



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 11:58 PM UTC

PDB ID : 1FQV / pdb_00001fqv
Title : Insights into scf ubiquitin ligases from the structure of the skp1-skp2 complex
Authors : Schulman, B.A.; Carrano, A.C.; Jeffrey, P.D.; Bowen, Z.; Kinnucan, E.R.;
Finnin, M.S.; Elledge, S.J.; Harper, J.W.; Pagano, M.; Pavletich, N.P.
Deposited on : 2000-09-06
Resolution : 2.80 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

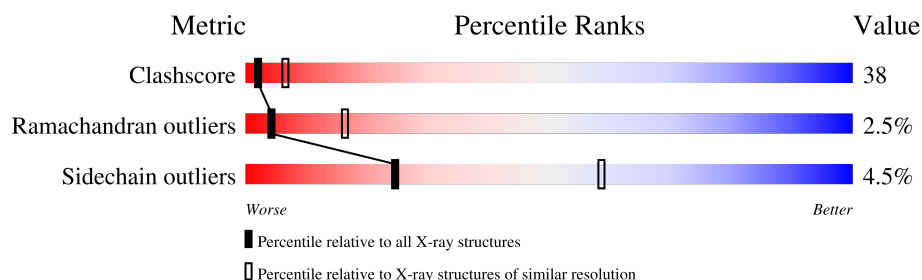
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)

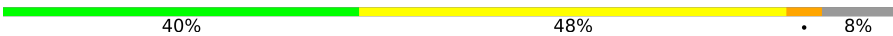
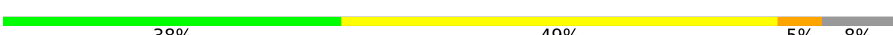
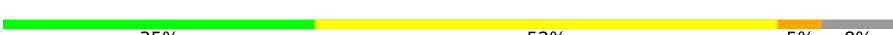
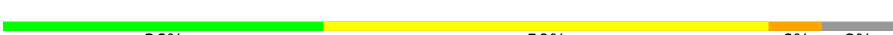
The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	336	44% 45% 7% ..
1	C	336	48% 41% 7% ..
1	E	336	44% 46% 6% ..
1	G	336	46% 44% 5% ..
1	I	336	47% 44% 5% ..
1	K	336	44% 46% 7% ..
1	M	336	46% 44% 6% ..
1	O	336	46% 42% 8% ..

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Mol	Chain	Length	Quality of chain
2	B	149	
2	D	149	
2	F	149	
2	H	149	
2	J	149	
2	L	149	
2	N	149	
2	P	149	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 29256 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 SKP2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	325	Total	C	N	O	S	4	0	0
			2563	1630	441	476	16			
1	C	325	Total	C	N	O	S	4	0	0
			2563	1630	441	476	16			
1	E	325	Total	C	N	O	S	4	0	0
			2563	1630	441	476	16			
1	G	325	Total	C	N	O	S	4	0	0
			2563	1630	441	476	16			
1	I	325	Total	C	N	O	S	4	0	0
			2563	1630	441	476	16			
1	K	325	Total	C	N	O	S	4	0	0
			2563	1630	441	476	16			
1	M	325	Total	C	N	O	S	4	0	0
			2563	1630	441	476	16			
1	O	325	Total	C	N	O	S	4	0	0
			2563	1630	441	476	16			

- Molecule 2 is a protein called SKP1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	137	Total	C	N	O	S	0	0	0
			1094	697	178	213	6			
2	D	137	Total	C	N	O	S	0	0	0
			1094	697	178	213	6			
2	F	137	Total	C	N	O	S	0	0	0
			1094	697	178	213	6			
2	H	137	Total	C	N	O	S	0	0	0
			1094	697	178	213	6			
2	J	137	Total	C	N	O	S	0	0	0
			1094	697	178	213	6			
2	L	137	Total	C	N	O	S	0	0	0
			1094	697	178	213	6			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	N	137	Total	C	N	O	S	0	0	0
			1094	697	178	213	6			
2	P	137	Total	C	N	O	S	0	0	0
			1094	697	178	213	6			

There are 192 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	?	-	ASP	deletion	UNP P63208
B	?	-	ASP	deletion	UNP P63208
B	?	-	GLU	deletion	UNP P63208
B	?	-	GLY	deletion	UNP P63208
B	?	-	ASP	deletion	UNP P63208
B	?	-	ASP	deletion	UNP P63208
B	?	-	PRO	deletion	UNP P63208
B	?	-	PRO	deletion	UNP P63208
B	?	-	PRO	deletion	UNP P63208
B	?	-	GLU	deletion	UNP P63208
B	?	-	ASP	deletion	UNP P63208
B	?	-	ASP	deletion	UNP P63208
B	?	-	GLU	deletion	UNP P63208
B	?	-	ASN	deletion	UNP P63208
B	?	-	LYS	deletion	UNP P63208
B	?	-	GLU	deletion	UNP P63208
B	?	-	LYS	deletion	UNP P63208
B	?	-	ARG	deletion	UNP P63208
B	?	-	THR	deletion	UNP P63208
B	78	GLY	-	SEE REMARK 999	UNP P63208
B	79	GLY	-	SEE REMARK 999	UNP P63208
B	80	SER	-	SEE REMARK 999	UNP P63208
B	81	GLY	-	SEE REMARK 999	UNP P63208
B	82	THR	-	SEE REMARK 999	UNP P63208
D	?	-	ASP	deletion	UNP P63208
D	?	-	ASP	deletion	UNP P63208
D	?	-	GLU	deletion	UNP P63208
D	?	-	GLY	deletion	UNP P63208
D	?	-	ASP	deletion	UNP P63208
D	?	-	ASP	deletion	UNP P63208
D	?	-	PRO	deletion	UNP P63208
D	?	-	PRO	deletion	UNP P63208
D	?	-	PRO	deletion	UNP P63208
D	?	-	GLU	deletion	UNP P63208

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Chain	Residue	Modelled	Actual	Comment	Reference
D	?	-	ASP	deletion	UNP P63208
D	?	-	ASP	deletion	UNP P63208
D	?	-	GLU	deletion	UNP P63208
D	?	-	ASN	deletion	UNP P63208
D	?	-	LYS	deletion	UNP P63208
D	?	-	GLU	deletion	UNP P63208
D	?	-	LYS	deletion	UNP P63208
D	?	-	ARG	deletion	UNP P63208
D	?	-	THR	deletion	UNP P63208
D	78	GLY	-	SEE REMARK 999	UNP P63208
D	79	GLY	-	SEE REMARK 999	UNP P63208
D	80	SER	-	SEE REMARK 999	UNP P63208
D	81	GLY	-	SEE REMARK 999	UNP P63208
D	82	THR	-	SEE REMARK 999	UNP P63208
F	?	-	ASP	deletion	UNP P63208
F	?	-	ASP	deletion	UNP P63208
F	?	-	GLU	deletion	UNP P63208
F	?	-	GLY	deletion	UNP P63208
F	?	-	ASP	deletion	UNP P63208
F	?	-	ASP	deletion	UNP P63208
F	?	-	PRO	deletion	UNP P63208
F	?	-	PRO	deletion	UNP P63208
F	?	-	PRO	deletion	UNP P63208
F	?	-	GLU	deletion	UNP P63208
F	?	-	ASP	deletion	UNP P63208
F	?	-	ASP	deletion	UNP P63208
F	?	-	GLU	deletion	UNP P63208
F	?	-	ASN	deletion	UNP P63208
F	?	-	LYS	deletion	UNP P63208
F	?	-	GLU	deletion	UNP P63208
F	?	-	LYS	deletion	UNP P63208
F	?	-	ARG	deletion	UNP P63208
F	?	-	THR	deletion	UNP P63208
F	78	GLY	-	SEE REMARK 999	UNP P63208
F	79	GLY	-	SEE REMARK 999	UNP P63208
F	80	SER	-	SEE REMARK 999	UNP P63208
F	81	GLY	-	SEE REMARK 999	UNP P63208
F	82	THR	-	SEE REMARK 999	UNP P63208
H	?	-	ASP	deletion	UNP P63208
H	?	-	ASP	deletion	UNP P63208
H	?	-	GLU	deletion	UNP P63208
H	?	-	GLY	deletion	UNP P63208

