.82
1:E:120:ALA:O	1:E:121:LEU:HD23	1.80	0.81
1:D:38:ASN:ND2	1:D:71:THR:HG23	1.95	0.81
3:L:20:MET:H	3:L:20:MET:HE2	1.45	0.81
1:F:30:ARG:HH11	1:F:30:ARG:CG	1.92	0.81
2:H:85:VAL:HG12	2:H:97:ALA:HB1	1.62	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:J:10:TYR:HE1	2:J:78:ASN:HB3	1.45	0.81
1:E:64:LEU:O	1:E:66:THR:N	2.13	0.81
3:M:32:ALA:HB1	3:M:42:LEU:HD11	1.62	0.81
3:M:20:MET:HB2	3:M:57:THR:HG21	1.61	0.81
1:C:41:THR:HG22	2:G:72:PHE:CE2	2.16	0.81
3:K:64:TRP:HA	3:K:67:ARG:HE	1.44	0.80
2:J:4:PRO:O	2:J:6:THR:HG22	1.81	0.80
3:M:49:GLU:HB2	3:M:56:ILE:HD12	1.63	0.80
1:E:98:ARG:NH1	1:E:100:ASP:OD2	2.15	0.80
2:H:95:ALA:CA	3:L:68:GLN:HE22	1.93	0.80
1:D:20:LEU:HD12	1:D:20:LEU:H	1.46	0.80
2:J:37:VAL:HG11	3:M:11:TRP:N	1.97	0.80
3:M:45:ILE:HA	3:M:48:MET:HE2	1.60	0.80
2:I:92:GLU:HA	2:I:94:PHE:CE2	2.17	0.80
1:C:29:VAL:HG13	1:C:105:LEU:HD11	1.64	0.80
1:F:102:LEU:HD12	1:F:102:LEU:H	1.46	0.80
2:H:41:LYS:HB3	2:H:42:PRO:HD2	1.64	0.80
1:A:29:VAL:HG13	1:A:105:LEU:HD22	1.62	0.79
1:E:31:ARG:HH11	1:E:31:ARG:CG	1.95	0.79
2:G:85:VAL:HG23	2:G:97:ALA:HB1	1.64	0.79
1:C:22:LEU:O	1:C:25:LEU:HD23	1.82	0.79
1:F:20:LEU:HD12	1:F:20:LEU:H	1.47	0.79
1:F:56:PHE:CE1	1:F:74:LEU:HD11	2.18	0.79
3:N:35:VAL:CG1	3:N:37:THR:OG1	2.31	0.79
1:E:25:LEU:HD13	1:E:86:VAL:CG2	2.12	0.79
2:H:85:VAL:HG21	2:H:101:LEU:HD11	1.63	0.79
3:M:16:LEU:HB2	3:M:89:ILE:HD11	1.64	0.79
3:M:20:MET:HB3	3:M:57:THR:HG21	1.63	0.79
1:E:20:LEU:HD12	1:E:20:LEU:N	1.98	0.79
3:N:56:ILE:HD12	3:N:56:ILE:O	1.83	0.79
2:G:100:LEU:HD22	2:G:101:LEU:CD2	2.09	0.79
2:I:10:TYR:HB2	2:I:13:CYS:SG	2.24	0.78
3:L:16:LEU:HD12	3:L:90:ILE:HG21	1.65	0.78
3:L:86:ALA:O	3:L:89:ILE:HG12	1.83	0.78
2:H:23:SER:O	2:H:27:ASP:HB2	1.83	0.78
1:B:40:ARG:HH21	1:E:31:ARG:HH22	1.30	0.78
1:A:81:ARG:HB3	1:A:82:PRO:HD2	1.65	0.78
2:G:96:PRO:HD3	3:K:66:MET:HE1	1.66	0.78
3:N:63:TRP:CE2	3:N:67:ARG:NH2	2.52	0.77
2:H:11:VAL:HG23	2:H:79:CYS:O	1.84	0.77
1:D:28:ARG:HG3	1:D:29:VAL:N	1.99	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:M:49:GLU:HB2	3:M:56:ILE:CD1	2.14	0.77
3:K:20:MET:HE3	3:K:20:MET:H	1.48	0.77
2:H:89:ILE:HD12	2:H:90:GLN:N	1.99	0.77
1:A:20:LEU:HB2	1:A:21:PRO:HD2	1.65	0.77
1:C:40:ARG:HH11	1:C:40:ARG:HB2	1.50	0.77
1:E:62:ARG:HH11	1:E:62:ARG:HB2	1.48	0.77
2:H:84:LEU:HD23	2:H:100:LEU:HD11	1.67	0.76
2:G:77:THR:HG22	2:G:77:THR:O	1.85	0.76
3:M:70:THR:HG21	3:M:72:GLN:HE21	1.50	0.76
1:E:67:GLN:H	1:E:67:GLN:NE2	1.83	0.76
2:G:96:PRO:HD3	3:K:66:MET:HE2	1.68	0.76
1:C:41:THR:HG22	2:G:72:PHE:HE2	1.51	0.76
2:G:50:GLN:HA	2:J:16:VAL:HG22	1.68	0.76
2:I:101:LEU:N	2:I:101:LEU:HD23	2.01	0.76
3:K:80:ARG:O	3:K:81:LEU:HD22	1.85	0.76
3:N:13:LEU:CD2	3:N:61:LEU:HD23	2.11	0.76
2:H:95:ALA:HA	3:L:68:GLN:NE2	2.01	0.76
1:D:64:LEU:O	1:D:66:THR:N	2.16	0.76
1:D:32:ARG:HH11	1:D:32:ARG:CG	1.99	0.75
3:K:70:THR:H	3:K:73:GLN:HE21	1.33	0.75
1:E:19:MET:HE3	1:E:87:GLY:HA3	1.67	0.75
1:E:59:LEU:O	1:E:59:LEU:HD23	1.85	0.75
1:F:40:ARG:HD3	1:F:46:ASP:OD2	1.85	0.75
1:B:42:GLN:HG3	1:B:42:GLN:O	1.86	0.75
1:A:93:LEU:HB2	1:A:102:LEU:HD11	1.67	0.75
2:H:80:THR:HG23	2:H:83:ASP:H	1.52	0.75
2:J:10:TYR:CE1	2:J:78:ASN:HB3	2.21	0.75
3:K:55:SER:CB	3:K:58:ARG:HB2	2.15	0.75
3:N:80:ARG:HB2	3:N:80:ARG:CZ	2.15	0.75
1:C:37:LEU:HB2	1:C:71:THR:HG21	1.68	0.74
3:L:23:LEU:HD11	3:L:83:LEU:HD21	1.67	0.74
2:J:85:VAL:HG12	2:J:97:ALA:HB1	1.67	0.74
1:A:69:ASP:N	1:A:70:PRO:HD3	2.02	0.74
1:D:35:LEU:O	1:E:59:LEU:HD21	1.87	0.74
1:C:40:ARG:HB2	1:C:40:ARG:NH1	2.02	0.74
1:F:25:LEU:HD21	1:F:86:VAL:CG2	2.17	0.74
1:B:81:ARG:HB3	1:B:82:PRO:CD	2.18	0.74
1:D:43:VAL:HG21	2:H:48:TYR:HE1	1.49	0.74
3:N:28:TRP:HZ2	3:N:56:ILE:HG13	1.50	0.74
3:N:87:ALA:O	3:N:91:LEU:HD23	1.88	0.74
1:C:102:LEU:HD12	1:C:102:LEU:H	1.52	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:58:TYR:HE1	1:E:30:ARG:CZ	2.00	0.73
2:J:7:PRO:HG3	2:J:101:LEU:CD1	2.18	0.73
1:B:79:GLN:HE21	1:B:80:GLY:N	1.87	0.73
2:J:25:PHE:CB	2:J:96:PRO:HG3	2.18	0.73
3:M:23:LEU:HD23	3:M:23:LEU:O	1.89	0.73
2:G:22:LEU:O	2:G:22:LEU:HG	1.88	0.73
2:J:18:LEU:C	2:J:18:LEU:CD1	2.59	0.73
3:L:45:ILE:HD11	3:L:59:GLU:HB3	1.70	0.73
3:M:3:CYS:O	3:M:71:VAL:HG12	1.89	0.73
1:E:107:PRO:HG2	1:E:108:SER:H	1.54	0.73
3:M:11:TRP:HA	3:M:14:ASP:HB2	1.71	0.72
1:C:54:MET:HB3	1:C:78:TRP:CD1	2.23	0.72
1:E:37:LEU:HB2	1:E:71:THR:CG2	2.18	0.72
1:E:81:ARG:CB	1:E:82:PRO:HD2	2.19	0.72
2:G:85:VAL:HA	2:G:88:LEU:HB2	1.71	0.72
1:C:94:THR:HG23	1:C:102:LEU:HD22	1.70	0.72
1:E:73:ARG:HA	1:E:73:ARG:NE	2.04	0.72
2:J:107:LEU:HD12	2:J:107:LEU:H	1.54	0.72
3:L:24:SER:HB3	3:L:27:ASP:HB2	1.71	0.72
1:F:30:ARG:HG3	1:F:30:ARG:NH1	1.98	0.72
1:D:39:VAL:HG23	1:D:39:VAL:O	1.88	0.72
2:J:48:TYR:OH	2:J:74:TRP:HB2	1.90	0.72
3:M:67:ARG:HB2	3:M:67:ARG:CZ	2.18	0.72
1:B:59:LEU:H	1:B:59:LEU:HD12	1.55	0.71
1:D:32:ARG:HG2	1:D:32:ARG:NH1	1.97	0.71
1:F:73:ARG:HA	1:F:73:ARG:NE	2.05	0.71
1:F:101:VAL:HG12	1:F:102:LEU:HD12	1.72	0.71
3:L:3:CYS:O	3:L:71:VAL:HG12	1.91	0.71
1:B:69:ASP:N	1:B:70:PRO:HD3	2.04	0.71
1:E:40:ARG:HB2	1:E:40:ARG:NH1	2.04	0.71
1:F:25:LEU:O	1:F:30:ARG:NH1	2.23	0.71
2:H:6:THR:HG23	2:H:9:THR:HG23	1.71	0.71
2:H:20:ARG:O	2:H:20:ARG:HG2	1.88	0.71
1:D:28:ARG:HG3	1:D:29:VAL:H	1.54	0.71
2:H:50:GLN:HE21	3:K:58:ARG:NH2	1.89	0.70
1:C:93:LEU:HD13	1:C:101:VAL:CG1	2.21	0.70
3:K:55:SER:HB3	3:K:58:ARG:CB	2.19	0.70
1:A:32:ARG:HG2	1:A:105:LEU:HD21	1.72	0.70
1:E:31:ARG:HH11	1:E:31:ARG:HG3	1.55	0.70
2:H:92:GLU:HA	2:H:94:PHE:CE2	2.27	0.70
2:I:4:PRO:O	2:I:6:THR:HG22	1.92	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:26:ASN:HD22	1:A:26:ASN:C	1.99	0.70
2:G:87:LEU:HA	2:G:90:GLN:HG2	1.73	0.70
2:J:67:THR:CG2	2:J:71:LEU:HD12	2.21	0.70
3:L:64:TRP:CZ3	3:L:74:LEU:HD23	2.26	0.70
2:I:27:ASP:OD2	2:I:66:PRO:HD2	1.92	0.70
3:L:20:MET:HB3	3:L:57:THR:CG2	2.17	0.70
3:L:55:SER:CB	3:L:58:ARG:HB2	2.22	0.70
3:K:48:MET:O	3:K:51:VAL:HG12	1.92	0.69
3:N:13:LEU:HD21	3:N:61:LEU:CD2	2.14	0.69
3:M:86:ALA:O	3:M:89:ILE:HG23	1.91	0.69
1:D:43:VAL:CG2	2:H:48:TYR:HE1	2.04	0.69
1:F:43:VAL:HG11	2:J:48:TYR:CD1	2.27	0.69
3:K:66:MET:HE3	3:K:66:MET:HA	1.75	0.69
1:B:66:THR:HG22	1:B:67:GLN:N	2.07	0.69
2:I:25:PHE:CE1	2:J:51:PHE:HZ	2.11	0.69
2:J:67:THR:HG22	2:J:71:LEU:HD12	1.74	0.69
2:J:89:ILE:HG13	2:J:90:GLN:N	2.08	0.69
2:H:56:PHE:CE2	2:H:70:LEU:HB2	2.28	0.69
1:D:27:MET:O	1:D:28:ARG:C	2.36	0.69
1:F:85:SER:HB3	1:F:88:ARG:CB	2.20	0.69
3:N:17:CYS:O	3:N:21:ASP:HB2	1.93	0.69
1:B:29:VAL:O	1:B:33:LEU:HB2	1.93	0.69
3:N:45:ILE:O	3:N:56:ILE:CD1	2.39	0.69
2:H:11:VAL:O	2:H:14:LEU:HD12	1.92	0.68
2:I:21:LYS:HD3	2:I:99:LEU:HD21	1.75	0.68
2:J:23:SER:HA	2:J:67:THR:HG21	1.75	0.68
3:K:43:ARG:NH2	3:N:58:ARG:HD3	2.09	0.68
1:D:73:ARG:HE	1:D:73:ARG:HA	1.58	0.68
2:G:9:THR:HG22	2:G:10:TYR:H	1.58	0.68
2:G:53:ILE:HG13	2:J:16:VAL:HG21	1.75	0.68
2:J:26:ILE:HG21	2:J:35:LEU:HD23	1.74	0.68
3:L:52:GLN:O	3:L:54:VAL:HG13	1.93	0.68
3:M:9:PRO:HG2	3:M:93:TRP:CZ2	2.28	0.68
2:I:85:VAL:HA	2:I:88:LEU:HB2	1.76	0.68
3:L:19:ASN:HB3	3:L:20:MET:HE2	1.75	0.68
3:N:28:TRP:CE3	3:N:29:MET:HA	2.28	0.68
1:C:42:GLN:HA	1:F:31:ARG:HH21	1.58	0.68
1:D:38:ASN:ND2	1:D:70:PRO:HD2	2.09	0.68
1:D:25:LEU:O	1:D:30:ARG:NH1	2.27	0.68
3:L:16:LEU:HD23	3:L:16:LEU:O	1.94	0.67
3:N:71:VAL:O	3:N:75:VAL:HG23	1.94	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:122:GLU:OE1	1:D:123:HIS:CE1	2.47	0.67
3:M:17:CYS:O	3:M:21:ASP:HB2	1.94	0.67
3:N:58:ARG:NH2	3:N:62:TRP:HB2	2.10	0.67
1:D:100:ASP:HB2	2:H:12:ARG:NH1	2.09	0.67
1:E:27:MET:O	1:E:28:ARG:C	2.36	0.67
1:B:22:LEU:HD23	1:B:79:GLN:HA	1.77	0.67
3:K:9:PRO:HG2	3:K:93:TRP:CH2	2.30	0.67
1:A:111:GLU:OE2	1:A:115:LYS:HE3	1.94	0.67
1:C:103:LEU:HD23	1:C:103:LEU:C	2.18	0.67
1:D:43:VAL:HG11	2:H:48:TYR:HD1	1.60	0.67
2:J:89:ILE:HG13	2:J:90:GLN:H	1.60	0.67
1:C:22:LEU:HD23	1:C:79:GLN:HG3	1.77	0.67
3:N:28:TRP:CZ2	3:N:56:ILE:CG1	2.78	0.67
1:A:20:LEU:H	1:A:20:LEU:HD23	1.60	0.66
1:D:54:MET:HG2	1:D:78:TRP:CE2	2.29	0.66
1:B:81:ARG:CB	1:B:82:PRO:HD2	2.22	0.66
1:E:50:LEU:O	1:E:53:GLU:N	2.28	0.66
2:J:18:LEU:HD12	2:J:19:ILE:N	2.09	0.66
2:J:38:ALA:CB	2:J:87:LEU:HD21	2.26	0.66
3:M:76:ASP:HB3	3:M:80:ARG:NH1	2.09	0.66
1:D:85:SER:HB3	1:D:88:ARG:H	1.61	0.66
1:F:67:GLN:H	1:F:67:GLN:HE21	1.44	0.66
1:C:22:LEU:H	1:C:22:LEU:CD2	2.00	0.66
1:E:82:PRO:O	1:E:84:ALA:N	2.29	0.66
2:J:33:LYS:O	2:J:37:VAL:HG23	1.95	0.66
3:M:72:GLN:HA	3:M:75:VAL:HG22	1.75	0.66
1:D:43:VAL:HG11	2:H:48:TYR:CD1	2.30	0.66
2:G:41:LYS:HB3	2:G:42:PRO:CD	2.24	0.66
2:H:25:PHE:HB3	2:H:96:PRO:HG3	1.78	0.66
3:L:82:GLU:HA	3:L:84:TYR:CE1	2.31	0.66
2:J:15:ASN:ND2	2:J:18:LEU:HB3	2.10	0.66
2:J:37:VAL:HG13	3:M:10:SER:HB2	1.77	0.66
3:M:70:THR:HG21	3:M:72:GLN:NE2	2.10	0.66
3:N:84:TYR:O	3:N:88:GLN:HG2	1.95	0.66
3:K:63:TRP:O	3:K:67:ARG:HD3	1.96	0.66
1:C:20:LEU:HB2	1:C:21:PRO:HD2	1.77	0.66
1:D:67:GLN:HE21	1:D:73:ARG:HG2	1.60	0.66
1:D:20:LEU:HD12	1:D:20:LEU:N	2.10	0.65
2:H:93:PHE:HD2	3:L:66:MET:SD	2.19	0.65
2:G:55:ARG:HG3	2:G:55:ARG:NH1	2.11	0.65
3:L:63:TRP:CD1	3:L:67:ARG:CZ	2.78	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:85:SER:HB3	1:E:88:ARG:H	1.61	0.65
2:I:89:ILE:HG13	2:I:90:GLN:N	2.11	0.65
3:L:63:TRP:NE1	3:L:67:ARG:NH1	2.44	0.65
1:A:37:LEU:HD13	1:A:50:LEU:HD22	1.79	0.65
3:L:35:VAL:O	3:L:35:VAL:HG13	1.95	0.65
1:F:43:VAL:HG11	2:J:48:TYR:HE1	1.57	0.65
3:K:80:ARG:O	3:K:80:ARG:HG2	1.97	0.65
3:N:30:GLU:HG3	3:N:34:TYR:CE1	2.32	0.65
1:A:67:GLN:NE2	1:A:73:ARG:HG2	2.11	0.65
2:J:18:LEU:HD12	2:J:19:ILE:HD13	1.77	0.65
3:L:78:LEU:HD13	3:L:86:ALA:HB3	1.78	0.65
2:G:23:SER:O	2:G:27:ASP:HB2	1.97	0.65
1:D:100:ASP:HB2	2:H:12:ARG:HH12	1.61	0.65
3:L:16:LEU:CD1	3:L:90:ILE:HG21	2.27	0.65
3:M:87:ALA:O	3:M:91:LEU:HD23	1.97	0.65
1:A:33:LEU:HD13	1:A:75:LEU:HD11	1.78	0.65
1:D:62:ARG:C	1:D:64:LEU:H	2.05	0.64
3:M:24:SER:HB3	3:M:27:ASP:HB2	1.79	0.64
1:D:78:TRP:O	1:D:84:ALA:HB2	1.97	0.64
3:K:63:TRP:CD1	3:K:67:ARG:NH1	2.65	0.64
2:H:93:PHE:O	2:H:96:PRO:HD2	1.98	0.64
2:I:85:VAL:HG11	2:I:101:LEU:HD21	1.80	0.64
2:H:80:THR:CG2	2:H:83:ASP:H	2.10	0.64
3:K:20:MET:H	3:K:20:MET:CE	2.10	0.64
2:H:48:TYR:OH	2:H:74:TRP:HB2	1.98	0.64
1:E:38:ASN:ND2	1:E:71:THR:H	1.92	0.64
3:K:77:LEU:O	3:K:77:LEU:HD23	1.98	0.64
3:N:54:VAL:HG12	3:N:59:GLU:OE2	1.97	0.64
2:G:37:VAL:HG11	2:J:17:GLY:H	1.62	0.64
3:N:20:MET:O	3:N:28:TRP:CD1	2.45	0.64
3:N:28:TRP:HE3	3:N:29:MET:HA	1.60	0.64
3:N:35:VAL:HG12	3:N:37:THR:OG1	1.97	0.64
1:F:79:GLN:C	1:F:81:ARG:H	2.06	0.64
1:D:37:LEU:HB2	1:D:71:THR:CG2	2.29	0.63
1:E:53:GLU:HA	1:E:53:GLU:OE1	1.96	0.63
1:D:89:LEU:HD23	1:D:90:LEU:HD23	1.80	0.63
3:L:20:MET:H	3:L:20:MET:HE3	1.61	0.63
1:A:43:VAL:HG13	1:E:60:GLU:HG2	1.79	0.63
1:C:75:LEU:N	1:C:75:LEU:HD23	2.12	0.63
2:G:95:ALA:HA	3:K:68:GLN:NE2	2.13	0.63
1:B:58:TYR:HB2	1:E:26:ASN:HB3	1.80	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:I:33:LYS:O	2:I:37:VAL:HG23	1.98	0.63
1:A:81:ARG:HB3	1:A:82:PRO:CD	2.29	0.63
1:B:56:PHE:CE1	1:B:74:LEU:HD11	2.34	0.63
1:E:40:ARG:HB2	1:E:40:ARG:HH11	1.62	0.63
3:K:14:ASP:O	3:K:18:ARG:HG2	1.98	0.63
3:K:13:LEU:HD11	3:K:58:ARG:HG2	1.79	0.63
1:E:78:TRP:O	1:E:84:ALA:HB2	1.99	0.62
3:L:5:ILE:HD13	3:L:70:THR:O	1.99	0.62
3:N:29:MET:HE1	3:N:46:LYS:HE2	1.81	0.62
2:I:100:LEU:HB3	2:I:101:LEU:HD23	1.81	0.62
3:L:32:ALA:HB1	3:L:42:LEU:HD11	1.81	0.62
3:N:76:ASP:HB3	3:N:77:LEU:HD12	1.81	0.62
1:C:77:ALA:O	1:C:81:ARG:HD2	2.00	0.62
1:F:39:VAL:HG21	2:J:72:PHE:HZ	1.64	0.62