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Chain	Residue	Modelled	Actual	Comment	Reference
H	?	-	ASP	deletion	UNP P63208
H	?	-	ASP	deletion	UNP P63208
H	?	-	PRO	deletion	UNP P63208
H	?	-	PRO	deletion	UNP P63208
H	?	-	PRO	deletion	UNP P63208
H	?	-	GLU	deletion	UNP P63208
H	?	-	ASP	deletion	UNP P63208
H	?	-	ASP	deletion	UNP P63208
H	?	-	GLU	deletion	UNP P63208
H	?	-	ASN	deletion	UNP P63208
H	?	-	LYS	deletion	UNP P63208
H	?	-	GLU	deletion	UNP P63208
H	?	-	LYS	deletion	UNP P63208
H	?	-	ARG	deletion	UNP P63208
H	?	-	THR	deletion	UNP P63208
H	78	GLY	-	SEE REMARK 999	UNP P63208
H	79	GLY	-	SEE REMARK 999	UNP P63208
H	80	SER	-	SEE REMARK 999	UNP P63208
H	81	GLY	-	SEE REMARK 999	UNP P63208
H	82	THR	-	SEE REMARK 999	UNP P63208
J	?	-	ASP	deletion	UNP P63208
J	?	-	ASP	deletion	UNP P63208
J	?	-	GLU	deletion	UNP P63208
J	?	-	GLY	deletion	UNP P63208
J	?	-	ASP	deletion	UNP P63208
J	?	-	ASP	deletion	UNP P63208
J	?	-	PRO	deletion	UNP P63208
J	?	-	PRO	deletion	UNP P63208
J	?	-	PRO	deletion	UNP P63208
J	?	-	GLU	deletion	UNP P63208
J	?	-	ASP	deletion	UNP P63208
J	?	-	ASP	deletion	UNP P63208
J	?	-	GLU	deletion	UNP P63208
J	?	-	ASN	deletion	UNP P63208
J	?	-	LYS	deletion	UNP P63208
J	?	-	GLU	deletion	UNP P63208
J	?	-	LYS	deletion	UNP P63208
J	?	-	ARG	deletion	UNP P63208
J	?	-	THR	deletion	UNP P63208
J	78	GLY	-	SEE REMARK 999	UNP P63208
J	79	GLY	-	SEE REMARK 999	UNP P63208
J	80	SER	-	SEE REMARK 999	UNP P63208

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Chain	Residue	Modelled	Actual	Comment	Reference
J	81	GLY	-	SEE REMARK 999	UNP P63208
J	82	THR	-	SEE REMARK 999	UNP P63208
L	?	-	ASP	deletion	UNP P63208
L	?	-	ASP	deletion	UNP P63208
L	?	-	GLU	deletion	UNP P63208
L	?	-	GLY	deletion	UNP P63208
L	?	-	ASP	deletion	UNP P63208
L	?	-	ASP	deletion	UNP P63208
L	?	-	PRO	deletion	UNP P63208
L	?	-	PRO	deletion	UNP P63208
L	?	-	PRO	deletion	UNP P63208
L	?	-	GLU	deletion	UNP P63208
L	?	-	ASP	deletion	UNP P63208
L	?	-	ASP	deletion	UNP P63208
L	?	-	GLU	deletion	UNP P63208
L	?	-	ASN	deletion	UNP P63208
L	?	-	LYS	deletion	UNP P63208
L	?	-	GLU	deletion	UNP P63208
L	?	-	LYS	deletion	UNP P63208
L	?	-	ARG	deletion	UNP P63208
L	?	-	THR	deletion	UNP P63208
L	78	GLY	-	SEE REMARK 999	UNP P63208
L	79	GLY	-	SEE REMARK 999	UNP P63208
L	80	SER	-	SEE REMARK 999	UNP P63208
L	81	GLY	-	SEE REMARK 999	UNP P63208
L	82	THR	-	SEE REMARK 999	UNP P63208
N	?	-	ASP	deletion	UNP P63208
N	?	-	ASP	deletion	UNP P63208
N	?	-	GLU	deletion	UNP P63208
N	?	-	GLY	deletion	UNP P63208
N	?	-	ASP	deletion	UNP P63208
N	?	-	ASP	deletion	UNP P63208
N	?	-	PRO	deletion	UNP P63208
N	?	-	PRO	deletion	UNP P63208
N	?	-	PRO	deletion	UNP P63208
N	?	-	GLU	deletion	UNP P63208
N	?	-	ASP	deletion	UNP P63208
N	?	-	ASP	deletion	UNP P63208
N	?	-	GLU	deletion	UNP P63208
N	?	-	ASN	deletion	UNP P63208
N	?	-	LYS	deletion	UNP P63208
N	?	-	GLU	deletion	UNP P63208

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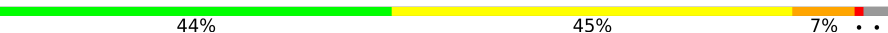
Chain	Residue	Modelled	Actual	Comment	Reference
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N	?	-	ARG	deletion	UNP P63208
N	?	-	THR	deletion	UNP P63208
N	78	GLY	-	SEE REMARK 999	UNP P63208
N	79	GLY	-	SEE REMARK 999	UNP P63208
N	80	SER	-	SEE REMARK 999	UNP P63208
N	81	GLY	-	SEE REMARK 999	UNP P63208
N	82	THR	-	SEE REMARK 999	UNP P63208
P	?	-	ASP	deletion	UNP P63208
P	?	-	ASP	deletion	UNP P63208
P	?	-	GLU	deletion	UNP P63208
P	?	-	GLY	deletion	UNP P63208
P	?	-	ASP	deletion	UNP P63208
P	?	-	ASP	deletion	UNP P63208
P	?	-	PRO	deletion	UNP P63208
P	?	-	PRO	deletion	UNP P63208
P	?	-	PRO	deletion	UNP P63208
P	?	-	GLU	deletion	UNP P63208
P	?	-	ASP	deletion	UNP P63208
P	?	-	ASP	deletion	UNP P63208
P	?	-	GLU	deletion	UNP P63208
P	?	-	ASN	deletion	UNP P63208
P	?	-	LYS	deletion	UNP P63208
P	?	-	GLU	deletion	UNP P63208
P	?	-	LYS	deletion	UNP P63208
P	?	-	ARG	deletion	UNP P63208
P	?	-	THR	deletion	UNP P63208
P	78	GLY	-	SEE REMARK 999	UNP P63208
P	79	GLY	-	SEE REMARK 999	UNP P63208
P	80	SER	-	SEE REMARK 999	UNP P63208
P	81	GLY	-	SEE REMARK 999	UNP P63208
P	82	THR	-	SEE REMARK 999	UNP P63208

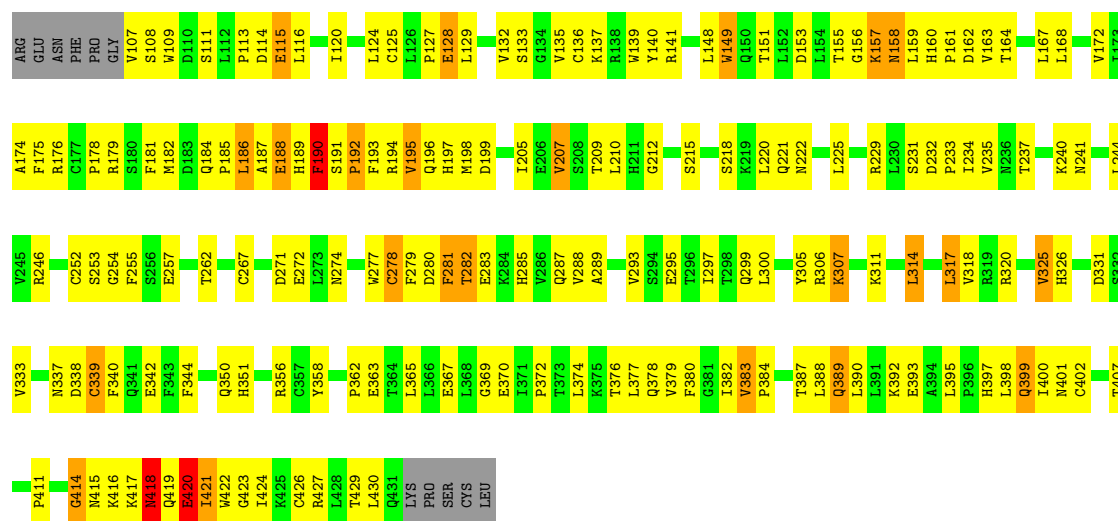
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

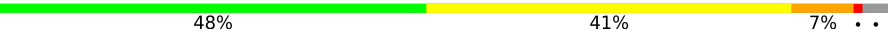
Note EDS was not executed.