1:C:45:ALA:HB1	1:C:49:ALA:HB3	1.82	0.62
2:I:20:ARG:O	2:I:20:ARG:HG2	1.98	0.62
3:M:4:TYR:HB2	3:M:7:GLN:HG3	1.82	0.62
1:A:89:LEU:O	1:A:93:LEU:HG	1.99	0.62
1:D:82:PRO:O	1:D:84:ALA:N	2.33	0.62
1:F:58:TYR:CD2	2:I:16:VAL:HG22	2.35	0.62
2:I:11:VAL:HG12	2:I:79:CYS:O	1.99	0.62
1:A:43:VAL:HG21	1:E:64:LEU:CD2	2.30	0.62
2:H:50:GLN:HG2	3:K:58:ARG:HH22	1.64	0.62
2:H:85:VAL:HG12	2:H:97:ALA:CB	2.30	0.62
1:A:22:LEU:HB2	1:A:79:GLN:HG3	1.81	0.62
1:A:30:ARG:HH11	1:A:30:ARG:HG2	1.64	0.62
3:L:5:ILE:HD11	3:L:69:ALA:HB3	1.80	0.62
1:E:78:TRP:HA	1:E:81:ARG:HG3	1.82	0.62
1:E:93:LEU:HD13	1:E:101:VAL:CG1	2.29	0.62
1:B:93:LEU:HD13	1:B:101:VAL:CG2	2.30	0.61
2:H:93:PHE:HB3	2:H:96:PRO:HG2	1.81	0.61
2:J:78:ASN:O	2:J:79:CYS:O	2.17	0.61
1:A:26:ASN:HD21	1:A:28:ARG:HB3	1.64	0.61
3:M:20:MET:HE1	3:M:78:LEU:HD11	1.82	0.61
1:C:78:TRP:HA	1:C:81:ARG:HD2	1.81	0.61
2:G:101:LEU:HD23	2:G:101:LEU:O	2.00	0.61
1:E:113:CYS:C	1:E:115:LYS:H	2.08	0.61
2:H:90:GLN:HG3	2:H:91:ASN:HD22	1.65	0.61
2:I:100:LEU:CB	2:I:101:LEU:HD23	2.30	0.61
3:L:17:CYS:O	3:L:21:ASP:HB2	2.01	0.61
3:M:23:LEU:HD21	3:M:83:LEU:HD21	1.82	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:58:TYR:CD1	1:E:30:ARG:NH1	2.69	0.61
2:G:85:VAL:HB	2:G:100:LEU:HD13	1.81	0.61
3:K:79:CYS:C	3:K:81:LEU:H	2.08	0.61
2:I:22:LEU:HD12	2:I:22:LEU:C	2.26	0.61
3:N:55:SER:HB3	3:N:58:ARG:HB2	1.80	0.61
1:C:38:ASN:ND2	1:C:70:PRO:HD2	2.15	0.61
2:G:100:LEU:CD2	2:G:101:LEU:HD22	2.14	0.61
2:J:85:VAL:HG12	2:J:97:ALA:CB	2.30	0.61
3:L:87:ALA:O	3:L:91:LEU:HD23	2.00	0.61
2:G:85:VAL:HG11	2:G:101:LEU:HD21	1.83	0.61
3:K:17:CYS:O	3:K:21:ASP:HB2	2.00	0.61
1:A:29:VAL:O	1:A:33:LEU:HB2	2.01	0.60
1:A:43:VAL:HG21	1:E:64:LEU:HD23	1.83	0.60
1:C:38:ASN:ND2	1:C:71:THR:HG23	2.16	0.60
3:L:57:THR:O	3:L:61:LEU:HB2	2.01	0.60
1:C:34:SER:HA	1:C:71:THR:OG1	2.01	0.60
1:C:38:ASN:ND2	1:C:71:THR:H	1.95	0.60
2:G:92:GLU:CD	3:K:67:ARG:HG3	2.26	0.60
3:L:38:ASP:OD1	3:L:40:THR:HG22	2.00	0.60
3:N:40:THR:O	3:N:43:ARG:HB2	2.00	0.60
1:A:54:MET:HG2	1:A:92:LEU:CD1	2.31	0.60
1:D:37:LEU:HB2	1:D:71:THR:HG21	1.83	0.60
3:K:78:LEU:HD13	3:K:86:ALA:HB3	1.83	0.60
3:N:16:LEU:CD2	3:N:61:LEU:HD21	2.31	0.60
3:N:70:THR:HG22	3:N:73:GLN:NE2	2.16	0.60
1:C:85:SER:CB	1:C:88:ARG:HB2	2.28	0.60
1:F:67:GLN:O	1:F:70:PRO:HD3	2.02	0.60
3:N:35:VAL:HG13	3:N:37:THR:OG1	2.00	0.60
3:N:20:MET:HE3	3:N:20:MET:H	1.67	0.60
2:I:18:LEU:HD23	2:I:19:ILE:N	2.16	0.60
1:D:93:LEU:CD1	1:D:101:VAL:HG11	2.30	0.60
1:E:46:ASP:OD2	1:E:48:THR:OG1	2.15	0.60
1:F:37:LEU:N	1:F:37:LEU:HD23	2.16	0.60
2:H:2:MET:CG	2:H:3:GLY:N	2.48	0.60
1:A:25:LEU:HD21	1:A:109:ILE:HG23	1.83	0.60
1:D:81:ARG:HB3	1:D:82:PRO:HD2	1.84	0.60
2:G:78:ASN:O	2:G:79:CYS:O	2.20	0.60
1:B:106:GLY:N	1:B:107:PRO:HD2	2.15	0.60
2:G:77:THR:O	2:G:77:THR:CG2	2.49	0.60
3:N:18:ARG:HG2	3:N:18:ARG:HH11	1.67	0.60
3:N:70:THR:HG23	3:N:73:GLN:H	1.66	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:32:ARG:HG2	1:A:105:LEU:CD2	2.32	0.59
1:A:54:MET:HE2	1:A:74:LEU:HD11	1.84	0.59
3:M:63:TRP:O	3:M:67:ARG:NH2	2.30	0.59
1:C:101:VAL:O	1:C:105:LEU:HB2	2.03	0.59
3:L:5:ILE:CG1	3:L:69:ALA:HB3	2.32	0.59
1:A:98:ARG:O	1:A:101:VAL:HG22	2.02	0.59
1:B:62:ARG:HH11	1:B:62:ARG:HG3	1.67	0.59
1:C:93:LEU:HD13	1:C:101:VAL:HG13	1.85	0.59
3:L:28:TRP:HZ3	3:L:46:LYS:HD3	1.67	0.59
1:D:62:ARG:O	1:D:64:LEU:N	2.34	0.59
2:H:33:LYS:O	2:H:37:VAL:HG22	2.02	0.59
3:N:4:TYR:HE2	3:N:68:GLN:HG3	1.67	0.59
1:C:50:LEU:HA	1:C:96:LEU:HD11	1.85	0.59
1:F:89:LEU:O	1:F:89:LEU:HD12	2.03	0.59
2:I:37:VAL:HG11	3:L:11:TRP:N	2.16	0.59
3:M:70:THR:HG22	3:M:71:VAL:H	1.67	0.59
1:A:54:MET:HE1	1:A:74:LEU:HG	1.84	0.59
3:N:37:THR:HG22	3:N:38:ASP:N	2.13	0.59
1:C:22:LEU:HD11	1:C:78:TRP:CZ3	2.38	0.59
3:K:70:THR:HG23	3:K:73:GLN:NE2	2.18	0.59
2:G:12:ARG:NH1	2:G:13:CYS:SG	2.73	0.58
3:M:48:MET:SD	3:M:59:GLU:HG2	2.43	0.58
1:F:20:LEU:HD12	1:F:20:LEU:N	2.18	0.58
3:M:20:MET:HB3	3:M:57:THR:CG2	2.33	0.58
1:A:52:GLU:OE2	1:D:28:ARG:HB3	2.03	0.58
3:L:70:THR:OG1	3:L:73:GLN:HG3	2.02	0.58
3:N:8:LEU:HD13	3:N:12:VAL:HG12	1.86	0.58
2:J:85:VAL:HA	2:J:88:LEU:HB2	1.86	0.58
1:B:120:ALA:O	1:B:121:LEU:HD23	2.03	0.58
1:D:40:ARG:CD	1:D:46:ASP:OD2	2.52	0.58
2:H:93:PHE:C	2:H:96:PRO:HD2	2.28	0.58
2:J:6:THR:HG23	2:J:9:THR:HG23	1.85	0.58
3:N:58:ARG:HB3	3:N:59:GLU:OE1	2.03	0.58
2:G:20:ARG:HG2	2:H:55:ARG:HH21	1.69	0.58
2:G:78:ASN:O	2:G:79:CYS:C	2.46	0.58
3:K:32:ALA:O	3:K:35:VAL:HG12	2.04	0.58
1:A:69:ASP:H	1:A:70:PRO:HD3	1.68	0.58
2:H:73:ASP:OD2	2:H:73:ASP:C	2.46	0.58
3:K:90:ILE:O	3:K:90:ILE:HG13	2.04	0.58
3:M:14:ASP:O	3:M:18:ARG:HG2	2.02	0.58
3:N:9:PRO:HG3	3:N:93:TRP:CH2	2.38	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:20:LEU:H	1:D:20:LEU:CD1	2.06	0.57
2:H:92:GLU:HA	2:H:94:PHE:HE2	1.67	0.57
2:H:98:SER:HA	2:H:105:VAL:HG21	1.85	0.57
2:I:90:GLN:NE2	2:I:91:ASN:HD22	2.02	0.57
1:E:73:ARG:HH11	1:E:73:ARG:HG2	1.68	0.57
1:E:107:PRO:HG2	1:E:108:SER:N	2.19	0.57
2:I:11:VAL:CG1	2:I:79:CYS:SG	2.91	0.57
2:I:78:ASN:O	2:I:79:CYS:C	2.46	0.57
3:L:63:TRP:NE1	3:L:67:ARG:CZ	2.67	0.57
3:M:72:GLN:HA	3:M:75:VAL:CG2	2.34	0.57
2:H:81:VAL:HG12	2:H:101:LEU:HD21	1.86	0.57
2:G:19:ILE:C	2:G:21:LYS:H	2.12	0.57
3:N:45:ILE:HG21	3:N:60:LEU:HD12	1.86	0.57
1:C:35:LEU:O	1:D:59:LEU:CD2	2.52	0.57
1:D:38:ASN:ND2	1:D:71:THR:H	1.99	0.57
1:D:50:LEU:O	1:D:53:GLU:N	2.37	0.57
1:E:43:VAL:HG11	2:I:52:HIS:CD2	2.39	0.57
3:L:90:ILE:C	3:L:91:LEU:HD22	2.30	0.57
1:B:58:TYR:OH	1:E:76:ASP:OD2	2.23	0.57
1:F:53:GLU:C	1:F:55:ASP:H	2.12	0.57
2:G:95:ALA:N	2:G:96:PRO:HD2	2.20	0.57
2:H:47:ARG:HG2	2:H:48:TYR:CE2	2.40	0.57
2:H:59:LEU:HA	2:H:62:THR:HG23	1.87	0.57
3:M:4:TYR:HD1	3:M:7:GLN:OE1	1.87	0.57
1:D:50:LEU:HD21	1:D:54:MET:CE	2.35	0.57
3:K:9:PRO:HG2	3:K:93:TRP:CZ3	2.40	0.57
3:N:78:LEU:HD13	3:N:86:ALA:HB3	1.86	0.57
1:B:58:TYR:CE1	1:E:30:ARG:NH1	2.72	0.57
1:F:46:ASP:OD1	1:F:48:THR:OG1	2.20	0.57
2:G:50:GLN:HA	2:J:16:VAL:CG2	2.32	0.57
2:H:78:ASN:O	2:H:79:CYS:O	2.22	0.57
2:H:85:VAL:O	2:H:89:ILE:HG13	2.04	0.57
2:J:50:GLN:NE2	3:M:58:ARG:HH22	1.99	0.57
3:L:54:VAL:O	3:L:54:VAL:HG23	2.05	0.57
2:J:15:ASN:OD1	2:J:15:ASN:N	2.38	0.57
2:J:18:LEU:CD1	2:J:19:ILE:HD13	2.35	0.57
3:N:23:LEU:HD13	3:N:23:LEU:O	2.05	0.57
1:E:62:ARG:C	1:E:64:LEU:H	2.11	0.56
1:E:83:GLY:O	1:E:85:SER:N	2.38	0.56
2:I:19:ILE:C	2:I:21:LYS:H	2.12	0.56
3:K:43:ARG:HH21	3:N:58:ARG:HD3	1.69	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:69:ASP:N	1:C:70:PRO:HD3	2.20	0.56
2:G:6:THR:HG22	2:G:7:PRO:HD2	1.86	0.56
2:I:54:ARG:HH12	3:L:58:ARG:HD3	1.70	0.56
3:L:63:TRP:HE1	3:L:67:ARG:NH1	2.02	0.56
1:A:56:PHE:CD2	1:A:60:GLU:HG2	2.40	0.56
1:B:39:VAL:HG23	1:B:39:VAL:O	2.05	0.56
1:C:93:LEU:HD13	1:C:101:VAL:HG11	1.86	0.56
2:I:74:TRP:HZ3	2:I:84:LEU:HD13	1.70	0.56
3:N:81:LEU:O	3:N:81:LEU:HD22	2.06	0.56
3:N:87:ALA:HA	3:N:90:ILE:HG13	1.87	0.56
1:D:93:LEU:HD13	1:D:101:VAL:CG1	2.32	0.56
1:E:22:LEU:C	1:E:24:ALA:H	2.14	0.56
2:I:89:ILE:HG13	2:I:90:GLN:H	1.71	0.56
1:B:41:THR:HG23	1:B:41:THR:O	2.06	0.56
2:J:78:ASN:O	2:J:79:CYS:C	2.47	0.56
3:L:4:TYR:HB2	3:L:7:GLN:OE1	2.05	0.56
3:L:79:CYS:C	3:L:81:LEU:H	2.14	0.56
3:N:55:SER:HB3	3:N:58:ARG:H	1.71	0.56
1:B:40:ARG:NH2	1:E:31:ARG:HH22	2.00	0.56
1:B:98:ARG:HD3	1:F:80:GLY:CA	2.36	0.56
2:H:47:ARG:HG2	2:H:48:TYR:CD2	2.41	0.56
2:H:50:GLN:HG2	3:K:58:ARG:NH2	2.21	0.56
1:A:52:GLU:OE1	1:D:27:MET:HB2	2.05	0.56
1:A:105:LEU:O	1:A:109:ILE:HG13	2.06	0.56
1:C:90:LEU:HD12	1:C:102:LEU:HD23	1.88	0.56
1:F:36:PHE:CD2	1:F:101:VAL:HG23	2.41	0.56
2:I:6:THR:HG23	2:I:9:THR:CG2	2.35	0.56
2:I:78:ASN:O	2:I:79:CYS:O	2.23	0.56
3:L:36:ILE:HG22	3:L:36:ILE:O	2.06	0.56
3:N:19:ASN:HB2	3:N:20:MET:CE	2.35	0.56
1:C:58:TYR:CB	1:F:26:ASN:HB3	2.33	0.56
1:E:57:GLU:HG3	1:E:60:GLU:H	1.70	0.56
1:F:71:THR:O	1:F:75:LEU:HG	2.05	0.56
2:G:24:ASP:HB3	2:H:51:PHE:HD1	1.71	0.56
2:H:78:ASN:O	2:H:79:CYS:C	2.49	0.56
2:J:38:ALA:HB1	2:J:87:LEU:HD21	1.88	0.56
3:K:84:TYR:O	3:K:88:GLN:CG	2.52	0.56
1:C:79:GLN:C	1:C:81:ARG:H	2.14	0.56
1:D:113:CYS:C	1:D:115:LYS:H	2.14	0.56
1:F:39:VAL:O	1:F:39:VAL:CG2	2.49	0.56
2:G:39:ILE:HD13	2:G:74:TRP:CD2	2.41	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:27:ASP:OD2	2:H:66:PRO:HD2	2.06	0.55
3:M:64:TRP:HA	3:M:67:ARG:NH2	2.20	0.55
1:B:27:MET:O	1:B:28:ARG:C	2.49	0.55
1:C:61:ILE:HD12	1:F:26:ASN:HD22	1.71	0.55
1:D:33:LEU:HD13	1:D:105:LEU:CD1	2.32	0.55
1:D:54:MET:HB3	1:D:78:TRP:CD1	2.41	0.55
1:E:122:GLU:HB2	1:E:123:HIS:CE1	2.41	0.55
3:L:63:TRP:CD1	3:L:67:ARG:NH2	2.75	0.55
2:J:2:MET:HE3	2:J:4:PRO:HD2	1.88	0.55
3:M:35:VAL:HG12	3:M:37:THR:OG1	2.06	0.55
2:I:25:PHE:O	2:I:28:PRO:HD2	2.07	0.55
2:J:11:VAL:HG21	2:J:79:CYS:SG	2.46	0.55
3:M:79:CYS:C	3:M:81:LEU:H	2.14	0.55
1:B:93:LEU:HD13	1:B:101:VAL:HG23	1.89	0.55
1:D:40:ARG:HB2	1:D:40:ARG:CZ	2.36	0.55
2:G:54:ARG:HH21	2:J:72:PHE:HB2	1.72	0.55
2:G:74:TRP:CG	2:G:74:TRP:O	2.59	0.55
3:L:2:ALA:O	3:L:3:CYS:HB2	2.07	0.55
3:N:70:THR:HG22	3:N:73:GLN:CD	2.32	0.55
1:B:79:GLN:NE2	1:B:80:GLY:N	2.55	0.55
1:E:54:MET:HE1	1:E:74:LEU:HG	1.87	0.55
1:F:54:MET:HE2	1:F:78:TRP:CE3	2.42	0.55
1:F:122:GLU:C	1:F:123:HIS:ND1	2.65	0.55
2:G:14:LEU:HD11	2:G:81:VAL:HG22	1.88	0.55
2:H:19:ILE:C	2:H:21:LYS:H	2.15	0.55
3:K:49:GLU:HB2	3:K:56:ILE:HD13	1.88	0.55
3:M:5:ILE:HG12	3:M:70:THR:O	2.06	0.55
3:N:64:TRP:HA	3:N:67:ARG:HH21	1.70	0.55
2:J:54:ARG:HH21	3:M:62:TRP:CG	2.24	0.55
3:K:11:TRP:CE2	3:K:12:VAL:HG23	2.42	0.55
1:F:51:ALA:HB1	1:F:61:ILE:HD13	1.87	0.55
2:G:28:PRO:HG2	3:K:62:TRP:CZ2	2.41	0.55
1:B:50:LEU:HD11	1:B:54:MET:HE3	1.87	0.55
1:D:35:LEU:HB3	1:E:59:LEU:HD22	1.89	0.55
3:M:5:ILE:HG13	3:M:69:ALA:HB3	1.89	0.55
3:K:24:SER:HB2	3:K:27:ASP:CG	2.32	0.54
1:A:101:VAL:HG23	1:A:102:LEU:HD12	1.88	0.54
1:C:62:ARG:C	1:C:64:LEU:H	2.16	0.54
1:F:40:ARG:CD	1:F:46:ASP:OD2	2.56	0.54
3:M:25:GLU:O	3:M:29:MET:HB2	2.07	0.54
1:A:46:ASP:OD2	1:A:48:THR:OG1	2.24	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:28:ARG:CG	1:D:29:VAL:N	2.70	0.54
1:D:90:LEU:HD21	1:D:109:ILE:HD13	1.89	0.54
2:H:73:ASP:OD2	2:H:73:ASP:O	2.25	0.54
2:J:19:ILE:C	2:J:21:LYS:H	2.14	0.54
1:C:43:VAL:HG12	2:G:73:ASP:OD2	2.07	0.54
3:L:81:LEU:HD13	3:L:81:LEU:O	2.07	0.54
3:L:85:ARG:HA	3:L:88:GLN:NE2	2.21	0.54
1:A:88:ARG:HA	1:A:91:GLU:HG2	1.90	0.54
1:C:20:LEU:HB2	1:C:21:PRO:CD	2.35	0.54
2:H:23:SER:OG	2:H:65:SER:OG	2.22	0.54
3:N:79:CYS:C	3:N:81:LEU:H	2.15	0.54
1:A:76:ASP:C	1:A:78:TRP:H	2.16	0.54
1:F:19:MET:HE2	1:F:86:VAL:HG12	1.90	0.54
1:F:83:GLY:O	1:F:85:SER:N	2.41	0.54
2:H:95:ALA:CA	3:L:68:GLN:NE2	2.64	0.54
2:I:59:LEU:HD13	2:I:69:GLU:HG3	1.90	0.54
1:A:98:ARG:CZ	1:E:80:GLY:HA3	2.38	0.54
2:J:2:MET:CE	2:J:4:PRO:HD2	2.38	0.54
3:K:64:TRP:HA	3:K:67:ARG:NE	2.20	0.54
3:N:11:TRP:C	3:N:11:TRP:CD1	2.86	0.54
1:B:50:LEU:O	1:B:53:GLU:N	2.39	0.54
1:C:103:LEU:HD23	1:C:104:GLU:N	2.23	0.54
2:G:85:VAL:CG2	2:G:97:ALA:HB1	2.37	0.54
1:A:68:ALA:O	1:A:69:ASP:HB3	2.06	0.54
1:D:39:VAL:O	1:D:39:VAL:CG2	2.56	0.54
1:E:93:LEU:HD13	1:E:101:VAL:HG13	1.90	0.54
3:K:63:TRP:NE1	3:K:67:ARG:NH1	2.56	0.54
1:E:77:ALA:O	1:E:81:ARG:HD3	2.07	0.53
3:L:5:ILE:CD1	3:L:69:ALA:HB3	2.36	0.53
3:L:76:ASP:HB3	3:L:80:ARG:HH21	1.73	0.53
3:N:15:ASP:HA	3:N:18:ARG:HG3	1.90	0.53
1:C:47:TRP:CB	1:C:71:THR:HG22	2.37	0.53
1:C:73:ARG:NE	1:C:73:ARG:CA	2.70	0.53
1:D:46:ASP:CG	1:D:48:THR:HG23	2.32	0.53
1:F:39:VAL:HG21	2:J:72:PHE:CZ	2.42	0.53
3:L:24:SER:O	3:L:26:TRP:N	2.42	0.53