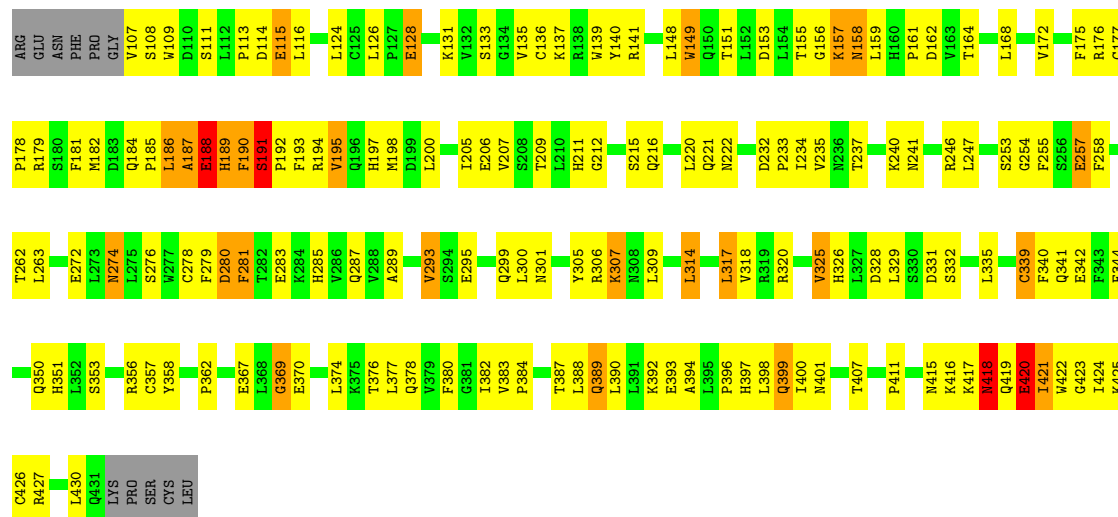
• Molecule 1: SKP2

Chain A: 



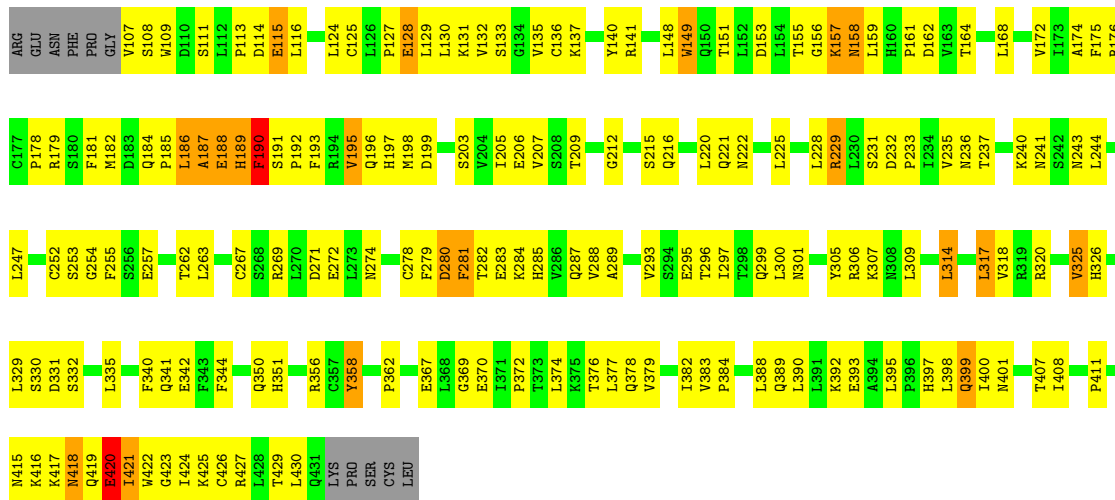
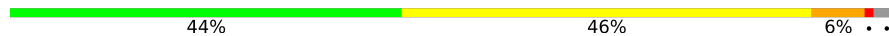
• Molecule 1: SKP2

Chain C: 



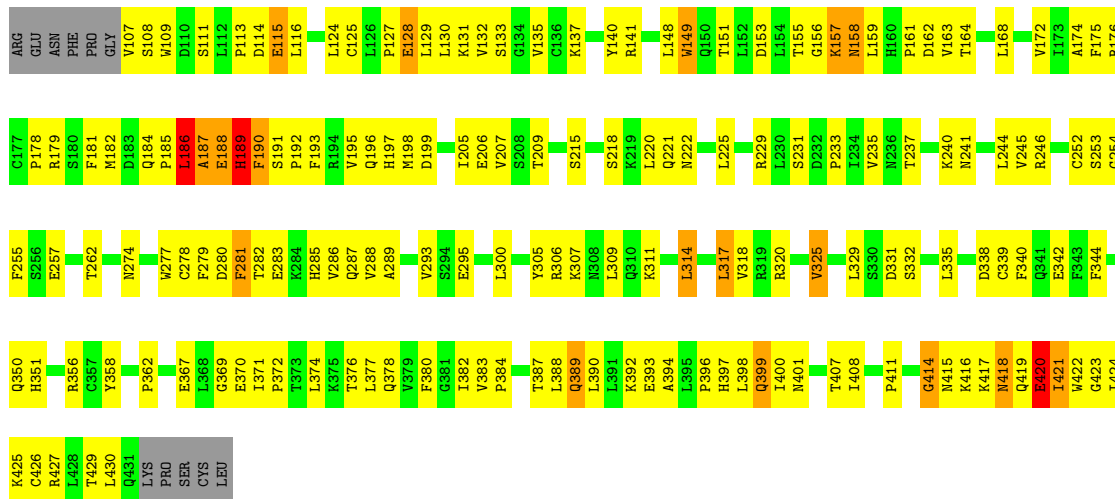
- Molecule 1: SKP2

Chain E:



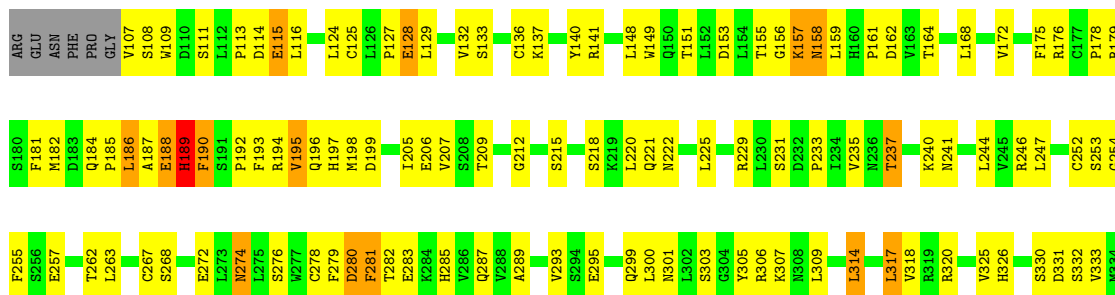
- Molecule 1: SKP2

Chain G:



- Molecule 1: SKP2

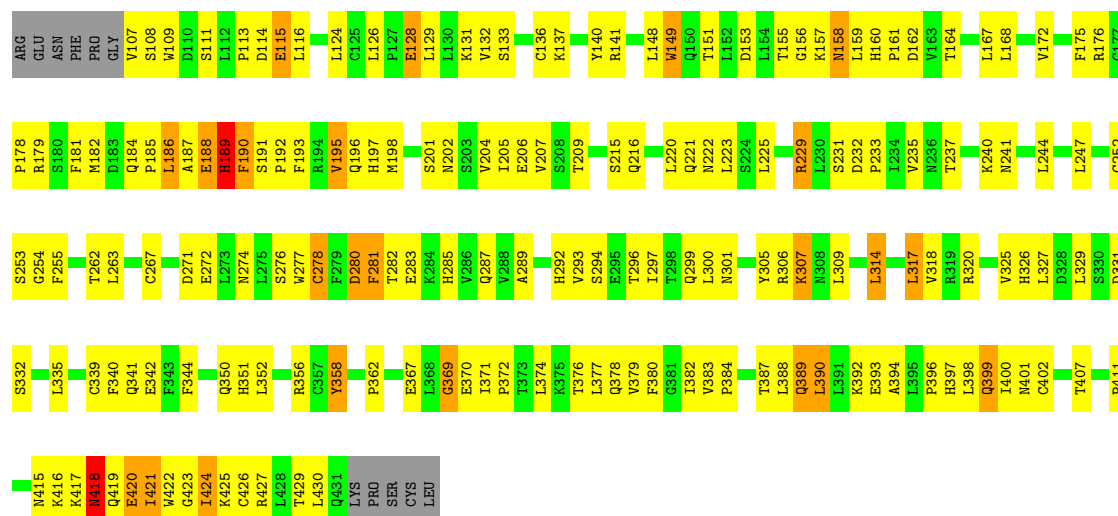
Chain I:





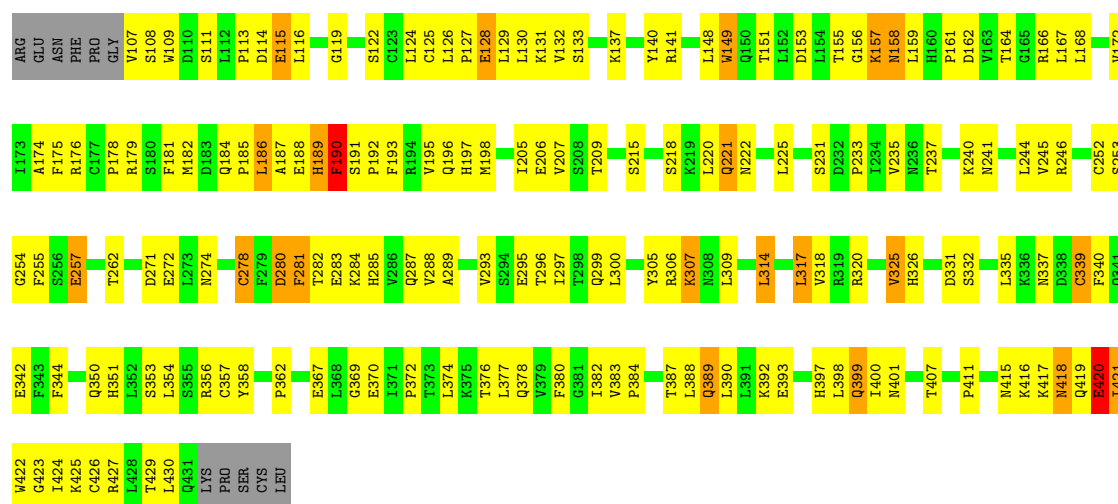
• Molecule 1: SKP2

Chain K: 44% 46% 7% ..