1:B:73:ARG:HA	1:B:73:ARG:NE	2.24	0.53
1:D:58:TYR:OH	2:G:72:PHE:HB2	2.07	0.53
2:H:94:PHE:N	2:H:94:PHE:CD2	2.76	0.53
3:K:64:TRP:CZ3	3:K:74:LEU:HD23	2.44	0.53
3:M:28:TRP:CZ3	3:M:46:LYS:HD2	2.43	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:27:MET:O	1:A:28:ARG:C	2.51	0.53
1:B:59:LEU:HD11	1:E:23:ALA:HB1	1.90	0.53
1:F:36:PHE:C	1:F:37:LEU:HD23	2.34	0.53
2:H:94:PHE:N	2:H:94:PHE:HD2	2.07	0.53
2:J:38:ALA:HB3	2:J:87:LEU:HD21	1.90	0.53
1:B:24:ALA:O	1:B:115:LYS:HE2	2.08	0.53
1:C:109:ILE:C	1:C:111:GLU:H	2.17	0.53
1:F:58:TYR:CE2	2:I:16:VAL:HG13	2.44	0.53
2:G:25:PHE:CE1	2:H:51:PHE:HZ	2.26	0.53
1:C:73:ARG:NE	1:C:73:ARG:HA	2.24	0.53
1:E:109:ILE:O	1:E:110:GLU:C	2.51	0.53
1:E:122:GLU:C	1:E:123:HIS:ND1	2.67	0.53
3:K:16:LEU:CB	3:K:89:ILE:HD11	2.28	0.53
3:N:80:ARG:HB2	3:N:80:ARG:NH1	2.23	0.53
1:C:22:LEU:HD21	1:C:79:GLN:HA	1.90	0.53
1:D:78:TRP:HZ2	1:D:88:ARG:HG2	1.74	0.53
3:N:18:ARG:HG2	3:N:18:ARG:NH1	2.22	0.53
1:F:73:ARG:NE	1:F:73:ARG:CA	2.72	0.53
1:F:98:ARG:NH1	1:F:100:ASP:OD2	2.42	0.53
2:J:10:TYR:HB2	2:J:13:CYS:SG	2.49	0.53
1:B:35:LEU:HD23	1:B:35:LEU:N	2.23	0.53
1:E:53:GLU:OE2	1:E:95:LYS:HE2	2.09	0.53
1:F:54:MET:HB3	1:F:78:TRP:CD1	2.43	0.53
1:F:98:ARG:HH11	1:F:98:ARG:HB3	1.73	0.53
2:H:50:GLN:HE21	3:K:58:ARG:HH21	1.55	0.53
3:N:9:PRO:HG3	3:N:93:TRP:CZ3	2.43	0.53
1:A:38:ASN:C	1:A:38:ASN:HD22	2.16	0.53
1:F:100:ASP:OD2	1:F:101:VAL:N	2.42	0.53
3:K:16:LEU:HD21	3:K:61:LEU:HD21	1.91	0.53
3:M:15:ASP:HA	3:M:18:ARG:CG	2.38	0.53
1:B:76:ASP:C	1:B:78:TRP:H	2.17	0.52
1:C:53:GLU:C	1:C:55:ASP:H	2.17	0.52
1:E:25:LEU:HD13	1:E:86:VAL:HG21	1.87	0.52
2:H:85:VAL:HA	2:H:88:LEU:HB2	1.91	0.52
2:H:74:TRP:HZ3	2:H:84:LEU:HD12	1.74	0.52
3:K:31:PHE:C	3:K:31:PHE:CD2	2.87	0.52
3:M:92:ASN:N	3:M:92:ASN:HD22	2.05	0.52
1:C:38:ASN:HD21	1:C:71:THR:N	2.00	0.52
1:C:40:ARG:HD2	1:C:46:ASP:OD2	2.10	0.52
1:E:67:GLN:H	1:E:67:GLN:CD	2.17	0.52
3:K:43:ARG:NH2	3:N:58:ARG:CD	2.72	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:N:28:TRP:HE3	3:N:29:MET:CA	2.22	0.52
1:B:98:ARG:HD3	1:F:80:GLY:HA2	1.91	0.52
1:C:42:GLN:NE2	2:G:59:LEU:HD11	2.23	0.52
2:J:107:LEU:HD12	2:J:107:LEU:N	2.23	0.52
3:K:66:MET:CE	3:K:66:MET:HA	2.38	0.52
3:N:24:SER:HB3	3:N:27:ASP:HB2	1.92	0.52
1:E:122:GLU:HB2	1:E:123:HIS:ND1	2.25	0.52
2:G:10:TYR:HB2	2:G:13:CYS:SG	2.50	0.52
2:G:95:ALA:HB3	3:K:66:MET:CE	2.39	0.52
2:H:95:ALA:HB3	2:H:96:PRO:HD3	1.92	0.52
3:M:48:MET:O	3:M:51:VAL:HG12	2.10	0.52
1:A:22:LEU:CB	1:A:79:GLN:HG3	2.39	0.52
1:A:90:LEU:C	1:A:90:LEU:HD13	2.35	0.52
1:D:61:ILE:O	1:D:64:LEU:HB2	2.09	0.52
1:D:120:ALA:O	1:D:121:LEU:HD23	2.10	0.52
1:E:93:LEU:HB3	1:E:102:LEU:HD21	1.90	0.52
1:F:27:MET:O	1:F:28:ARG:C	2.52	0.52
2:H:26:ILE:HG22	2:H:32:TRP:HB3	1.90	0.52
2:H:74:TRP:CG	2:H:74:TRP:O	2.63	0.52
1:B:64:LEU:C	1:B:66:THR:H	2.17	0.52
1:D:117:ILE:O	1:D:118:ALA:C	2.53	0.52
2:I:96:PRO:O	2:I:99:LEU:HB3	2.10	0.52
3:L:5:ILE:HG13	3:L:69:ALA:HB3	1.91	0.52
1:C:73:ARG:HA	1:C:73:ARG:HE	1.75	0.52
3:K:20:MET:CB	3:K:57:THR:HG21	2.40	0.52
3:K:34:TYR:CD2	3:K:81:LEU:HD21	2.45	0.52
3:L:5:ILE:O	3:L:8:LEU:HD23	2.10	0.52
1:A:58:TYR:HB2	1:D:26:ASN:HB3	1.91	0.52
1:A:106:GLY:N	1:A:107:PRO:HD2	2.25	0.52
1:C:33:LEU:O	1:C:37:LEU:HD12	2.10	0.52
1:E:25:LEU:O	1:E:30:ARG:NH1	2.43	0.52
1:B:73:ARG:NH1	1:B:73:ARG:HB2	2.25	0.52
1:C:109:ILE:C	1:C:111:GLU:N	2.65	0.52
2:G:19:ILE:HD11	2:G:71:LEU:HB3	1.92	0.52
2:G:27:ASP:OD2	2:G:66:PRO:HD2	2.09	0.52
3:N:9:PRO:HG2	3:N:12:VAL:HB	1.92	0.52
3:N:29:MET:HE1	3:N:46:LYS:CE	2.40	0.52
1:C:58:TYR:CE1	1:F:30:ARG:NH2	2.78	0.51
1:D:117:ILE:O	1:D:119:ALA:O	2.27	0.51
2:G:45:ASP:OD2	2:G:45:ASP:N	2.42	0.51
3:K:89:ILE:HG13	3:K:90:ILE:N	2.25	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:106:GLY:N	1:E:107:PRO:HD2	2.25	0.51
1:F:43:VAL:HG13	2:J:73:ASP:OD1	2.11	0.51
1:F:54:MET:HE2	1:F:78:TRP:CZ3	2.44	0.51
3:M:63:TRP:C	3:M:67:ARG:HH21	2.17	0.51
3:M:76:ASP:HB3	3:M:80:ARG:CZ	2.39	0.51
3:N:75:VAL:CG1	3:N:91:LEU:HD21	2.40	0.51
1:A:114:GLN:HG2	1:A:118:ALA:HB2	1.93	0.51
1:C:67:GLN:HG3	1:C:73:ARG:HG2	1.90	0.51
1:D:61:ILE:HD12	2:G:16:VAL:HG11	1.92	0.51
1:D:62:ARG:C	1:D:64:LEU:N	2.66	0.51
1:D:122:GLU:HB2	1:D:123:HIS:ND1	2.25	0.51
2:G:53:ILE:HG13	2:J:16:VAL:CG2	2.40	0.51
2:J:22:LEU:HD23	2:J:22:LEU:C	2.34	0.51
1:B:99:ASP:O	1:B:100:ASP:C	2.53	0.51
1:D:51:ALA:HB2	1:D:74:LEU:HD22	1.93	0.51
1:D:73:ARG:HA	1:D:73:ARG:NE	2.26	0.51
1:F:37:LEU:O	1:F:47:TRP:HD1	1.94	0.51
1:F:78:TRP:CG	1:F:78:TRP:O	2.63	0.51
3:K:71:VAL:O	3:K:75:VAL:HG23	2.10	0.51
3:L:37:THR:HG21	3:L:42:LEU:HD23	1.91	0.51
3:L:84:TYR:N	3:L:84:TYR:CD1	2.76	0.51
1:A:39:VAL:HG21	1:B:59:LEU:HD23	1.92	0.51
1:B:49:ALA:O	1:B:53:GLU:HG2	2.10	0.51
3:M:62:TRP:CZ2	3:M:66:MET:HE3	2.46	0.51
1:A:101:VAL:HG23	1:A:102:LEU:CD1	2.40	0.51
1:B:58:TYR:CE1	1:E:30:ARG:NE	2.79	0.51
1:E:64:LEU:C	1:E:66:THR:N	2.69	0.51
1:E:117:ILE:O	1:E:118:ALA:C	2.52	0.51
1:F:56:PHE:CE1	1:F:74:LEU:CD1	2.93	0.51
3:N:5:ILE:HG22	3:N:6:TYR:N	2.25	0.51
1:A:30:ARG:HG2	1:A:30:ARG:NH1	2.24	0.51
1:A:30:ARG:HH22	1:A:73:ARG:HH21	1.57	0.51
1:A:122:GLU:HG3	1:A:123:HIS:N	2.20	0.51
1:D:101:VAL:HG12	1:D:102:LEU:N	2.25	0.51
1:E:83:GLY:O	1:E:84:ALA:C	2.54	0.51
2:I:2:MET:CG	2:I:3:GLY:H	2.23	0.51
1:B:54:MET:SD	1:B:74:LEU:HD21	2.51	0.51
1:C:41:THR:HG22	2:G:72:PHE:CZ	2.45	0.51
1:C:64:LEU:C	1:C:66:THR:H	2.19	0.51
1:D:50:LEU:HD21	1:D:54:MET:HE1	1.92	0.51
3:N:2:ALA:HB2	3:N:93:TRP:HE1	1.75	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:47:TRP:HB3	1:D:71:THR:HG22	1.92	0.51
1:A:56:PHE:HD2	1:A:60:GLU:HG2	1.76	0.50
1:C:85:SER:OG	1:C:88:ARG:N	2.39	0.50
2:I:85:VAL:CG1	2:I:101:LEU:HD21	2.41	0.50
3:M:74:LEU:O	3:M:78:LEU:HG	2.12	0.50
1:D:22:LEU:C	1:D:24:ALA:H	2.18	0.50
1:D:83:GLY:O	1:D:85:SER:N	2.43	0.50
1:E:53:GLU:OE1	1:E:53:GLU:CA	2.60	0.50
1:E:81:ARG:CB	1:E:82:PRO:CD	2.89	0.50
1:E:107:PRO:CG	1:E:108:SER:H	2.23	0.50
1:E:22:LEU:C	1:E:24:ALA:N	2.67	0.50
1:F:22:LEU:HD23	1:F:86:VAL:HG23	1.92	0.50
2:G:11:VAL:HG22	2:G:84:LEU:HD23	1.93	0.50
2:H:35:LEU:HD22	2:H:70:LEU:HD21	1.92	0.50
2:H:95:ALA:N	2:H:96:PRO:CD	2.74	0.50
2:J:50:GLN:NE2	3:M:58:ARG:NH2	2.51	0.50
3:M:15:ASP:HA	3:M:18:ARG:HG3	1.94	0.50
1:B:51:ALA:CB	1:B:61:ILE:HD13	2.42	0.50
1:B:103:LEU:O	1:B:104:GLU:HG2	2.11	0.50
1:D:40:ARG:HH11	1:D:40:ARG:CB	2.15	0.50
1:D:64:LEU:C	1:D:66:THR:N	2.70	0.50
1:F:25:LEU:HD22	1:F:109:ILE:HG23	1.94	0.50
1:F:82:PRO:O	1:F:84:ALA:N	2.44	0.50
2:G:92:GLU:HG2	2:G:94:PHE:HE2	1.77	0.50
2:G:93:PHE:C	2:G:96:PRO:HD2	2.36	0.50
3:K:5:ILE:HG12	3:K:70:THR:O	2.12	0.50
1:A:64:LEU:C	1:A:66:THR:H	2.18	0.50
1:B:38:ASN:HD21	1:C:63:GLN:HG3	1.77	0.50
1:E:62:ARG:O	1:E:64:LEU:N	2.44	0.50
1:F:67:GLN:CG	1:F:73:ARG:HG3	2.42	0.50
1:B:62:ARG:HG3	1:B:62:ARG:NH1	2.26	0.50
1:C:78:TRP:CG	1:C:78:TRP:O	2.64	0.50
1:C:82:PRO:O	1:C:84:ALA:N	2.45	0.50
1:F:98:ARG:NH1	1:F:98:ARG:HB3	2.26	0.50
3:L:66:MET:CA	3:L:66:MET:HE2	2.42	0.50
3:M:89:ILE:HG12	3:M:90:ILE:HG22	1.93	0.50
1:C:37:LEU:HB2	1:C:71:THR:CG2	2.41	0.50
1:E:73:ARG:HG2	1:E:73:ARG:NH1	2.27	0.50
1:F:64:LEU:C	1:F:66:THR:H	2.19	0.50
2:I:101:LEU:N	2:I:101:LEU:CD2	2.74	0.50
3:L:28:TRP:CZ3	3:L:46:LYS:HD3	2.46	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:M:5:ILE:O	3:M:8:LEU:HD23	2.12	0.50
1:A:20:LEU:HD23	1:A:20:LEU:N	2.26	0.50
1:F:34:SER:HB3	2:G:55:ARG:HH21	1.76	0.50
2:G:37:VAL:HG11	2:J:17:GLY:N	2.26	0.50
2:I:85:VAL:HG11	2:I:101:LEU:CD2	2.42	0.50
2:J:10:TYR:HD1	2:J:78:ASN:O	1.95	0.50
2:J:67:THR:HG23	2:J:71:LEU:HD12	1.94	0.50
3:K:5:ILE:CD1	3:K:69:ALA:HB3	2.42	0.50
2:G:37:VAL:HG21	2:J:17:GLY:CA	2.42	0.49
1:B:73:ARG:HB2	1:B:73:ARG:CZ	2.41	0.49
1:C:114:GLN:HA	1:C:114:GLN:OE1	2.11	0.49
1:D:22:LEU:C	1:D:24:ALA:N	2.69	0.49
2:G:85:VAL:HG21	2:G:101:LEU:HD21	1.94	0.49
1:F:50:LEU:O	1:F:51:ALA:C	2.55	0.49
1:F:53:GLU:HG3	1:F:96:LEU:HD21	1.92	0.49
1:F:56:PHE:CZ	1:F:78:TRP:HB2	2.47	0.49
2:G:92:GLU:HG2	2:G:94:PHE:CE2	2.48	0.49
2:H:91:ASN:O	2:H:92:GLU:HB2	2.12	0.49
3:K:38:ASP:OD2	3:K:38:ASP:C	2.55	0.49
3:K:70:THR:N	3:K:73:GLN:HE21	2.06	0.49
3:L:55:SER:HB2	3:L:58:ARG:CB	2.35	0.49
3:N:62:TRP:CH2	3:N:66:MET:HE3	2.47	0.49
1:E:20:LEU:H	1:E:20:LEU:CD1	1.98	0.49
1:F:34:SER:HB3	2:G:55:ARG:NH2	2.27	0.49
3:M:46:LYS:HZ2	3:M:50:TRP:HZ2	1.61	0.49
3:M:70:THR:CG2	3:M:72:GLN:NE2	2.74	0.49
3:N:63:TRP:CZ2	3:N:67:ARG:NH1	2.80	0.49
1:A:69:ASP:N	1:A:70:PRO:CD	2.75	0.49
1:E:35:LEU:O	1:F:59:LEU:CD2	2.61	0.49
2:G:98:SER:HA	2:G:105:VAL:HG21	1.94	0.49
1:F:109:ILE:C	1:F:111:GLU:H	2.19	0.49
2:H:80:THR:HG23	2:H:83:ASP:N	2.23	0.49
2:I:93:PHE:HB3	2:I:96:PRO:CG	2.43	0.49
2:J:37:VAL:HG11	3:M:10:SER:C	2.37	0.49
2:J:41:LYS:HA	2:J:41:LYS:CE	2.34	0.49
2:J:101:LEU:HD23	2:J:101:LEU:N	2.27	0.49
3:M:51:VAL:O	3:M:52:GLN:C	2.54	0.49
3:N:20:MET:H	3:N:20:MET:CE	2.24	0.49
1:B:73:ARG:NE	1:B:73:ARG:CA	2.75	0.49
1:C:27:MET:O	1:C:28:ARG:C	2.55	0.49
1:E:93:LEU:HD13	1:E:101:VAL:HG11	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:109:ILE:C	1:F:111:GLU:N	2.70	0.49
2:I:90:GLN:NE2	2:I:91:ASN:ND2	2.60	0.49
1:A:19:MET:HB2	1:A:20:LEU:HD23	1.95	0.49
1:E:104:GLU:C	1:E:105:LEU:HD23	2.37	0.49
1:F:67:GLN:H	1:F:67:GLN:CD	2.19	0.49
2:H:68:SER:OG	2:H:69:GLU:N	2.46	0.49
2:H:82:GLY:O	2:H:84:LEU:N	2.46	0.49
2:J:11:VAL:C	2:J:13:CYS:H	2.21	0.49
1:C:49:ALA:O	1:C:52:GLU:HB3	2.12	0.49
3:K:5:ILE:HD11	3:K:69:ALA:CB	2.42	0.49
3:L:49:GLU:CA	3:L:56:ILE:HD13	2.43	0.49
3:L:91:LEU:C	3:L:93:TRP:H	2.21	0.49
3:M:2:ALA:O	3:M:3:CYS:HB2	2.12	0.49
1:A:22:LEU:HD22	1:A:79:GLN:HA	1.95	0.48
1:E:98:ARG:O	1:E:101:VAL:HG12	2.13	0.48
2:H:101:LEU:C	2:H:103:ASP:H	2.20	0.48
3:L:19:ASN:O	3:L:21:ASP:N	2.46	0.48
1:B:107:PRO:HG2	1:B:108:SER:H	1.78	0.48
2:I:10:TYR:O	2:I:13:CYS:HB2	2.13	0.48
2:I:52:HIS:HA	2:I:55:ARG:HG3	1.95	0.48
3:N:59:GLU:O	3:N:63:TRP:HB3	2.13	0.48
1:B:64:LEU:C	1:B:66:THR:N	2.69	0.48
1:C:119:ALA:C	1:C:121:LEU:N	2.72	0.48
1:F:122:GLU:C	1:F:123:HIS:CG	2.91	0.48
2:I:85:VAL:HG23	2:I:86:ASP:N	2.27	0.48
2:J:25:PHE:HD2	2:J:96:PRO:HA	1.78	0.48
3:L:5:ILE:CD1	3:L:70:THR:O	2.60	0.48
1:A:80:GLY:O	1:A:81:ARG:HG2	2.13	0.48
1:C:40:ARG:CD	1:C:46:ASP:OD2	2.62	0.48
1:E:43:VAL:CG2	2:I:48:TYR:CE1	2.90	0.48
1:E:47:TRP:HB3	1:E:71:THR:CG2	2.43	0.48
1:E:119:ALA:C	1:E:121:LEU:N	2.70	0.48
2:G:55:ARG:NH1	2:G:55:ARG:CG	2.76	0.48
2:I:93:PHE:HB3	2:I:96:PRO:HG3	1.95	0.48
1:B:93:LEU:HD13	1:B:101:VAL:HG21	1.95	0.48
1:C:78:TRP:O	1:C:78:TRP:CD2	2.66	0.48
1:C:119:ALA:C	1:C:121:LEU:H	2.22	0.48
2:G:25:PHE:O	2:G:28:PRO:HD2	2.12	0.48
3:K:49:GLU:CA	3:K:56:ILE:HD13	2.43	0.48
1:E:39:VAL:HG11	1:F:62:ARG:HG2	1.95	0.48
2:H:6:THR:C	2:H:8:SER:H	2.20	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:J:6:THR:HG23	2:J:9:THR:CG2	2.43	0.48
3:K:20:MET:HB3	3:K:57:THR:CG2	2.44	0.48
3:K:79:CYS:C	3:K:81:LEU:N	2.71	0.48
3:N:51:VAL:O	3:N:52:GLN:C	2.56	0.48
1:F:101:VAL:O	1:F:105:LEU:HB2	2.12	0.48
3:M:32:ALA:HB1	3:M:42:LEU:CD1	2.41	0.48
2:I:83:ASP:O	2:I:87:LEU:HB2	2.13	0.48
3:L:19:ASN:O	3:L:20:MET:C	2.56	0.48
3:M:20:MET:HE1	3:M:78:LEU:CD1	2.44	0.48
1:A:75:LEU:O	1:A:79:GLN:HB2	2.13	0.48
1:C:53:GLU:HB3	1:C:92:LEU:HD13	1.95	0.48
1:D:67:GLN:NE2	1:D:73:ARG:HG2	2.29	0.48
2:I:41:LYS:HB3	2:I:42:PRO:CD	2.44	0.48
2:J:11:VAL:CG2	2:J:79:CYS:O	2.45	0.48
3:L:3:CYS:HB3	3:L:7:GLN:HE22	1.78	0.48
3:N:17:CYS:HB3	3:N:55:SER:OG	2.14	0.48
3:N:23:LEU:HD21	3:N:85:ARG:NH2	2.29	0.48
1:F:22:LEU:O	1:F:24:ALA:N	2.47	0.48
1:F:67:GLN:HG3	1:F:73:ARG:HG3	1.96	0.48
2:G:101:LEU:C	2:G:103:ASP:H	2.21	0.48
2:I:6:THR:HG23	2:I:9:THR:HG23	1.96	0.48
2:I:93:PHE:C	2:I:96:PRO:HD2	2.38	0.48
1:A:64:LEU:C	1:A:66:THR:N	2.71	0.47
1:E:62:ARG:C	1:E:64:LEU:N	2.71	0.47
1:F:62:ARG:C	1:F:64:LEU:H	2.22	0.47
2:G:85:VAL:O	2:G:89:ILE:HG13	2.14	0.47
3:K:52:GLN:O	3:K:53:GLY:C	2.57	0.47
1:D:45:ALA:HB1	1:D:49:ALA:HB3	1.96	0.47
2:G:69:GLU:O	2:G:73:ASP:HB2	2.14	0.47
2:I:100:LEU:HB3	2:I:101:LEU:CD2	2.44	0.47
3:K:2:ALA:O	3:K:3:CYS:HB2	2.13	0.47
3:N:5:ILE:HD11	3:N:64:TRP:CE3	2.49	0.47
1:B:64:LEU:O	1:B:66:THR:N	2.47	0.47
1:B:114:GLN:HA	1:B:117:ILE:CG2	2.18	0.47
1:B:59:LEU:HD11	1:E:23:ALA:CB	2.44	0.47
1:E:117:ILE:O	1:E:119:ALA:O	2.31	0.47
1:F:119:ALA:C	1:F:121:LEU:H	2.23	0.47
2:J:21:LYS:HD3	2:J:99:LEU:HD21	1.95	0.47
3:K:45:ILE:CD1	3:K:60:LEU:HA	2.44	0.47
3:L:76:ASP:HB3	3:L:80:ARG:NH2	2.30	0.47
3:M:24:SER:O	3:M:26:TRP:N	2.47	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:50:LEU:CD1	1:C:54:MET:HE3	2.35	0.47
1:D:37:LEU:HB2	1:D:71:THR:HG22	1.97	0.47
1:E:34:SER:HA	1:E:71:THR:OG1	2.14	0.47
2:I:6:THR:C	2:I:8:SER:H	2.22	0.47
2:I:49:ASN:OD1	2:I:49:ASN:C	2.56	0.47
2:I:73:ASP:OD2	2:I:73:ASP:C	2.57	0.47
3:K:20:MET:HB2	3:K:57:THR:HG21	1.96	0.47
3:K:51:VAL:HG13	3:K:54:VAL:CG2	2.44	0.47
3:L:51:VAL:O	3:L:52:GLN:C	2.57	0.47
2:I:18:LEU:C	2:I:18:LEU:CD2	2.73	0.47
3:L:16:LEU:HD12	3:L:90:ILE:HG23	1.95	0.47
3:N:57:THR:O	3:N:61:LEU:HD22	2.13	0.47
1:C:47:TRP:HB3	1:C:71:THR:CG2	2.39	0.47
1:E:25:LEU:HD13	1:E:86:VAL:HG22	1.94	0.47
1:F:119:ALA:C	1:F:121:LEU:N	2.72	0.47
2:I:25:PHE:CE1	2:J:51:PHE:CZ	2.99	0.47
2:I:49:ASN:O	2:I:51:PHE:N	2.48	0.47
2:J:19:ILE:HA	2:J:22:LEU:HD22	1.96	0.47
3:L:11:TRP:HA	3:L:14:ASP:HB2	1.96	0.47
3:L:48:MET:O	3:L:51:VAL:HG12	2.15	0.47
3:L:64:TRP:HZ3	3:L:74:LEU:CD2	2.21	0.47
3:M:28:TRP:HZ3	3:M:46:LYS:HD2	1.78	0.47
3:N:29:MET:HE2	3:N:29:MET:HB2	1.73	0.47
1:C:35:LEU:O	1:D:59:LEU:HD22	2.13	0.47
1:C:83:GLY:O	1:C:85:SER:N	2.48	0.47
1:F:22:LEU:C	1:F:24:ALA:N	2.68	0.47
2:H:29:GLN:HE21	3:L:59:GLU:HG3	1.80	0.47
1:C:105:LEU:HD12	1:C:109:ILE:HD11	1.97	0.47
2:G:9:THR:HG22	2:G:10:TYR:N	2.30	0.47
2:J:65:SER:OG	2:J:67:THR:HB	2.15	0.47
1:C:40:ARG:NH2	1:C:65:GLU:OE2	2.44	0.47
1:F:101:VAL:CG1	1:F:102:LEU:N	2.78	0.47
2:G:34:LYS:HZ3	3:K:63:TRP:HZ3	1.61	0.47
3:L:75:VAL:HG12	3:L:76:ASP:N	2.30	0.47
3:N:35:VAL:C	3:N:36:ILE:HG12	2.40	0.47
3:N:45:ILE:HD11	3:N:59:GLU:C	2.40	0.47
1:A:47:TRP:CZ3	1:A:61:ILE:HG23	2.50	0.46
1:C:29:VAL:CG1	1:C:105:LEU:HD11	2.39	0.46
1:D:78:TRP:HA	1:D:81:ARG:HD2	1.97	0.46
1:F:62:ARG:O	1:F:62:ARG:HG3	2.13	0.46
2:H:88:LEU:HD13	2:H:96:PRO:HB2	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:22:LEU:O	1:E:24:ALA:N	2.48	0.46
1:E:42:GLN:HG3	2:I:56:PHE:CE2	2.50	0.46
2:G:19:ILE:C	2:G:21:LYS:N	2.72	0.46
3:K:51:VAL:O	3:K:52:GLN:C	2.58	0.46
3:L:13:LEU:HD22	3:L:13:LEU:HA	1.70	0.46
1:D:34:SER:HA	1:D:71:THR:OG1	2.15	0.46
1:D:39:VAL:HG21	2:H:72:PHE:CZ	2.50	0.46
1:E:22:LEU:HA	1:E:22:LEU:HD12	1.51	0.46
1:F:42:GLN:O	2:J:55:ARG:NH1	2.40	0.46
2:I:51:PHE:O	2:I:54:ARG:HB2	2.15	0.46
2:J:85:VAL:O	2:J:89:ILE:HG12	2.15	0.46
3:M:70:THR:CG2	3:M:72:GLN:HE21	2.25	0.46
1:A:106:GLY:O	1:A:110:GLU:HG2	2.16	0.46
2:G:93:PHE:HB3	2:G:96:PRO:HG3	1.97	0.46
2:I:75:GLY:C	2:I:77:THR:H	2.23	0.46
3:L:5:ILE:H	3:L:5:ILE:HG12	1.50	0.46
3:L:49:GLU:HA	3:L:56:ILE:HD13	1.96	0.46
3:N:14:ASP:O	3:N:18:ARG:HG3	2.16	0.46
1:C:85:SER:OG	1:C:86:VAL:N	2.49	0.46
1:C:88:ARG:NH2	1:C:92:LEU:HD21	2.31	0.46
1:F:30:ARG:CG	1:F:30:ARG:NH1	2.62	0.46
2:I:18:LEU:O	2:I:21:LYS:HB3	2.16	0.46
1:D:119:ALA:C	1:D:121:LEU:N	2.73	0.46
1:E:105:LEU:HD23	1:E:105:LEU:N	2.30	0.46
3:K:63:TRP:CD1	3:K:67:ARG:CZ	2.99	0.46
3:L:85:ARG:HA	3:L:88:GLN:CD	2.41	0.46
3:M:46:LYS:HA	3:M:56:ILE:HD11	1.97	0.46
3:M:70:THR:HG21	3:M:72:GLN:HG2	1.98	0.46
3:N:4:TYR:CE2	3:N:68:GLN:HG3	2.49	0.46
3:N:90:ILE:C	3:N:91:LEU:HD22	2.40	0.46
1:C:56:PHE:CE1	1:C:74:LEU:CD1	2.98	0.46
1:D:122:GLU:OE1	1:D:122:GLU:N	2.49	0.46
2:G:19:ILE:HD11	2:G:71:LEU:CB	2.45	0.46
2:G:38:ALA:HB1	2:G:87:LEU:HD11	1.97	0.46
2:I:23:SER:O	2:I:27:ASP:HB2	2.16	0.46
3:N:70:THR:H	3:N:73:GLN:NE2	2.14	0.46
1:F:38:ASN:ND2	1:F:71:THR:OG1	2.49	0.46
2:H:105:VAL:HG12	2:H:105:VAL:O	2.15	0.46
2:J:28:PRO:O	2:J:29:GLN:HB2	2.15	0.46
3:K:16:LEU:HD21	3:K:61:LEU:CD2	2.46	0.46
3:K:70:THR:OG1	3:K:73:GLN:HG3	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:M:89:ILE:HG12	3:M:90:ILE:CG2	2.46	0.46
1:A:114:GLN:HG2	1:A:118:ALA:CB	2.46	0.46
2:J:91:ASN:O	2:J:92:GLU:HB2	2.15	0.46
3:K:86:ALA:O	3:K:89:ILE:HG12	2.15	0.46
3:L:24:SER:C	3:L:26:TRP:N	2.74	0.46
1:F:29:VAL:O	1:F:30:ARG:C	2.58	0.46
1:F:85:SER:OG	1:F:86:VAL:N	2.47	0.46
3:M:31:PHE:HA	3:M:81:LEU:CD1	2.45	0.46
1:A:88:ARG:HG3	1:A:92:LEU:HG	1.99	0.45
1:A:99:ASP:O	1:A:100:ASP:C	2.58	0.45
1:F:81:ARG:HE	1:F:81:ARG:HB3	1.54	0.45
1:F:89:LEU:HD12	1:F:89:LEU:C	2.41	0.45
1:F:103:LEU:C	1:F:103:LEU:HD23	2.41	0.45
2:G:95:ALA:N	2:G:96:PRO:CD	2.79	0.45
2:H:40:LYS:HD3	2:H:44:GLY:O	2.16	0.45
2:I:90:GLN:HE22	2:I:91:ASN:ND2	2.14	0.45
3:M:3:CYS:C	3:M:71:VAL:HG12	2.41	0.45
1:A:58:TYR:CE1	1:D:30:ARG:CZ	2.99	0.45
1:C:88:ARG:HH21	1:C:92:LEU:HD21	1.80	0.45
1:D:19:MET:HG3	1:D:20:LEU:HD12	1.98	0.45
1:D:70:PRO:O	1:D:71:THR:C	2.59	0.45
1:E:33:LEU:HD22	1:E:105:LEU:HD12	1.98	0.45
1:F:19:MET:HB2	1:F:20:LEU:HD12	1.98	0.45
2:G:16:VAL:O	2:G:16:VAL:CG2	2.61	0.45
2:H:93:PHE:CD2	3:L:66:MET:SD	3.06	0.45
2:I:101:LEU:C	2:I:103:ASP:H	2.25	0.45
2:J:18:LEU:O	2:J:21:LYS:HB3	2.15	0.45
2:J:49:ASN:OD1	2:J:51:PHE:HB2	2.16	0.45
3:M:36:ILE:HG22	3:M:36:ILE:O	2.16	0.45
1:A:78:TRP:O	1:A:80:GLY:N	2.50	0.45
1:E:51:ALA:O	1:E:56:PHE:HB2	2.16	0.45
2:J:70:LEU:HD23	2:J:70:LEU:C	2.41	0.45
3:L:91:LEU:HA	3:L:91:LEU:HD13	1.47	0.45
3:N:21:ASP:OD1	3:N:57:THR:HG22	2.16	0.45
3:N:91:LEU:C	3:N:93:TRP:H	2.23	0.45
1:A:26:ASN:HD22	1:A:27:MET:N	2.14	0.45
1:A:64:LEU:O	1:A:66:THR:N	2.49	0.45
2:H:21:LYS:CD	2:H:99:LEU:HD21	2.46	0.45
2:I:73:ASP:OD2	2:I:73:ASP:O	2.34	0.45
2:J:19:ILE:C	2:J:21:LYS:N	2.74	0.45
3:L:56:ILE:HA	3:L:56:ILE:HD12	1.76	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:M:35:VAL:O	3:M:35:VAL:HG22	2.16	0.45
3:N:16:LEU:HG	3:N:20:MET:HE1	1.99	0.45
1:D:81:ARG:HB3	1:D:82:PRO:CD	2.46	0.45
1:E:43:VAL:CG1	2:I:52:HIS:CD2	2.99	0.45
2:H:4:PRO:O	2:H:6:THR:HG22	2.17	0.45
2:H:41:LYS:HB3	2:H:42:PRO:CD	2.41	0.45
2:H:76:THR:O	2:H:76:THR:OG1	2.33	0.45
3:M:16:LEU:CD2	3:M:61:LEU:HD11	2.47	0.45
1:A:39:VAL:CG2	1:B:59:LEU:HD23	2.46	0.45
1:C:26:ASN:O	1:C:28:ARG:N	2.50	0.45
1:C:62:ARG:HD2	1:C:65:GLU:OE1	2.17	0.45
1:D:83:GLY:O	1:D:84:ALA:C	2.59	0.45
2:G:70:LEU:HD23	2:G:70:LEU:C	2.42	0.45
2:H:19:ILE:C	2:H:21:LYS:N	2.75	0.45
2:I:50:GLN:HA	3:L:10:SER:OG	2.16	0.45
3:L:11:TRP:CG	3:L:12:VAL:N	2.85	0.45
3:L:20:MET:HE3	3:L:20:MET:N	2.30	0.45
3:L:52:GLN:O	3:L:53:GLY:C	2.60	0.45
3:M:72:GLN:O	3:M:76:ASP:CG	2.60	0.45
1:B:63:GLN:O	1:B:66:THR:HB	2.17	0.45
1:C:82:PRO:HB2	1:C:83:GLY:H	1.52	0.45
2:H:12:ARG:C	2:H:12:ARG:HD2	2.41	0.45
2:I:67:THR:O	2:I:71:LEU:HB2	2.16	0.45
2:J:23:SER:O	2:J:27:ASP:HB2	2.16	0.45
3:K:8:LEU:HD13	3:K:8:LEU:HA	1.74	0.45
3:K:19:ASN:HD22	3:K:19:ASN:HA	1.51	0.45
1:C:22:LEU:C	1:C:24:ALA:N	2.73	0.45
1:E:103:LEU:HD23	1:E:104:GLU:HG2	1.98	0.45
1:F:22:LEU:C	1:F:24:ALA:H	2.23	0.45
2:H:25:PHE:CE1	2:H:95:ALA:HB1	2.51	0.45
2:H:29:GLN:NE2	3:L:59:GLU:HG3	2.31	0.45
3:K:14:ASP:O	3:K:18:ARG:HD3	2.16	0.45
3:K:63:TRP:O	3:K:67:ARG:CD	2.65	0.45
3:K:91:LEU:CD2	3:K:94:LYS:HG2	2.47	0.45
3:M:31:PHE:HA	3:M:81:LEU:HD11	1.99	0.45
3:N:63:TRP:NE1	3:N:67:ARG:NH2	2.65	0.45
1:C:56:PHE:HD2	1:C:60:GLU:OE2	2.00	0.45
1:D:34:SER:O	1:D:35:LEU:C	2.59	0.45
1:D:73:ARG:NE	1:D:73:ARG:CA	2.79	0.45
1:F:96:LEU:HD23	1:F:96:LEU:N	2.32	0.45
2:I:14:LEU:HD22	2:I:18:LEU:HD22	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:122:GLU:CG	1:A:123:HIS:H	2.22	0.45
1:E:107:PRO:CG	1:E:108:SER:N	2.78	0.45
1:F:26:ASN:O	1:F:28:ARG:N	2.50	0.45
2:G:62:THR:OG1	2:G:64:LYS:HG3	2.17	0.45
3:L:35:VAL:HG12	3:L:37:THR:HG1	1.80	0.45
3:M:54:VAL:O	3:M:54:VAL:CG2	2.64	0.45
1:A:54:MET:CE	1:A:74:LEU:HG	2.47	0.44
1:E:119:ALA:O	1:E:120:ALA:C	2.60	0.44
2:I:60:LEU:N	2:I:60:LEU:HD22	2.31	0.44
2:I:88:LEU:HA	2:I:88:LEU:HD23	1.71	0.44
2:J:52:HIS:ND1	2:J:52:HIS:N	2.65	0.44
2:J:82:GLY:O	2:J:84:LEU:N	2.50	0.44
1:A:37:LEU:HD12	1:A:71:THR:CG2	2.47	0.44
1:E:96:LEU:HA	1:E:96:LEU:HD23	1.58	0.44
2:H:81:VAL:O	2:H:85:VAL:HG22	2.17	0.44
2:I:50:GLN:NE2	3:L:6:TYR:HA	2.33	0.44
2:J:100:LEU:HB3	2:J:101:LEU:CD2	2.37	0.44
2:J:107:LEU:O	2:J:108:GLU:O	2.35	0.44
3:K:35:VAL:HG12	3:K:37:THR:OG1	2.18	0.44
3:N:5:ILE:HG21	3:N:61:LEU:O	2.17	0.44
1:E:22:LEU:O	1:E:25:LEU:CD2	2.65	0.44
1:F:70:PRO:HA	1:F:73:ARG:HB2	1.99	0.44
2:I:70:LEU:C	2:I:70:LEU:HD23	2.43	0.44
2:I:91:ASN:O	2:I:92:GLU:HB2	2.16	0.44
2:J:75:GLY:C	2:J:77:THR:H	2.25	0.44
1:B:88:ARG:HA	1:B:88:ARG:HD3	1.58	0.44
1:D:88:ARG:HA	1:D:91:GLU:OE1	2.17	0.44
1:F:33:LEU:HD12	1:F:33:LEU:HA	1.80	0.44
2:H:39:ILE:O	2:H:47:ARG:HB3	2.18	0.44
2:H:54:ARG:HH21	3:K:58:ARG:NE	2.16	0.44
3:M:32:ALA:O	3:M:35:VAL:HG12	2.16	0.44
3:M:48:MET:HE1	3:M:59:GLU:HB3	1.98	0.44
3:M:91:LEU:C	3:M:93:TRP:H	2.25	0.44
3:N:28:TRP:HZ2	3:N:56:ILE:CG1	2.20	0.44
1:D:71:THR:O	1:D:74:LEU:HB3	2.17	0.44
2:G:11:VAL:O	2:G:14:LEU:HD12	2.18	0.44
2:J:2:MET:HG2	2:J:3:GLY:N	2.32	0.44
2:J:73:ASP:C	2:J:73:ASP:OD2	2.59	0.44
3:K:4:TYR:CE2	3:K:68:GLN:HG3	2.52	0.44
3:L:52:GLN:H	3:L:52:GLN:HG2	1.54	0.44
3:L:63:TRP:O	3:L:67:ARG:NE	2.51	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:N:92:ASN:N	3:N:92:ASN:HD22	2.14	0.44
1:E:39:VAL:HG23	1:E:39:VAL:O	2.17	0.44
3:M:79:CYS:C	3:M:81:LEU:N	2.76	0.44
1:B:90:LEU:HD12	1:B:90:LEU:HA	1.36	0.44
2:H:84:LEU:HD23	2:H:100:LEU:CD1	2.44	0.44
2:J:98:SER:HA	2:J:105:VAL:HG11	1.99	0.44
3:K:85:ARG:C	3:K:88:GLN:HG3	2.42	0.44
3:L:8:LEU:HA	3:L:9:PRO:HD2	1.73	0.44
1:B:66:THR:CG2	1:B:67:GLN:N	2.78	0.44
1:C:43:VAL:HG23	2:G:52:HIS:NE2	2.32	0.44
1:E:58:TYR:HA	2:H:16:VAL:HG12	2.00	0.44
1:F:102:LEU:H	1:F:102:LEU:CD1	2.22	0.44
2:J:30:GLU:OE2	2:J:30:GLU:N	2.51	0.44
3:M:19:ASN:O	3:M:21:ASP:N	2.51	0.44
3:M:52:GLN:O	3:M:53:GLY:C	2.60	0.44
1:A:51:ALA:HA	1:A:74:LEU:HD11	1.99	0.44
1:A:52:GLU:CD	1:D:28:ARG:HB3	2.43	0.44
1:B:33:LEU:O	1:B:33:LEU:HD22	2.17	0.44
1:B:76:ASP:C	1:B:78:TRP:N	2.76	0.44
1:C:103:LEU:C	1:C:103:LEU:CD2	2.86	0.44
1:E:20:LEU:HD23	1:E:117:ILE:HD13	2.00	0.44
1:E:59:LEU:HD23	1:E:59:LEU:C	2.43	0.44
1:F:69:ASP:N	1:F:70:PRO:CD	2.80	0.44
2:I:82:GLY:O	2:I:84:LEU:N	2.51	0.44
2:J:25:PHE:HD2	2:J:96:PRO:CA	2.31	0.44
3:K:43:ARG:HD3	3:K:43:ARG:HA	1.89	0.44
3:L:37:THR:HG22	3:L:38:ASP:H	1.83	0.44
3:M:63:TRP:CE2	3:M:67:ARG:NH2	2.85	0.44
3:M:75:VAL:CG1	3:M:90:ILE:HD11	2.48	0.44
3:N:24:SER:O	3:N:26:TRP:N	2.50	0.44
3:N:90:ILE:O	3:N:91:LEU:HD22	2.17	0.44
1:B:37:LEU:HD23	1:B:37:LEU:HA	1.66	0.43
1:F:83:GLY:O	1:F:84:ALA:C	2.61	0.43
3:K:44:LYS:HE3	3:K:48:MET:CE	2.47	0.43
3:K:91:LEU:C	3:K:93:TRP:H	2.25	0.43
3:N:64:TRP:HA	3:N:67:ARG:NH2	2.32	0.43
1:C:29:VAL:O	1:C:30:ARG:C	2.60	0.43
1:D:121:LEU:O	1:D:122:GLU:HG3	2.17	0.43
2:G:2:MET:HB3	2:G:3:GLY:H	1.71	0.43
2:G:93:PHE:O	2:G:96:PRO:HG2	2.17	0.43
2:H:71:LEU:HD23	2:H:71:LEU:HA	1.31	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:95:ALA:CB	3:L:68:GLN:HE22	2.31	0.43
2:I:49:ASN:O	2:I:50:GLN:C	2.61	0.43