• Molecule 1: SKP2

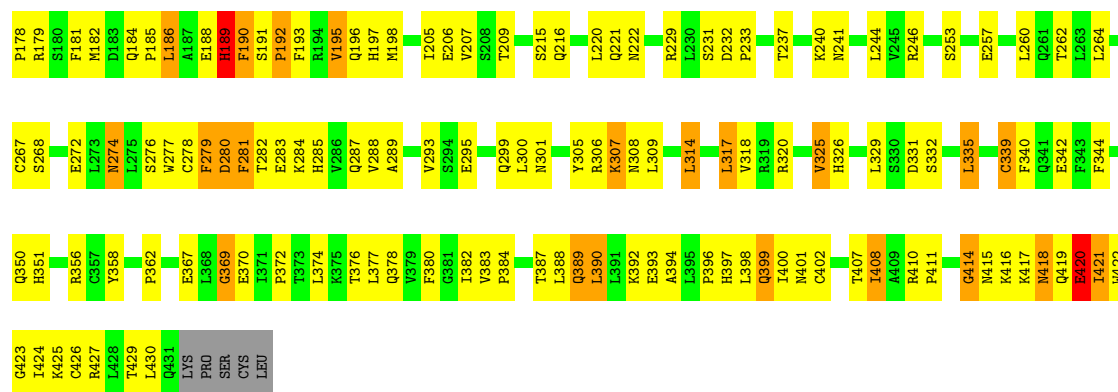
Chain M: 46% 44% 6% ..



• Molecule 1: SKP2

Chain O: 46% 42% 8% ..

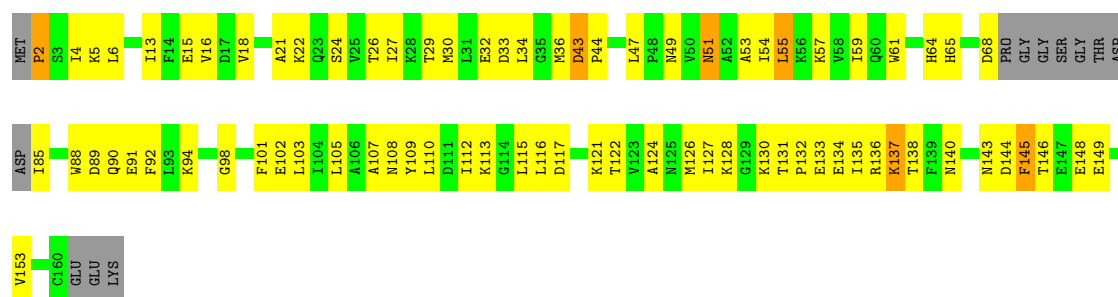




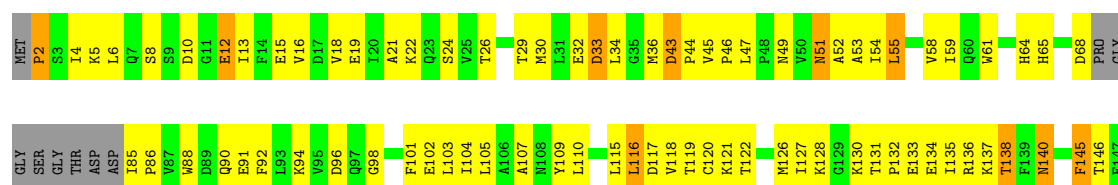
• Molecule 2: SKP1



• Molecule 2: SKP1



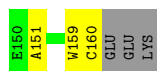
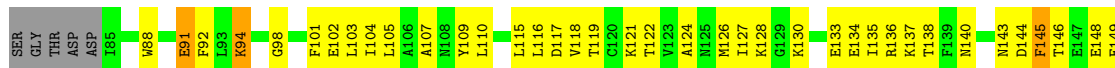
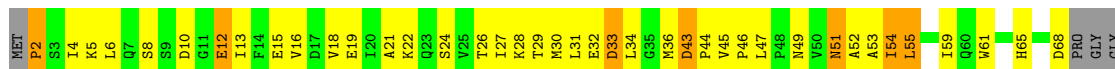
• Molecule 2: SKP1





• Molecule 2: SKP1

Chain H: 38% 48% 7% 8%



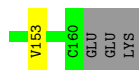
• Molecule 2: SKP1

Chain J: 37% 49% 6% 8%



• Molecule 2: SKP1

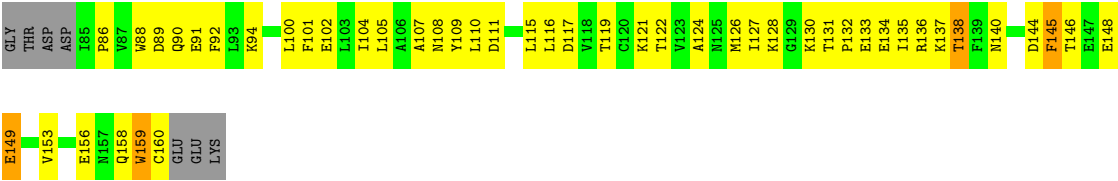
Chain L: 38% 49% 5% 8%



• Molecule 2: SKP1

Chain N: 35% 52% 5% 8%





• Molecule 2: SKP1



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	262.70Å 148.20Å 133.30Å 90.00° 120.03° 90.00°	Depositor
Resolution (Å)	8.00 – 2.80	Depositor
% Data completeness (in resolution range)	(Not available) (8.00-2.80)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS	Depositor
R, R_{free}	0.275 , 0.314	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	29256	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	A	0.73	3/2612 (0.1%)	1.16	15/3545 (0.4%)
1	C	0.81	5/2612 (0.2%)	1.26	21/3545 (0.6%)
1	E	0.64	1/2612 (0.0%)	1.14	14/3545 (0.4%)
1	G	0.86	3/2612 (0.1%)	1.45	20/3545 (0.6%)
1	I	0.81	2/2612 (0.1%)	1.40	17/3545 (0.5%)
1	K	0.77	1/2612 (0.0%)	1.43	26/3545 (0.7%)
1	M	0.75	4/2612 (0.2%)	1.21	18/3545 (0.5%)
1	O	0.69	2/2612 (0.1%)	1.44	23/3545 (0.6%)
2	B	0.41	0/1111	0.92	3/1502 (0.2%)
2	D	0.50	0/1111	0.97	2/1502 (0.1%)
2	F	0.45	0/1111	0.96	3/1502 (0.2%)
2	H	0.72	0/1111	1.09	7/1502 (0.5%)
2	J	0.52	0/1111	1.02	6/1502 (0.4%)
2	L	0.52	0/1111	1.01	5/1502 (0.3%)
2	N	0.53	0/1111	0.97	3/1502 (0.2%)
2	P	0.68	0/1111	1.01	4/1502 (0.3%)
All	All	0.70	21/29784 (0.1%)	1.23	187/40376 (0.5%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	I	0	1

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	G	186	LEU	C-N	26.45	1.70	1.33
1	I	186	LEU	C-N	21.26	1.63	1.33
1	G	189	HIS	C-N	-15.10	1.12	1.33
1	K	189	HIS	C-N	-14.44	1.13	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	M	186	LEU	C-N	12.73	1.51	1.33
1	A	190	PHE	N-CA	-11.83	1.30	1.46
1	O	189	HIS	C-N	-11.31	1.17	1.33
1	C	189	HIS	C-N	-10.80	1.19	1.33
1	M	189	HIS	CA-C	-10.63	1.38	1.52
1	M	190	PHE	N-CA	-8.95	1.34	1.46
1	I	189	HIS	C-N	-8.95	1.21	1.33
1	A	189	HIS	CA-C	-8.10	1.43	1.52
1	C	189	HIS	N-CA	-7.60	1.37	1.46
1	E	190	PHE	N-CA	-7.47	1.36	1.46
1	C	188	GLU	C-N	-6.88	1.24	1.33
1	C	190	PHE	CA-C	-6.26	1.44	1.53
1	C	190	PHE	C-N	-6.00	1.21	1.33
1	A	189	HIS	C-N	-5.96	1.25	1.33
1	M	189	HIS	C-N	-5.60	1.25	1.33
1	O	186	LEU	C-N	5.29	1.41	1.33
1	G	191	SER	CA-C	5.07	1.59	1.53

All (187) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	I	189	HIS	CA-C-N	-28.95	66.25	121.54
1	I	189	HIS	C-N-CA	-28.95	66.25	121.54
1	G	189	HIS	O-C-N	27.78	159.53	122.59
1	K	186	LEU	O-C-N	26.28	157.54	122.59
1	O	189	HIS	CA-C-N	-25.57	72.70	121.54
1	O	189	HIS	C-N-CA	-25.57	72.70	121.54
1	G	189	HIS	CA-C-N	-25.03	73.73	121.54
1	G	189	HIS	C-N-CA	-25.03	73.73	121.54
1	G	186	LEU	O-C-N	24.42	155.06	122.59
1	K	189	HIS	CA-C-N	-23.77	76.14	121.54
1	K	189	HIS	C-N-CA	-23.77	76.14	121.54
1	O	186	LEU	O-C-N	23.70	154.11	122.59
1	M	186	LEU	O-C-N	23.66	154.06	122.59
1	O	189	HIS	O-C-N	22.76	152.87	122.59
1	I	189	HIS	O-C-N	22.58	152.62	122.59
1	K	189	HIS	O-C-N	22.16	152.06	122.59
1	C	190	PHE	N-CA-CB	-19.24	84.09	110.25
1	I	186	LEU	O-C-N	14.66	142.09	122.59
1	A	281	PHE	N-CA-C	14.62	132.19	109.81
1	K	281	PHE	N-CA-C	14.17	131.50	109.81
1	C	281	PHE	N-CA-C	14.05	131.31	109.81

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	281	PHE	N-CA-C	13.92	131.10	109.81
1	I	281	PHE	N-CA-C	13.59	130.60	109.81
1	M	281	PHE	N-CA-C	13.29	130.14	109.81
1	G	281	PHE	N-CA-C	13.28	130.83	109.85
1	C	189	HIS	N-CA-C	13.23	129.61	112.41
1	O	281	PHE	N-CA-C	12.90	129.54	109.81
1	O	186	LEU	CA-C-N	-12.32	101.12	122.07
1	O	186	LEU	C-N-CA	-12.32	101.12	122.07
1	G	186	LEU	CA-C-N	-10.76	100.98	121.54
1	G	186	LEU	C-N-CA	-10.76	100.98	121.54
2	N	145	PHE	N-CA-C	10.10	122.83	110.41
1	K	186	LEU	CA-C-N	-9.92	102.60	121.54
1	K	186	LEU	C-N-CA	-9.92	102.60	121.54
2	F	145	PHE	N-CA-C	9.11	121.61	110.41
2	L	11	GLY	N-CA-C	8.62	125.45	114.25
1	A	420	GLU	N-CA-C	-8.52	99.80	111.56
1	I	158	ASN	N-CA-C	-8.52	97.68	110.28
1	C	191	SER	N-CA-C	8.26	128.06	109.81
1	C	189	HIS	CB-CA-C	-8.19	97.75	111.02
1	A	158	ASN	N-CA-C	-8.15	97.32	110.20
1	I	420	GLU	N-CA-C	-8.15	100.31	111.56
1	E	158	ASN	N-CA-C	-8.12	98.27	110.28
2	L	144	ASP	N-CA-C	8.05	119.83	111.14
1	M	186	LEU	CA-C-N	-8.03	106.20	121.54
1	M	186	LEU	C-N-CA	-8.03	106.20	121.54
1	G	420	GLU	N-CA-C	-7.99	100.54	111.56
1	M	420	GLU	N-CA-C	-7.94	100.60	111.56
2	H	12	GLU	N-CA-C	7.93	121.73	110.23
2	H	145	PHE	N-CA-C	7.88	122.02	110.59
1	C	188	GLU	O-C-N	7.84	133.02	122.59
1	O	420	GLU	N-CA-C	-7.77	100.84	111.56
1	E	420	GLU	N-CA-C	-7.75	100.86	111.56
1	M	407	THR	N-CA-C	-7.70	104.38	114.31
1	C	420	GLU	N-CA-C	-7.66	100.98	111.56
1	O	158	ASN	N-CA-C	-7.64	98.12	110.20
1	G	158	ASN	N-CA-C	-7.63	98.14	110.20
2	D	145	PHE	N-CA-C	7.51	120.22	110.53
1	K	158	ASN	N-CA-C	-7.43	98.46	110.20
1	M	158	ASN	N-CA-C	-7.39	98.53	110.20
2	B	145	PHE	N-CA-C	7.28	120.38	110.55
1	K	420	GLU	N-CA-C	-7.01	101.88	111.56
1	C	158	ASN	N-CA-C	-6.99	99.16	110.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	189	HIS	N-CA-C	-6.94	104.09	114.64
1	C	325	VAL	N-CA-C	-6.89	106.32	111.90
1	A	229	ARG	N-CA-C	-6.88	100.10	110.28
1	C	149	TRP	N-CA-C	6.69	120.75	112.59
1	A	218	SER	N-CA-C	-6.67	107.01	114.62
1	C	190	PHE	CA-CB-CG	6.63	120.43	113.80
2	J	145	PHE	N-CA-C	6.61	120.27	110.48
1	O	229	ARG	N-CA-C	-6.60	100.52	110.28
1	C	407	THR	N-CA-C	-6.60	104.95	114.12
1	K	282	THR	N-CA-C	6.59	118.67	109.31
1	K	149	TRP	N-CA-C	6.56	120.59	112.59
1	I	192	PRO	CB-CA-C	6.55	119.77	111.39
2	N	2	PRO	N-CA-CB	6.54	110.20	103.00
1	M	280	ASP	CA-C-N	-6.50	111.39	123.03
1	M	280	ASP	C-N-CA	-6.50	111.39	123.03
2	D	2	PRO	N-CA-CB	6.42	110.07	103.00
1	E	149	TRP	N-CA-C	6.40	120.40	112.59
1	E	407	THR	N-CA-C	-6.37	105.45	113.72
1	A	189	HIS	N-CA-C	-6.36	104.75	114.16
2	F	2	PRO	N-CA-CB	6.28	109.91	103.00
2	J	2	PRO	N-CA-CB	6.28	109.91	103.00
2	L	2	PRO	N-CA-CB	6.27	109.90	103.00
2	B	2	PRO	N-CA-CB	6.27	109.89	103.00
1	O	339	CYS	N-CA-C	-6.26	105.78	113.41
1	O	325	VAL	N-CA-C	-6.21	106.75	112.96
2	J	85	ILE	CA-C-N	6.17	126.19	119.89
2	J	85	ILE	C-N-CA	6.17	126.19	119.89
1	G	218	SER	N-CA-C	-6.09	107.67	114.62
1	G	192	PRO	CB-CA-C	6.08	119.28	111.56
2	H	91	GLU	N-CA-C	-6.03	104.79	111.36
1	A	278	CYS	N-CA-C	-6.02	97.03	108.17
2	H	2	PRO	N-CA-CB	6.01	109.61	103.00
1	I	407	THR	N-CA-C	-6.00	105.78	114.12
1	O	268	SER	N-CA-C	6.00	120.33	113.19
2	P	2	PRO	N-CA-CB	5.96	109.55	103.00
1	K	339	CYS	N-CA-C	-5.95	106.15	113.41
1	G	229	ARG	N-CA-C	-5.95	101.47	110.28
1	O	282	THR	N-CA-C	5.94	119.07	109.86
1	A	339	CYS	N-CA-C	-5.92	106.21	113.50
1	O	369	GLY	N-CA-C	-5.92	107.56	115.32
1	M	218	SER	N-CA-C	-5.91	107.88	114.62
1	A	189	HIS	O-C-N	5.88	127.91	121.85