2:J:49:ASN:O	2:J:51:PHE:N	2.52	0.43
3:K:19:ASN:O	3:K:21:ASP:N	2.52	0.43
3:M:13:LEU:HD23	3:M:13:LEU:HA	1.84	0.43
1:B:98:ARG:NH1	1:B:100:ASP:OD1	2.51	0.43
1:C:115:LYS:O	1:C:119:ALA:HB3	2.18	0.43
1:E:113:CYS:C	1:E:115:LYS:N	2.74	0.43
2:I:50:GLN:HG2	3:L:58:ARG:HH12	1.83	0.43
2:J:39:ILE:HD12	2:J:74:TRP:CD2	2.53	0.43
3:K:19:ASN:O	3:K:20:MET:C	2.60	0.43
3:K:44:LYS:O	3:K:46:LYS:N	2.52	0.43
3:M:5:ILE:HD12	3:M:64:TRP:HB3	1.99	0.43
3:M:57:THR:O	3:M:61:LEU:HB2	2.17	0.43
1:B:92:LEU:HD23	1:B:92:LEU:HA	1.74	0.43
2:I:73:ASP:O	2:I:76:THR:HG23	2.18	0.43
2:I:84:LEU:HD23	2:I:100:LEU:HD11	2.00	0.43
2:J:6:THR:C	2:J:8:SER:H	2.26	0.43
2:J:101:LEU:C	2:J:103:ASP:H	2.26	0.43
3:M:24:SER:C	3:M:26:TRP:N	2.75	0.43
3:N:19:ASN:O	3:N:20:MET:C	2.61	0.43
3:N:52:GLN:O	3:N:53:GLY:C	2.60	0.43
1:B:117:ILE:CG2	1:B:118:ALA:N	2.81	0.43
1:C:69:ASP:N	1:C:70:PRO:CD	2.80	0.43
1:C:71:THR:O	1:C:74:LEU:HB3	2.19	0.43
1:F:122:GLU:HB2	1:F:123:HIS:CE1	2.52	0.43
2:H:48:TYR:HA	2:H:52:HIS:ND1	2.33	0.43
2:I:26:ILE:H	2:I:26:ILE:HG12	1.46	0.43
3:N:19:ASN:HB2	3:N:20:MET:HE2	2.01	0.43
1:B:93:LEU:HD23	1:B:93:LEU:HA	1.73	0.43
1:C:85:SER:HB3	1:C:88:ARG:CB	2.33	0.43
1:F:29:VAL:HG23	1:F:30:ARG:N	2.32	0.43
2:H:26:ILE:CG2	2:H:32:TRP:HB3	2.47	0.43
2:I:71:LEU:HD23	2:I:71:LEU:HA	1.36	0.43
1:A:89:LEU:HD11	1:A:93:LEU:HD21	2.01	0.43
1:B:38:ASN:C	1:B:38:ASN:HD22	2.26	0.43
1:E:80:GLY:C	1:E:81:ARG:HG2	2.42	0.43
1:E:105:LEU:C	1:E:107:PRO:HD2	2.43	0.43
2:G:85:VAL:HG11	2:G:101:LEU:CD2	2.48	0.43
2:H:11:VAL:HA	2:H:14:LEU:HD12	2.01	0.43
2:H:60:LEU:HD13	2:H:60:LEU:HA	1.40	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:I:5:ILE:HD12	2:I:5:ILE:H	1.84	0.43
2:I:19:ILE:C	2:I:21:LYS:N	2.75	0.43
2:I:22:LEU:HD12	2:I:22:LEU:O	2.17	0.43
3:K:66:MET:HG3	3:K:66:MET:O	2.19	0.43
3:N:44:LYS:HG2	3:N:48:MET:HE3	2.01	0.43
1:B:51:ALA:HB1	1:B:61:ILE:HD13	2.01	0.43
1:B:58:TYR:CE2	1:B:62:ARG:HG2	2.53	0.43
1:C:70:PRO:O	1:C:74:LEU:N	2.51	0.43
1:C:98:ARG:HD3	1:C:98:ARG:HA	1.53	0.43
1:F:34:SER:C	1:F:36:PHE:H	2.27	0.43
1:F:110:GLU:O	1:F:114:GLN:NE2	2.52	0.43
3:K:55:SER:O	3:K:58:ARG:HB2	2.19	0.43
3:M:15:ASP:HA	3:M:18:ARG:HG2	2.00	0.43
3:N:16:LEU:HD22	3:N:61:LEU:HD21	2.01	0.43
3:N:25:GLU:O	3:N:29:MET:HB2	2.18	0.43
2:I:2:MET:CG	2:I:3:GLY:N	2.82	0.43
1:C:58:TYR:CE1	1:F:30:ARG:CZ	3.02	0.43
1:E:71:THR:O	1:E:75:LEU:HG	2.19	0.43
2:J:101:LEU:HA	2:J:102:PRO:HD2	1.75	0.43
3:K:24:SER:C	3:K:26:TRP:N	2.77	0.43
3:K:93:TRP:CD1	3:K:93:TRP:C	2.96	0.43
3:N:24:SER:C	3:N:26:TRP:N	2.77	0.43
1:B:26:ASN:HB2	1:B:27:MET:SD	2.59	0.42
1:B:69:ASP:H	1:B:70:PRO:HD3	1.79	0.42
1:B:113:CYS:O	1:B:117:ILE:HB	2.19	0.42
1:F:79:GLN:C	1:F:81:ARG:N	2.72	0.42
2:H:68:SER:HA	2:H:71:LEU:HB2	2.01	0.42
3:M:70:THR:CG2	3:M:72:GLN:HG2	2.49	0.42
1:A:43:VAL:HG21	1:E:64:LEU:HD21	1.98	0.42
1:A:50:LEU:HD23	1:A:74:LEU:HD21	2.00	0.42
1:B:52:GLU:HG3	1:E:26:ASN:HD21	1.84	0.42
1:B:105:LEU:C	1:B:107:PRO:HD2	2.44	0.42
2:H:35:LEU:HD21	2:H:39:ILE:HD11	2.01	0.42
2:J:73:ASP:O	2:J:76:THR:HG23	2.18	0.42
3:M:16:LEU:HD12	3:M:90:ILE:HG21	2.01	0.42
1:B:66:THR:O	1:B:67:GLN:C	2.61	0.42
1:F:25:LEU:CD2	1:F:86:VAL:CG2	2.93	0.42
1:F:29:VAL:CG2	1:F:30:ARG:N	2.81	0.42
3:K:89:ILE:HA	3:K:92:ASN:OD1	2.18	0.42
3:L:4:TYR:HB2	3:L:7:GLN:CD	2.45	0.42
3:N:77:LEU:HD12	3:N:77:LEU:N	2.34	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:58:TYR:CE2	1:D:30:ARG:HD2	2.55	0.42
1:B:63:GLN:O	1:B:64:LEU:C	2.63	0.42
1:D:78:TRP:CZ2	1:D:88:ARG:HG2	2.53	0.42
2:I:11:VAL:C	2:I:13:CYS:H	2.27	0.42
2:J:11:VAL:O	2:J:13:CYS:N	2.53	0.42
2:J:49:ASN:O	2:J:50:GLN:C	2.62	0.42
2:J:54:ARG:NH2	3:M:62:TRP:CG	2.87	0.42
3:M:19:ASN:O	3:M:20:MET:C	2.61	0.42
3:M:70:THR:HG21	3:M:72:GLN:CG	2.50	0.42
1:C:22:LEU:HD22	1:C:22:LEU:N	2.05	0.42
2:H:3:GLY:H	2:H:4:PRO:HD2	1.85	0.42
2:J:11:VAL:C	2:J:13:CYS:N	2.77	0.42
2:J:41:LYS:HB3	2:J:42:PRO:CD	2.49	0.42
3:K:5:ILE:HD11	3:K:69:ALA:HB1	2.02	0.42
3:N:70:THR:HG23	3:N:73:GLN:N	2.33	0.42
3:N:75:VAL:HG13	3:N:91:LEU:HD21	2.01	0.42
1:D:109:ILE:O	1:D:110:GLU:C	2.61	0.42
1:F:53:GLU:C	1:F:55:ASP:N	2.78	0.42
2:J:12:ARG:H	2:J:12:ARG:HG3	1.59	0.42
3:K:35:VAL:HG11	3:K:60:LEU:HD11	2.02	0.42
3:M:32:ALA:HA	3:M:60:LEU:HD21	2.02	0.42
1:B:46:ASP:OD2	1:B:46:ASP:N	2.52	0.42
1:C:111:GLU:C	1:C:113:CYS:H	2.28	0.42
1:D:98:ARG:HD2	1:D:100:ASP:OD2	2.19	0.42
1:E:62:ARG:HD2	1:E:65:GLU:OE1	2.19	0.42
1:E:122:GLU:OE2	1:E:123:HIS:CE1	2.73	0.42
2:H:25:PHE:O	2:H:28:PRO:HD2	2.19	0.42
2:H:89:ILE:CD1	2:H:90:GLN:N	2.77	0.42
3:M:30:GLU:O	3:M:34:TYR:HB2	2.20	0.42
1:B:62:ARG:NH1	1:E:76:ASP:CG	2.78	0.42
1:B:117:ILE:HG23	1:B:118:ALA:N	2.33	0.42
2:G:10:TYR:N	2:G:10:TYR:CD2	2.88	0.42
3:N:79:CYS:C	3:N:81:LEU:N	2.77	0.42
1:A:89:LEU:CD1	1:A:93:LEU:HD21	2.50	0.42
1:C:34:SER:C	1:C:36:PHE:H	2.28	0.42
1:D:78:TRP:O	1:D:78:TRP:CG	2.73	0.42
2:I:50:GLN:NE2	3:L:58:ARG:NH2	2.54	0.42
3:N:2:ALA:HB2	3:N:93:TRP:NE1	2.35	0.42
3:N:45:ILE:CD1	3:N:60:LEU:HA	2.49	0.42
3:N:83:LEU:HD23	3:N:83:LEU:HA	1.85	0.42
1:A:78:TRP:O	1:A:79:GLN:C	2.63	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:62:ARG:HH12	1:E:76:ASP:CG	2.28	0.42
1:B:69:ASP:N	1:B:70:PRO:CD	2.78	0.42
1:D:89:LEU:O	1:D:93:LEU:HG	2.20	0.42
1:E:31:ARG:HH11	1:E:31:ARG:HG2	1.82	0.42
1:E:33:LEU:O	1:E:33:LEU:HD12	2.20	0.42
1:E:78:TRP:O	1:E:78:TRP:CG	2.72	0.42
2:G:71:LEU:HD13	2:G:71:LEU:HA	1.70	0.42
2:H:11:VAL:C	2:H:13:CYS:H	2.28	0.42
3:K:44:LYS:HE3	3:K:48:MET:HE1	2.01	0.42
3:L:87:ALA:HA	3:L:90:ILE:HG12	2.02	0.42
3:M:16:LEU:CD1	3:M:90:ILE:HG21	2.50	0.42
1:A:66:THR:O	1:A:67:GLN:C	2.63	0.41
2:G:6:THR:C	2:G:8:SER:H	2.28	0.41
2:G:87:LEU:HD12	2:G:90:GLN:NE2	2.35	0.41
2:I:50:GLN:HE22	3:L:6:TYR:HA	1.84	0.41
3:K:75:VAL:HG22	3:K:90:ILE:HD11	2.02	0.41
3:L:32:ALA:C	3:L:34:TYR:N	2.76	0.41
3:M:78:LEU:HD13	3:M:86:ALA:HB3	2.01	0.41
1:A:48:THR:HG22	1:A:61:ILE:HG21	2.01	0.41
1:C:79:GLN:C	1:C:81:ARG:N	2.76	0.41
1:D:98:ARG:O	1:D:100:ASP:N	2.53	0.41
2:G:49:ASN:O	2:G:50:GLN:C	2.63	0.41
2:H:35:LEU:CD2	2:H:70:LEU:HD21	2.50	0.41
3:L:3:CYS:HB3	3:L:7:GLN:NE2	2.35	0.41
3:M:90:ILE:O	3:M:90:ILE:HG13	2.19	0.41
3:N:32:ALA:O	3:N:35:VAL:HG12	2.20	0.41
1:C:100:ASP:OD2	1:C:101:VAL:N	2.54	0.41
1:D:59:LEU:HA	1:D:59:LEU:HD23	1.82	0.41
2:G:82:GLY:O	2:G:84:LEU:N	2.53	0.41
2:H:107:LEU:HD23	2:H:107:LEU:HA	1.93	0.41
3:K:24:SER:O	3:K:26:TRP:N	2.52	0.41
3:M:45:ILE:CD1	3:M:60:LEU:HB2	2.50	0.41
1:B:34:SER:HA	1:B:71:THR:HG21	2.02	0.41
1:B:98:ARG:O	1:B:101:VAL:HG22	2.21	0.41
1:E:92:LEU:HA	1:E:92:LEU:HD23	1.48	0.41
1:F:92:LEU:HA	1:F:92:LEU:HD23	1.66	0.41
2:H:59:LEU:HD13	2:H:69:GLU:HG3	2.01	0.41
2:H:76:THR:H	2:H:76:THR:HG23	1.40	0.41
2:H:82:GLY:C	2:H:84:LEU:N	2.78	0.41
3:L:91:LEU:HD22	3:L:91:LEU:N	2.36	0.41
3:N:23:LEU:HD11	3:N:83:LEU:HD11	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:33:LEU:HA	1:B:33:LEU:HD23	1.45	0.41
1:B:44:ALA:HB2	1:F:77:ALA:O	2.21	0.41
1:B:100:ASP:HB2	1:F:80:GLY:HA2	2.02	0.41
1:B:106:GLY:N	1:B:107:PRO:CD	2.82	0.41
1:C:117:ILE:CG2	1:C:118:ALA:N	2.83	0.41
2:H:86:ASP:O	2:H:90:GLN:HB3	2.20	0.41
2:I:18:LEU:HD23	2:I:18:LEU:O	2.20	0.41
2:I:56:PHE:O	2:I:59:LEU:HG	2.20	0.41
2:I:85:VAL:CG2	2:I:86:ASP:N	2.83	0.41
2:I:93:PHE:HD2	3:M:66:MET:SD	2.42	0.41
3:N:5:ILE:HD11	3:N:64:TRP:HE3	1.85	0.41
1:A:26:ASN:ND2	1:A:28:ARG:HB3	2.35	0.41
1:D:122:GLU:HB2	1:D:123:HIS:CE1	2.55	0.41
2:I:93:PHE:O	2:I:96:PRO:HD2	2.21	0.41
1:A:50:LEU:HD23	1:A:74:LEU:CD2	2.51	0.41
1:B:110:GLU:OE1	1:B:110:GLU:HA	2.20	0.41
1:D:56:PHE:CE1	1:D:74:LEU:HD11	2.54	0.41
2:H:49:ASN:O	2:H:51:PHE:N	2.54	0.41
3:K:25:GLU:CD	3:N:53:GLY:O	2.63	0.41
3:L:12:VAL:HG12	3:L:13:LEU:N	2.34	0.41
3:L:24:SER:C	3:L:26:TRP:H	2.29	0.41
3:N:15:ASP:CA	3:N:18:ARG:HG3	2.50	0.41
1:A:114:GLN:CG	1:A:118:ALA:HB2	2.50	0.41
1:A:123:HIS:O	1:A:123:HIS:CG	2.73	0.41
1:B:48:THR:OG1	1:B:49:ALA:N	2.53	0.41
1:B:78:TRP:O	1:B:80:GLY:N	2.54	0.41
1:D:28:ARG:HE	1:D:28:ARG:HB2	1.49	0.41
1:F:58:TYR:HD2	2:I:16:VAL:HG22	1.84	0.41
2:G:35:LEU:HD23	2:G:70:LEU:HD21	2.03	0.41
2:G:87:LEU:O	2:G:91:ASN:HB2	2.20	0.41
2:H:79:CYS:O	2:H:79:CYS:SG	2.79	0.41
1:C:98:ARG:HD2	1:C:100:ASP:OD2	2.20	0.41
1:D:33:LEU:HA	1:D:33:LEU:HD12	1.68	0.41
1:F:36:PHE:O	1:F:39:VAL:CG2	2.57	0.41
2:G:95:ALA:HB3	3:K:66:MET:HE2	2.03	0.41
2:H:19:ILE:HG22	2:H:20:ARG:N	2.35	0.41
2:I:77:THR:O	2:I:79:CYS:N	2.47	0.41
2:J:18:LEU:CD1	2:J:18:LEU:O	2.68	0.41
2:J:76:THR:O	2:J:76:THR:OG1	2.33	0.41
3:K:65:GLY:C	3:K:67:ARG:H	2.29	0.41
3:L:61:LEU:HD12	3:L:61:LEU:HA	1.64	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:M:8:LEU:HD13	3:M:8:LEU:HA	1.87	0.41
3:N:8:LEU:HD22	3:N:12:VAL:HG11	2.02	0.41
3:N:54:VAL:CG1	3:N:59:GLU:OE2	2.66	0.41
1:D:50:LEU:HD21	1:D:54:MET:HE3	2.02	0.41
1:E:43:VAL:CG2	2:I:48:TYR:HE1	2.33	0.41
1:E:47:TRP:HB3	1:E:71:THR:HG22	2.03	0.41
1:F:89:LEU:CD1	1:F:93:LEU:HG	2.51	0.41
2:G:55:ARG:O	2:G:58:ALA:HB3	2.21	0.41
2:G:67:THR:O	2:G:71:LEU:HB2	2.20	0.41
2:J:47:ARG:HH21	2:J:74:TRP:HA	1.86	0.41
2:J:82:GLY:C	2:J:84:LEU:N	2.77	0.41
1:A:37:LEU:HD23	1:A:98:ARG:HD2	2.04	0.40
1:F:58:TYR:HA	2:I:16:VAL:CG2	2.51	0.40
2:H:80:THR:C	2:H:82:GLY:N	2.79	0.40
3:K:4:TYR:HB2	3:K:7:GLN:HG3	2.03	0.40
3:L:32:ALA:CB	3:L:42:LEU:HD11	2.50	0.40
1:B:88:ARG:O	1:B:92:LEU:HB2	2.21	0.40
1:C:19:MET:HB2	1:C:20:LEU:H	1.69	0.40
1:D:42:GLN:NE2	2:H:59:LEU:HD11	2.37	0.40
1:D:51:ALA:O	1:D:56:PHE:HD1	2.04	0.40
1:E:102:LEU:HD13	1:E:102:LEU:HA	1.88	0.40
2:H:60:LEU:O	2:I:62:THR:HG22	2.21	0.40
2:I:62:THR:OG1	2:I:64:LYS:HG3	2.21	0.40
2:I:96:PRO:HD3	3:M:66:MET:SD	2.60	0.40
2:I:99:LEU:O	2:I:99:LEU:HD12	2.22	0.40
2:J:82:GLY:O	2:J:83:ASP:C	2.65	0.40
2:J:90:GLN:O	2:J:90:GLN:HG2	2.21	0.40
3:K:64:TRP:HZ3	3:K:74:LEU:HD23	1.85	0.40
1:A:76:ASP:C	1:A:78:TRP:N	2.75	0.40
1:C:56:PHE:HE1	1:C:74:LEU:CD1	2.35	0.40
2:I:89:ILE:CG1	2:I:90:GLN:N	2.80	0.40
2:J:76:THR:HG23	2:J:76:THR:H	1.45	0.40
3:N:23:LEU:C	3:N:23:LEU:HD22	2.47	0.40
1:B:48:THR:HB	1:E:27:MET:HG2	2.03	0.40
1:B:73:ARG:CZ	1:B:73:ARG:CB	2.99	0.40
1:C:90:LEU:HD13	1:C:90:LEU:HA	1.55	0.40
2:G:85:VAL:HG23	2:G:97:ALA:CB	2.44	0.40
2:G:85:VAL:CG2	2:G:89:ILE:HG13	2.51	0.40
2:I:66:PRO:HA	2:I:69:GLU:HB2	2.04	0.40
3:K:63:TRP:NE1	3:K:67:ARG:HH11	2.18	0.40
3:K:78:LEU:HD13	3:K:87:ALA:N	2.37	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:5:ILE:HG12	3:L:69:ALA:O	2.22	0.40
3:L:20:MET:CE	3:L:20:MET:N	2.69	0.40
3:L:58:ARG:NH2	3:L:62:TRP:HB2	2.37	0.40
3:M:34:TYR:CD2	3:M:77:LEU:HD23	2.56	0.40
3:N:11:TRP:CG	3:N:12:VAL:N	2.89	0.40
3:N:23:LEU:HD23	3:N:85:ARG:NH1	2.37	0.40
1:A:69:ASP:O	1:A:69:ASP:CG	2.62	0.40
1:A:100:ASP:HB2	1:E:80:GLY:HA2	2.02	0.40
1:C:105:LEU:O	1:C:109:ILE:HG13	2.22	0.40
1:D:102:LEU:O	1:D:103:LEU:C	2.64	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all 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	103/110 (94%)	74 (72%)	21 (20%)	8 (8%)	1	4
1	B	103/110 (94%)	75 (73%)	22 (21%)	6 (6%)	1	8
1	C	103/110 (94%)	65 (63%)	30 (29%)	8 (8%)	1	4
1	D	103/110 (94%)	70 (68%)	18 (18%)	15 (15%)	0	0
1	E	103/110 (94%)	68 (66%)	22 (21%)	13 (13%)	0	1
1	F	103/110 (94%)	65 (63%)	30 (29%)	8 (8%)	1	4
2	G	105/113 (93%)	77 (73%)	19 (18%)	9 (9%)	0	4
2	H	105/113 (93%)	80 (76%)	17 (16%)	8 (8%)	1	5
2	I	105/113 (93%)	76 (72%)	20 (19%)	9 (9%)	0	4
2	J	105/113 (93%)	77 (73%)	17 (16%)	11 (10%)	0	2
3	K	91/111 (82%)	64 (70%)	13 (14%)	14 (15%)	0	0
3	L	91/111 (82%)	61 (67%)	16 (18%)	14 (15%)	0	0