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	O	414	GLY	N-CA-C	5.88	118.39	111.63
2	L	146	THR	N-CA-C	-5.86	100.15	109.59
1	K	369	GLY	N-CA-C	-5.82	107.69	115.32
1	G	339	CYS	N-CA-C	-5.81	106.32	113.41
2	P	145	PHE	N-CA-C	5.79	118.37	110.55
1	M	149	TRP	N-CA-C	5.79	119.94	112.41
1	O	407	THR	N-CA-C	-5.77	106.22	113.72
1	M	190	PHE	N-CA-C	-5.75	98.56	110.80
1	M	335	LEU	N-CA-C	5.72	118.62	110.10
1	K	390	LEU	N-CA-C	-5.70	105.15	111.36
1	G	149	TRP	N-CA-C	5.68	119.76	112.26
1	K	292	HIS	N-CA-C	5.67	120.93	112.94
1	G	325	VAL	N-CA-C	-5.66	107.30	112.96
1	K	407	THR	N-CA-C	-5.64	106.39	113.72
1	I	268	SER	N-CA-C	5.64	119.90	113.19
1	M	339	CYS	N-CA-C	-5.60	106.61	113.50
2	J	146	THR	N-CA-C	-5.60	101.16	110.17
2	P	12	GLU	N-CA-C	5.59	117.95	110.35
1	C	339	CYS	N-CA-C	-5.57	106.61	113.41
1	O	149	TRP	N-CA-C	5.56	119.60	112.26
1	K	280	ASP	CA-C-N	-5.55	113.09	123.03
1	K	280	ASP	C-N-CA	-5.55	113.09	123.03
2	H	54	ILE	N-CA-C	5.55	116.33	110.72
2	F	12	GLU	N-CA-C	5.54	117.44	110.24
2	H	94	LYS	N-CA-C	-5.51	101.35	109.62
2	P	146	THR	N-CA-C	-5.49	101.22	109.95
1	A	149	TRP	N-CA-C	5.49	119.28	112.59
1	A	282	THR	N-CA-C	5.47	118.35	109.86
1	I	188	GLU	N-CA-C	5.47	122.45	110.80
2	L	91	GLU	N-CA-C	-5.46	105.40	111.36
1	K	229	ARG	N-CA-C	-5.46	101.77	109.96
1	A	383	VAL	N-CA-C	5.46	113.95	107.73
1	E	325	VAL	N-CA-C	-5.45	107.51	112.96
1	E	280	ASP	CA-C-N	-5.39	113.38	123.03
1	E	280	ASP	C-N-CA	-5.39	113.38	123.03
1	E	335	LEU	N-CA-C	5.38	117.98	109.96
1	G	407	THR	N-CA-C	-5.34	106.78	113.72
1	E	358	TYR	N-CA-C	5.34	119.28	112.34
1	C	257	GLU	N-CA-C	-5.33	106.73	113.18
1	K	358	TYR	N-CA-C	5.33	118.88	112.38
1	K	201	SER	N-CA-C	5.32	117.95	110.23
2	H	116	LEU	N-CA-C	-5.32	105.57	111.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	414	GLY	N-CA-C	5.31	117.74	111.63
1	O	280	ASP	CA-C-N	-5.29	113.55	123.03
1	O	280	ASP	C-N-CA	-5.29	113.55	123.03
1	I	218	SER	N-CA-C	-5.29	108.09	114.75
1	K	371	ILE	CA-C-N	5.28	124.92	119.05
1	K	371	ILE	C-N-CA	5.28	124.92	119.05
1	E	229	ARG	N-CA-C	-5.28	102.47	110.28
1	O	390	LEU	N-CA-C	-5.28	105.58	112.23
1	C	190	PHE	N-CA-C	-5.27	103.66	110.19
1	G	371	ILE	CA-C-N	5.27	124.45	118.97
1	G	371	ILE	C-N-CA	5.27	124.45	118.97
2	B	12	GLU	N-CA-C	5.24	117.82	110.23
1	I	280	ASP	CA-C-N	-5.24	113.66	123.03
1	I	280	ASP	C-N-CA	-5.24	113.66	123.03
1	C	293	VAL	N-CA-C	-5.23	102.83	109.58
1	M	257	GLU	N-CA-C	-5.23	106.85	113.18
2	J	12	GLU	N-CA-C	5.22	117.61	110.24
1	A	407	THR	N-CA-C	-5.22	106.31	113.30
2	N	149	GLU	N-CA-C	-5.20	105.22	112.45
1	I	414	GLY	N-CA-C	5.19	117.60	111.63
1	O	192	PRO	CB-CA-C	5.19	118.16	111.56
1	C	280	ASP	CA-C-N	-5.18	113.75	123.03
1	C	280	ASP	C-N-CA	-5.18	113.75	123.03
1	C	369	GLY	N-CA-C	-5.16	108.56	115.32
1	K	335	LEU	N-CA-C	5.15	117.77	110.10
1	K	216	GLN	N-CA-C	-5.14	105.73	112.41
1	O	335	LEU	N-CA-C	5.14	117.75	110.10
1	C	216	GLN	N-CA-C	-5.13	105.71	112.94
1	G	335	LEU	N-CA-C	5.13	117.60	109.96
1	A	325	VAL	N-CA-C	-5.10	107.86	112.96
1	M	278	CYS	N-CA-C	-5.09	98.41	107.98
1	M	325	VAL	N-CA-C	-5.09	107.87	112.96
1	E	282	THR	N-CA-C	5.06	117.71	109.86
1	G	414	GLY	N-CA-C	5.06	118.15	110.97
1	E	189	HIS	O-C-N	5.06	127.29	121.93
1	I	229	ARG	N-CA-C	-5.05	102.58	110.10
1	I	257	GLU	N-CA-C	-5.05	107.07	113.18
1	C	357	CYS	N-CA-C	-5.04	99.11	108.24
1	M	221	GLN	N-CA-C	-5.04	107.20	113.55
1	K	278	CYS	N-CA-C	-5.03	98.86	108.17

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	I	189	HIS	Mainchain

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2563	0	2598	230	0
1	C	2563	0	2598	216	0
1	E	2563	0	2598	225	0
1	G	2563	0	2596	223	0
1	I	2563	0	2598	196	0
1	K	2563	0	2598	230	0
1	M	2563	0	2598	205	0
1	O	2563	0	2598	226	0
2	B	1094	0	1099	91	0
2	D	1094	0	1099	79	0
2	F	1094	0	1099	89	1
2	H	1094	0	1099	75	0
2	J	1094	0	1099	78	0
2	L	1094	0	1099	86	1
2	N	1094	0	1099	81	0
2	P	1094	0	1099	93	0
All	All	29256	0	29574	2224	1

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 38.

All (2224) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:186:LEU:C	1:G:187:ALA:N	1.70	1.44
1:M:189:HIS:O	1:M:190:PHE:O	1.58	1.17
1:K:417:LYS:CE	1:O:417:LYS:HG3	1.75	1.15
1:I:417:LYS:HG3	1:M:417:LYS:CE	1.77	1.14
2:L:43:ASP:HB2	2:L:44:PRO:HD3	1.29	1.14

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:416:LYS:HD2	1:I:417:LYS:HZ1	1.10	1.13
1:M:189:HIS:O	1:M:190:PHE:C	1.88	1.11
2:H:4:ILE:HG12	2:H:18:VAL:HG12	1.33	1.10
1:C:188:GLU:O	1:C:212:GLY:O	1.69	1.10
1:K:148:LEU:HA	1:K:427:ARG:HH22	1.14	1.09
1:A:417:LYS:HG3	1:E:417:LYS:CE	1.81	1.09
1:C:417:LYS:CE	1:G:417:LYS:HG3	1.84	1.07
1:E:215:SER:HA	1:E:237:THR:HG21	1.36	1.07
1:G:186:LEU:CA	1:G:187:ALA:N	2.18	1.07
2:N:65:HIS:HB3	2:N:68:ASP:HB2	1.31	1.07
1:O:416:LYS:HD2	1:O:417:LYS:HZ1	1.13	1.06
1:A:416:LYS:HB3	1:A:417:LYS:HE2	1.38	1.06
1:E:189:HIS:O	1:E:190:PHE:O	1.73	1.05
2:J:4:ILE:HG12	2:J:18:VAL:HG12	1.37	1.05
1:M:189:HIS:C	1:M:190:PHE:O	1.91	1.05
2:P:43:ASP:HB2	2:P:44:PRO:HD3	1.38	1.05
1:I:417:LYS:HD2	1:M:417:LYS:HG3	1.36	1.04
1:A:215:SER:HA	1:A:237:THR:HG21	1.39	1.04
2:B:43:ASP:HB2	2:B:44:PRO:HD3	1.40	1.04
1:C:417:LYS:HE3	1:G:417:LYS:HG3	1.05	1.04
2:B:65:HIS:HB3	2:B:68:ASP:HB2	1.35	1.04
2:N:43:ASP:HB2	2:N:44:PRO:HD3	1.38	1.04
2:F:65:HIS:HB3	2:F:68:ASP:HB2	1.40	1.03
2:J:43:ASP:HB2	2:J:44:PRO:HD3	1.40	1.03
1:K:416:LYS:HB3	1:K:417:LYS:HE2	1.37	1.03
1:O:416:LYS:HB3	1:O:417:LYS:HE2	1.39	1.03
2:L:65:HIS:HB3	2:L:68:ASP:HB2	1.41	1.03
1:A:416:LYS:HD2	1:A:417:LYS:HZ1	1.17	1.03
2:P:4:ILE:HG12	2:P:18:VAL:HG12	1.39	1.03
2:P:65:HIS:HB3	2:P:68:ASP:HB2	1.38	1.02
2:D:4:ILE:HG12	2:D:18:VAL:HG12	1.39	1.02
2:F:43:ASP:HB2	2:F:44:PRO:HD3	1.37	1.02
1:K:153:ASP:OD1	1:K:155:THR:HG23	1.60	1.02
2:D:43:ASP:HB2	2:D:44:PRO:HD3	1.41	1.02
2:F:96:ASP:OD1	1:O:284:LYS:HE3	1.59	1.02
1:K:148:LEU:HA	1:K:427:ARG:NH2	1.74	1.02
2:B:4:ILE:HG12	2:B:18:VAL:HG12	1.41	1.02
1:C:416:LYS:HB3	1:C:417:LYS:HE2	1.39	1.01
1:C:162:ASP:HA	1:C:190:PHE:HZ	1.26	1.01
1:A:289:ALA:O	1:A:293:VAL:HG23	1.59	1.01
1:C:148:LEU:HA	1:C:427:ARG:HH22	1.23	1.01