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	M	91/111 (82%)	64 (70%)	16 (18%)	11 (12%)	0	1
3	N	91/111 (82%)	63 (69%)	18 (20%)	10 (11%)	0	2
All	All	1402/1556 (90%)	979 (70%)	279 (20%)	144 (10%)	0	2

All (144) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	27	MET
1	C	82	PRO
1	C	84	ALA
1	D	28	ARG
1	D	65	GLU
1	D	82	PRO
1	D	83	GLY
1	D	84	ALA
1	E	27	MET
1	E	28	ARG
1	E	65	GLU
1	E	82	PRO
1	E	83	GLY
1	E	84	ALA
1	F	27	MET
1	F	82	PRO
1	F	84	ALA
2	G	79	CYS
2	H	79	CYS
2	I	79	CYS
2	J	79	CYS
3	K	19	ASN
3	K	21	ASP
3	K	24	SER
3	L	19	ASN
3	L	20	MET
3	L	21	ASP
3	M	19	ASN
3	M	21	ASP
3	N	19	ASN
3	N	21	ASP
3	N	24	SER
3	N	52	GLN
1	A	26	ASN

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Mol	Chain	Res	Type
1	A	28	ARG
1	A	67	GLN
1	B	26	ASN
1	B	28	ARG
1	C	33	LEU
1	C	83	GLY
1	D	27	MET
1	D	35	LEU
1	D	63	GLN
1	D	118	ALA
1	D	120	ALA
1	E	35	LEU
1	E	63	GLN
1	E	120	ALA
1	F	35	LEU
1	F	83	GLY
2	G	83	ASP
2	H	83	ASP
2	I	50	GLN
2	I	83	ASP
2	J	83	ASP
3	K	3	CYS
3	K	20	MET
3	K	52	GLN
3	K	53	GLY
3	K	66	MET
3	L	3	CYS
3	L	24	SER
3	L	25	GLU
3	L	36	ILE
3	L	52	GLN
3	L	53	GLY
3	L	66	MET
3	M	3	CYS
3	M	20	MET
3	M	24	SER
3	M	25	GLU
3	M	52	GLN
3	M	53	GLY
3	N	3	CYS
3	N	36	ILE
3	N	53	GLY

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Mol	Chain	Res	Type
1	A	65	GLU
1	A	79	GLN
1	A	82	PRO
1	B	65	GLU
1	D	51	ALA
1	D	112	ASP
1	D	119	ALA
1	E	114	GLN
1	E	118	ALA
1	E	119	ALA
2	G	20	ARG
2	G	100	LEU
2	I	106	PRO
2	J	50	GLN
3	K	36	ILE
3	K	80	ARG
3	L	80	ARG
3	M	36	ILE
3	M	80	ARG
3	N	85	ARG
1	B	67	GLN
1	B	82	PRO
1	C	35	LEU
1	E	102	LEU
1	F	33	LEU
2	G	103	ASP
2	H	20	ARG
2	H	50	GLN
2	H	103	ASP
2	I	20	ARG
2	I	76	THR
2	J	76	THR
2	J	100	LEU
3	K	25	GLU
3	K	85	ARG
3	L	55	SER
3	N	20	MET
3	N	80	ARG
1	A	104	GLU
1	C	103	LEU
1	D	114	GLN
1	F	32	ARG

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Mol	Chain	Res	Type
1	F	103	LEU
2	G	50	GLN
2	J	12	ARG
2	J	102	PRO
2	J	103	ASP
3	K	70	THR
3	L	70	THR
3	L	85	ARG
3	M	85	ARG
1	B	104	GLU
1	C	57	GLU
1	D	99	ASP
2	G	106	PRO
2	J	20	ARG
3	K	55	SER
2	G	7	PRO
2	H	106	PRO
2	I	7	PRO
2	J	7	PRO
2	J	106	PRO
1	A	69	ASP
2	H	102	PRO
2	H	7	PRO
2	I	102	PRO
2	I	65	SER
2	G	102	PRO