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:155:THR:HG22	1:G:178:PRO:HD2	1.42	1.01
1:K:417:LYS:HG3	1:O:417:LYS:HD2	1.43	1.01
2:L:4:ILE:HG12	2:L:18:VAL:HG12	1.40	1.01
1:C:215:SER:HA	1:C:237:THR:HG21	1.40	1.00
2:H:43:ASP:HB2	2:H:44:PRO:HD3	1.38	1.00
1:I:417:LYS:CG	1:M:417:LYS:HE3	1.90	1.00
1:M:215:SER:HA	1:M:237:THR:HG21	1.44	1.00
2:F:4:ILE:HG12	2:F:18:VAL:HG12	1.42	1.00
1:E:162:ASP:HA	1:E:190:PHE:HZ	1.23	1.00
1:E:416:LYS:HB3	1:E:417:LYS:HE2	1.38	0.99
1:K:417:LYS:HE3	1:O:417:LYS:CG	1.90	0.99
1:A:417:LYS:CG	1:E:417:LYS:HE3	1.92	0.99
1:A:417:LYS:HG3	1:E:417:LYS:HE3	1.00	0.99
2:N:4:ILE:HG12	2:N:18:VAL:HG12	1.40	0.99
1:M:155:THR:HG22	1:M:178:PRO:HD2	1.44	0.99
1:C:417:LYS:HE3	1:G:417:LYS:CG	1.92	0.98
1:C:148:LEU:HA	1:C:427:ARG:NH2	1.77	0.98
1:A:417:LYS:HD2	1:E:417:LYS:HG3	1.45	0.98
1:G:416:LYS:HB3	1:G:417:LYS:HE2	1.44	0.97
1:M:416:LYS:HB3	1:M:417:LYS:HE2	1.44	0.97
1:E:155:THR:HG22	1:E:178:PRO:HD2	1.45	0.97
1:O:215:SER:HA	1:O:237:THR:HG21	1.44	0.97
1:I:416:LYS:HB3	1:I:417:LYS:HE2	1.44	0.97
1:G:215:SER:HA	1:G:237:THR:HG21	1.45	0.96
1:K:417:LYS:HE3	1:O:417:LYS:HG3	0.97	0.96
1:G:176:ARG:CZ	1:G:419:GLN:H	1.79	0.95
1:G:416:LYS:HD2	1:G:417:LYS:HZ1	1.25	0.95
1:C:155:THR:HG22	1:C:178:PRO:HD2	1.48	0.95
1:G:195:VAL:HG11	1:G:198:MET:HE3	1.47	0.95
1:I:417:LYS:HG3	1:M:417:LYS:HE3	0.94	0.94
1:A:155:THR:HG22	1:A:178:PRO:HD2	1.48	0.94
1:I:215:SER:HA	1:I:237:THR:HG21	1.45	0.93
1:I:417:LYS:CD	1:M:417:LYS:HG3	1.97	0.93
1:K:399:GLN:HA	1:K:399:GLN:HE21	1.34	0.93
1:O:195:VAL:HG11	1:O:198:MET:HE3	1.50	0.93
2:D:65:HIS:HB3	2:D:68:ASP:HB2	1.48	0.92
1:K:116:LEU:HD21	2:L:108:ASN:CG	1.95	0.92
1:M:156:GLY:HA2	1:M:179:ARG:O	1.70	0.92
1:C:113:PRO:HD2	1:C:116:LEU:HD12	1.50	0.91
1:I:416:LYS:HD2	1:I:417:LYS:NZ	1.84	0.91
1:K:215:SER:HA	1:K:237:THR:HG21	1.52	0.91