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all 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	87/92 (95%)	73 (84%)	14 (16%)	2	10
1	B	87/92 (95%)	65 (75%)	22 (25%)	0	1
1	C	87/92 (95%)	54 (62%)	33 (38%)	0	0
1	D	87/92 (95%)	61 (70%)	26 (30%)	0	1

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	E	87/92 (95%)	66 (76%)	21 (24%)	1	2
1	F	87/92 (95%)	59 (68%)	28 (32%)	0	0
2	G	94/100 (94%)	69 (73%)	25 (27%)	0	1
2	H	94/100 (94%)	76 (81%)	18 (19%)	1	6
2	I	94/100 (94%)	69 (73%)	25 (27%)	0	1
2	J	94/100 (94%)	64 (68%)	30 (32%)	0	0
3	K	85/101 (84%)	55 (65%)	30 (35%)	0	0
3	L	85/101 (84%)	55 (65%)	30 (35%)	0	0
3	M	85/101 (84%)	58 (68%)	27 (32%)	0	0
3	N	85/101 (84%)	56 (66%)	29 (34%)	0	0
All	All	1238/1356 (91%)	880 (71%)	358 (29%)	0	1

All (358) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	LEU
1	A	26	ASN
1	A	27	MET
1	A	32	ARG
1	A	33	LEU
1	A	38	ASN
1	A	60	GLU
1	A	62	ARG
1	A	73	ARG
1	A	90	LEU
1	A	93	LEU
1	A	94	THR
1	A	98	ARG
1	A	112	ASP
1	B	20	LEU
1	B	22	LEU
1	B	27	MET
1	B	29	VAL
1	B	30	ARG
1	B	33	LEU
1	B	35	LEU
1	B	38	ASN
1	B	42	GLN

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Mol	Chain	Res	Type
1	B	50	LEU
1	B	60	GLU
1	B	61	ILE
1	B	64	LEU
1	B	66	THR
1	B	67	GLN
1	B	73	ARG
1	B	79	GLN
1	B	86	VAL
1	B	88	ARG
1	B	90	LEU
1	B	94	THR
1	B	112	ASP
1	C	20	LEU
1	C	22	LEU
1	C	29	VAL
1	C	32	ARG
1	C	35	LEU
1	C	37	LEU
1	C	40	ARG
1	C	42	GLN
1	C	46	ASP
1	C	48	THR
1	C	50	LEU
1	C	55	ASP
1	C	63	GLN
1	C	64	LEU
1	C	66	THR
1	C	67	GLN
1	C	73	ARG
1	C	75	LEU
1	C	85	SER
1	C	89	LEU
1	C	90	LEU
1	C	92	LEU
1	C	94	THR
1	C	98	ARG
1	C	102	LEU
1	C	103	LEU
1	C	105	LEU
1	C	108	SER
1	C	110	GLU

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Mol	Chain	Res	Type
1	C	111	GLU
1	C	114	GLN
1	C	117	ILE
1	C	121	LEU
1	D	20	LEU
1	D	22	LEU
1	D	25	LEU
1	D	28	ARG
1	D	29	VAL
1	D	32	ARG
1	D	38	ASN
1	D	40	ARG
1	D	43	VAL
1	D	48	THR
1	D	57	GLU
1	D	59	LEU
1	D	61	ILE
1	D	62	ARG
1	D	63	GLN
1	D	64	LEU
1	D	73	ARG
1	D	86	VAL
1	D	91	GLU
1	D	101	VAL
1	D	103	LEU
1	D	111	GLU
1	D	112	ASP
1	D	114	GLN
1	D	117	ILE
1	D	123	HIS
1	E	20	LEU
1	E	22	LEU
1	E	25	LEU
1	E	27	MET
1	E	31	ARG
1	E	33	LEU
1	E	40	ARG
1	E	52	GLU
1	E	59	LEU
1	E	61	ILE
1	E	62	ARG
1	E	63	GLN

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Mol	Chain	Res	Type
1	E	64	LEU
1	E	73	ARG
1	E	86	VAL
1	E	90	LEU
1	E	103	LEU
1	E	108	SER
1	E	114	GLN
1	E	117	ILE
1	E	123	HIS
1	F	20	LEU
1	F	25	LEU
1	F	27	MET
1	F	30	ARG
1	F	32	ARG
1	F	38	ASN
1	F	40	ARG
1	F	41	THR
1	F	42	GLN
1	F	43	VAL
1	F	57	GLU
1	F	59	LEU
1	F	61	ILE
1	F	62	ARG
1	F	63	GLN
1	F	66	THR
1	F	67	GLN
1	F	71	THR
1	F	82	PRO
1	F	86	VAL
1	F	88	ARG
1	F	89	LEU
1	F	90	LEU
1	F	91	GLU
1	F	96	LEU
1	F	102	LEU
1	F	114	GLN
1	F	123	HIS
2	G	2	MET
2	G	6	THR
2	G	9	THR
2	G	14	LEU
2	G	16	VAL

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Mol	Chain	Res	Type
2	G	22	LEU
2	G	27	ASP
2	G	37	VAL
2	G	45	ASP
2	G	51	PHE
2	G	53	ILE
2	G	55	ARG
2	G	61	GLN
2	G	69	GLU
2	G	70	LEU
2	G	71	LEU
2	G	77	THR
2	G	84	LEU
2	G	85	VAL
2	G	86	ASP
2	G	91	ASN
2	G	92	GLU
2	G	99	LEU
2	G	101	LEU
2	G	107	LEU
2	H	6	THR
2	H	13	CYS
2	H	14	LEU
2	H	15	ASN
2	H	16	VAL
2	H	19	ILE
2	H	20	ARG
2	H	22	LEU
2	H	37	VAL
2	H	43	SER
2	H	60	LEU
2	H	62	THR
2	H	64	LYS
2	H	68	SER
2	H	70	LEU
2	H	76	THR
2	H	85	VAL
2	H	101	LEU
2	I	2	MET
2	I	6	THR
2	I	9	THR
2	I	12	ARG

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Mol	Chain	Res	Type
2	I	16	VAL
2	I	18	LEU
2	I	19	ILE
2	I	21	LYS
2	I	22	LEU
2	I	26	ILE
2	I	33	LYS
2	I	37	VAL
2	I	43	SER
2	I	52	HIS
2	I	64	LYS
2	I	67	THR
2	I	70	LEU
2	I	71	LEU
2	I	76	THR
2	I	77	THR
2	I	81	VAL
2	I	87	LEU
2	I	91	ASN
2	I	101	LEU
2	I	105	VAL
2	J	5	ILE
2	J	6	THR
2	J	15	ASN
2	J	16	VAL
2	J	18	LEU
2	J	20	ARG
2	J	22	LEU
2	J	23	SER
2	J	24	ASP
2	J	35	LEU
2	J	37	VAL
2	J	39	ILE
2	J	41	LYS
2	J	43	SER
2	J	46	ASP
2	J	52	HIS
2	J	53	ILE
2	J	61	GLN
2	J	67	THR
2	J	71	LEU
2	J	76	THR

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Mol	Chain	Res	Type
2	J	77	THR
2	J	83	ASP
2	J	85	VAL
2	J	87	LEU
2	J	88	LEU
2	J	90	GLN
2	J	99	LEU
2	J	101	LEU
2	J	107	LEU
3	K	8	LEU
3	K	10	SER
3	K	16	LEU
3	K	18	ARG
3	K	20	MET
3	K	21	ASP
3	K	34	TYR
3	K	36	ILE
3	K	37	THR
3	K	39	LEU
3	K	42	LEU
3	K	45	ILE
3	K	47	SER
3	K	52	GLN
3	K	54	VAL
3	K	56	ILE
3	K	57	THR
3	K	58	ARG
3	K	59	GLU
3	K	61	LEU
3	K	66	MET
3	K	67	ARG
3	K	68	GLN
3	K	72	GLN
3	K	74	LEU
3	K	75	VAL
3	K	81	LEU
3	K	88	GLN
3	K	90	ILE
3	K	91	LEU
3	L	5	ILE
3	L	12	VAL
3	L	13	LEU

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Mol	Chain	Res	Type
3	L	14	ASP
3	L	20	MET
3	L	21	ASP
3	L	23	LEU
3	L	30	GLU
3	L	34	TYR
3	L	35	VAL
3	L	37	THR
3	L	39	LEU
3	L	42	LEU
3	L	44	LYS
3	L	52	GLN
3	L	56	ILE
3	L	58	ARG
3	L	59	GLU
3	L	61	LEU
3	L	68	GLN
3	L	71	VAL
3	L	72	GLN
3	L	74	LEU
3	L	75	VAL
3	L	77	LEU
3	L	78	LEU
3	L	81	LEU
3	L	83	LEU
3	L	84	TYR
3	L	91	LEU
3	M	5	ILE
3	M	13	LEU
3	M	14	ASP
3	M	21	ASP
3	M	37	THR
3	M	38	ASP
3	M	41	GLN
3	M	45	ILE
3	M	47	SER
3	M	54	VAL
3	M	56	ILE
3	M	57	THR
3	M	58	ARG
3	M	59	GLU
3	M	63	TRP

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Mol	Chain	Res	Type
3	M	67	ARG
3	M	70	THR
3	M	71	VAL
3	M	72	GLN
3	M	74	LEU
3	M	78	LEU
3	M	81	LEU
3	M	83	LEU
3	M	84	TYR
3	M	89	ILE
3	M	90	ILE
3	M	94	LYS
3	N	11	TRP
3	N	13	LEU
3	N	20	MET
3	N	21	ASP
3	N	23	LEU
3	N	26	TRP
3	N	38	ASP
3	N	39	LEU
3	N	40	THR
3	N	41	GLN
3	N	43	ARG
3	N	45	ILE
3	N	51	VAL
3	N	54	VAL
3	N	56	ILE
3	N	57	THR
3	N	58	ARG
3	N	59	GLU
3	N	60	LEU
3	N	61	LEU
3	N	63	TRP
3	N	74	LEU
3	N	78	LEU
3	N	80	ARG
3	N	81	LEU
3	N	84	TYR
3	N	85	ARG
3	N	89	ILE
3	N	90	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (51)

such sidechains are listed below:

Mol	Chain	Res	Type
1	A	26	ASN
1	A	38	ASN
1	A	67	GLN
1	A	79	GLN
1	A	123	HIS
1	B	42	GLN
1	B	67	GLN
1	B	79	GLN
1	C	26	ASN
1	C	38	ASN
1	C	42	GLN
1	D	38	ASN
1	D	42	GLN
1	D	79	GLN
1	E	26	ASN
1	E	38	ASN
1	E	67	GLN
1	E	79	GLN
1	F	38	ASN
1	F	42	GLN
1	F	63	GLN
1	F	67	GLN
1	F	114	GLN
2	G	29	GLN
2	G	61	GLN
2	G	90	GLN
2	H	50	GLN
2	H	91	ASN
2	I	50	GLN
2	I	61	GLN
2	I	90	GLN
2	I	91	ASN
2	J	29	GLN
2	J	50	GLN
2	J	61	GLN
3	K	19	ASN
3	K	68	GLN
3	K	73	GLN
3	K	88	GLN
3	L	19	ASN
3	L	68	GLN
3	L	72	GLN

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Mol	Chain	Res	Type
3	M	19	ASN
3	M	41	GLN
3	M	68	GLN
3	M	72	GLN
3	M	92	ASN
3	N	7	GLN
3	N	41	GLN
3	N	73	GLN
3	N	92	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	105/110 (95%)	-0.25	1 (0%) 79 66	103, 190, 331, 374	105 (100%)
1	B	105/110 (95%)	-0.86	0 100 100	80, 138, 223, 329	0
1	C	105/110 (95%)	-0.78	0 100 100	100, 163, 234, 335	0
1	D	105/110 (95%)	-0.68	0 100 100	87, 134, 192, 309	0
1	E	105/110 (95%)	-0.81	1 (0%) 79 66	69, 119, 182, 292	0
1	F	105/110 (95%)	-0.67	0 100 100	85, 162, 254, 320	0
2	G	107/113 (94%)	-0.32	0 100 100	106, 183, 297, 329	0
2	H	107/113 (94%)	-0.71	0 100 100	83, 141, 227, 297	0
2	I	107/113 (94%)	-0.70	0 100 100	91, 145, 253, 307	0
2	J	107/113 (94%)	-0.65	0 100 100	115, 184, 274, 332	0
3	K	93/111 (83%)	-0.38	2 (2%) 62 47	121, 199, 285, 322	0
3	L	93/111 (83%)	-0.56	0 100 100	116, 191, 261, 320	0
3	M	93/111 (83%)	-0.61	0 100 100	126, 218, 286, 317	0
3	N	93/111 (83%)	-0.49	0 100 100	138, 207, 258, 276	0
All	All	1430/1556 (91%)	-0.61	4 (0%) 90 84	69, 168, 278, 374	105 (7%)

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	K	48	MET	2.6
3	K	90	ILE	2.4
1	A	108	SER	2.4
1	E	80	GLY	2.2

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.