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:253:SER:HB2	1:K:280:ASP:HB2	1.53	0.91
1:K:197:HIS:HD2	1:K:222:ASN:HD22	1.19	0.90
1:K:417:LYS:HG3	1:O:417:LYS:CD	2.00	0.90
1:A:113:PRO:HD2	1:A:116:LEU:HD12	1.53	0.90
1:E:195:VAL:HG11	1:E:198:MET:HE3	1.54	0.90
1:G:186:LEU:C	1:G:187:ALA:CA	2.44	0.90
2:H:65:HIS:HB3	2:H:68:ASP:HB2	1.54	0.89
1:O:176:ARG:CZ	1:O:419:GLN:H	1.84	0.89
1:O:416:LYS:HD2	1:O:417:LYS:NZ	1.87	0.89
1:O:155:THR:HG22	1:O:178:PRO:HD2	1.55	0.89
1:A:156:GLY:HA2	1:A:179:ARG:O	1.72	0.88
1:C:253:SER:HB2	1:C:280:ASP:HB2	1.52	0.88
1:I:289:ALA:O	1:I:293:VAL:HG23	1.73	0.88
1:E:189:HIS:C	1:E:190:PHE:O	2.14	0.88
2:H:107:ALA:HB2	2:H:115:LEU:HD12	1.55	0.88
1:I:155:THR:HG22	1:I:178:PRO:HD2	1.51	0.88
1:G:289:ALA:O	1:G:293:VAL:HG23	1.74	0.88
1:M:253:SER:HB2	1:M:280:ASP:HB2	1.54	0.88
1:I:113:PRO:HD2	1:I:116:LEU:HD12	1.54	0.88
2:D:5:LYS:HG2	2:D:15:GLU:HG2	1.56	0.87
1:A:197:HIS:CD2	1:A:222:ASN:HD22	1.92	0.87
1:G:197:HIS:HD2	1:G:222:ASN:HD22	1.21	0.87
1:I:414:GLY:HA2	1:M:416:LYS:NZ	1.89	0.87
1:O:109:TRP:HA	2:P:101:PHE:HE2	1.38	0.87
2:B:107:ALA:HB2	2:B:115:LEU:HD12	1.54	0.87
1:C:153:ASP:OD1	1:C:155:THR:HG23	1.76	0.86
1:A:195:VAL:HG11	1:A:198:MET:HE3	1.56	0.86
1:I:414:GLY:HA2	1:M:416:LYS:HZ2	1.40	0.86
1:A:137:LYS:HE3	2:B:144:ASP:O	1.75	0.86
2:F:5:LYS:HG2	2:F:15:GLU:HG2	1.58	0.86
1:E:189:HIS:O	1:E:190:PHE:C	2.17	0.86
1:C:253:SER:CB	1:C:280:ASP:HB2	2.05	0.86
2:P:107:ALA:HB2	2:P:115:LEU:HD12	1.55	0.86
1:E:389:GLN:O	1:E:393:GLU:HG2	1.77	0.85
1:O:389:GLN:O	1:O:393:GLU:HG2	1.76	0.85
1:K:113:PRO:HD2	1:K:116:LEU:HD12	1.56	0.85
2:J:65:HIS:HB3	2:J:68:ASP:HB2	1.56	0.85
1:K:195:VAL:HG11	1:K:198:MET:HE3	1.57	0.85
1:O:128:GLU:HG2	2:P:128:LYS:HG2	1.59	0.85
1:C:215:SER:HA	1:C:237:THR:CG2	2.07	0.85
1:E:215:SER:HA	1:E:237:THR:CG2	2.07	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:415:ASN:O	1:I:417:LYS:HG2	1.76	0.84
1:K:155:THR:HG22	1:K:178:PRO:HD2	1.57	0.84
1:K:417:LYS:HG3	1:O:417:LYS:CE	2.07	0.84
1:M:197:HIS:HD2	1:M:222:ASN:HD22	1.22	0.84
1:M:289:ALA:O	1:M:293:VAL:HG23	1.77	0.84
1:A:220:LEU:H	1:A:241:ASN:HD22	1.26	0.84
1:C:289:ALA:O	1:C:293:VAL:HG23	1.77	0.84
1:E:156:GLY:HA2	1:E:179:ARG:O	1.78	0.84
1:G:416:LYS:HD2	1:G:417:LYS:NZ	1.92	0.84
1:O:220:LEU:H	1:O:241:ASN:HD22	1.26	0.84
1:O:415:ASN:O	1:O:417:LYS:HG2	1.78	0.84
1:I:195:VAL:HG11	1:I:198:MET:HE3	1.59	0.84
1:K:416:LYS:CB	1:K:417:LYS:HE2	2.07	0.84
1:E:283:GLU:O	1:E:287:GLN:HG3	1.78	0.83
1:I:156:GLY:HA2	1:I:179:ARG:O	1.78	0.83
1:A:197:HIS:HD2	1:A:222:ASN:HD22	1.26	0.83
1:C:195:VAL:HG11	1:C:198:MET:HE3	1.58	0.83
1:E:416:LYS:CB	1:E:417:LYS:HE2	2.08	0.83
1:I:253:SER:CB	1:I:280:ASP:HB2	2.08	0.83
1:M:195:VAL:HG11	1:M:198:MET:HE3	1.60	0.83
1:A:417:LYS:CD	1:E:417:LYS:HG3	2.09	0.83
1:M:215:SER:HA	1:M:237:THR:CG2	2.08	0.83
1:O:253:SER:HB2	1:O:280:ASP:HB2	1.58	0.83
1:E:253:SER:CB	1:E:280:ASP:HB2	2.09	0.83
1:K:253:SER:CB	1:K:280:ASP:HB2	2.09	0.83
1:A:416:LYS:HD2	1:A:417:LYS:NZ	1.94	0.82
1:A:253:SER:CB	1:A:280:ASP:HB2	2.09	0.82
1:E:253:SER:HB2	1:E:280:ASP:HB2	1.61	0.82
1:K:136:CYS:HB2	2:L:144:ASP:OD2	1.79	0.82
1:I:253:SER:HB2	1:I:280:ASP:HB2	1.61	0.82
1:K:162:ASP:HA	1:K:190:PHE:HZ	1.44	0.82
1:G:113:PRO:HD2	1:G:116:LEU:HD12	1.59	0.82
2:J:5:LYS:HG2	2:J:15:GLU:HG2	1.61	0.82
1:O:148:LEU:HA	1:O:427:ARG:HH22	1.44	0.81
1:A:389:GLN:O	1:A:393:GLU:HG2	1.80	0.81
1:A:415:ASN:O	1:A:417:LYS:HG2	1.80	0.81
1:C:162:ASP:HA	1:C:190:PHE:CZ	2.15	0.81
1:C:417:LYS:HG3	1:G:417:LYS:HD2	1.60	0.81
1:O:289:ALA:O	1:O:293:VAL:HG23	1.79	0.81
1:M:113:PRO:HD2	1:M:116:LEU:HD12	1.61	0.81
1:C:191:SER:HB3	1:C:192:PRO:HD2	1.62	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:283:GLU:O	1:K:287:GLN:HG3	1.80	0.81
1:K:281:PHE:HB2	1:K:285:HIS:CG	2.16	0.81
1:G:148:LEU:HA	1:G:427:ARG:HH22	1.46	0.81
1:G:176:ARG:CD	1:G:419:GLN:HA	2.11	0.81
1:C:416:LYS:CB	1:C:417:LYS:HE2	2.10	0.80
1:C:220:LEU:HB2	1:C:241:ASN:HD22	1.46	0.80
1:M:148:LEU:HA	1:M:427:ARG:HH22	1.46	0.80
1:A:253:SER:HB2	1:A:280:ASP:HB2	1.62	0.80
1:E:148:LEU:HA	1:E:427:ARG:HH22	1.44	0.80
2:N:5:LYS:HG2	2:N:15:GLU:HG2	1.60	0.80
1:G:162:ASP:HA	1:G:190:PHE:HZ	1.45	0.80
1:E:281:PHE:HB2	1:E:285:HIS:ND1	1.96	0.80
1:K:220:LEU:H	1:K:241:ASN:HD22	1.26	0.80
1:I:220:LEU:HB2	1:I:241:ASN:ND2	1.97	0.80
1:M:253:SER:CB	1:M:280:ASP:HB2	2.12	0.80
1:O:162:ASP:HA	1:O:190:PHE:HZ	1.47	0.80
1:K:151:THR:HB	1:K:426:CYS:HA	1.64	0.79
1:E:416:LYS:HD2	1:E:417:LYS:NZ	1.96	0.79
1:M:175:PHE:HD2	1:M:198:MET:HE2	1.47	0.79
1:A:414:GLY:HA2	1:E:416:LYS:NZ	1.97	0.79
1:I:215:SER:HA	1:I:237:THR:CG2	2.12	0.79
1:G:253:SER:CB	1:G:280:ASP:HB2	2.13	0.79
2:B:5:LYS:HG2	2:B:15:GLU:HG2	1.63	0.79
1:K:389:GLN:O	1:K:393:GLU:HG2	1.82	0.79
2:P:5:LYS:HG2	2:P:15:GLU:HG2	1.64	0.79
1:C:389:GLN:O	1:C:393:GLU:HG2	1.82	0.79
1:I:389:GLN:O	1:I:393:GLU:HG2	1.83	0.79
1:K:189:HIS:O	1:K:190:PHE:O	1.60	0.79
1:O:113:PRO:HD2	1:O:116:LEU:HD12	1.64	0.79
1:K:281:PHE:HB2	1:K:285:HIS:ND1	1.98	0.79
1:O:215:SER:HA	1:O:237:THR:CG2	2.13	0.79
1:C:137:LYS:HE3	2:D:144:ASP:O	1.83	0.79
1:E:197:HIS:HD2	1:E:222:ASN:HD22	1.27	0.79
1:O:351:HIS:HD2	1:O:376:THR:OG1	1.66	0.79
1:E:113:PRO:HD2	1:E:116:LEU:HD12	1.65	0.78
2:F:107:ALA:HB2	2:F:115:LEU:HD12	1.66	0.78
1:K:215:SER:HA	1:K:237:THR:CG2	2.12	0.78
1:E:162:ASP:HA	1:E:190:PHE:CZ	2.13	0.78
1:E:281:PHE:HB2	1:E:285:HIS:CG	2.19	0.78
1:G:159:LEU:HD23	1:G:164:THR:HG22	1.65	0.78
1:I:417:LYS:CE	1:M:417:LYS:HG3	2.12	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:415:ASN:O	1:M:417:LYS:HG2	1.84	0.78
1:E:159:LEU:HD23	1:E:164:THR:HG22	1.65	0.78
1:O:131:LYS:HD3	2:P:130:LYS:O	1.83	0.78
1:G:281:PHE:HB2	1:G:285:HIS:CG	2.19	0.78
1:I:351:HIS:HD2	1:I:376:THR:OG1	1.66	0.77
2:H:4:ILE:CG1	2:H:18:VAL:HG12	2.13	0.77
1:G:148:LEU:HA	1:G:427:ARG:NH2	1.99	0.77
1:M:389:GLN:O	1:M:393:GLU:HG2	1.84	0.77
1:A:281:PHE:HB2	1:A:285:HIS:CG	2.19	0.77
1:G:281:PHE:HB2	1:G:285:HIS:ND1	2.00	0.77
2:N:107:ALA:HB2	2:N:115:LEU:HD12	1.64	0.77
1:O:253:SER:CB	1:O:280:ASP:HB2	2.14	0.77
1:G:215:SER:HA	1:G:237:THR:CG2	2.12	0.77
1:K:197:HIS:CD2	1:K:222:ASN:HD22	2.03	0.77
2:L:146:THR:HG22	2:L:148:GLU:H	1.49	0.77
1:O:182:MET:O	1:O:205:ILE:HA	1.84	0.77
1:K:116:LEU:HD21	2:L:108:ASN:OD1	1.85	0.77
1:I:197:HIS:HD2	1:I:222:ASN:HD22	1.32	0.77
1:M:220:LEU:H	1:M:241:ASN:HD22	1.30	0.77
1:G:253:SER:HB2	1:G:280:ASP:HB2	1.67	0.77
1:C:220:LEU:HB2	1:C:241:ASN:ND2	1.99	0.76
2:P:4:ILE:CG1	2:P:18:VAL:HG12	2.15	0.76
1:G:153:ASP:OD1	1:G:155:THR:HG23	1.85	0.76
1:I:197:HIS:CD2	1:I:222:ASN:HD22	2.03	0.76
1:I:220:LEU:HB2	1:I:241:ASN:HD22	1.47	0.76
1:K:416:LYS:NZ	1:O:414:GLY:HA2	2.00	0.76
1:C:415:ASN:O	1:C:417:LYS:HG2	1.85	0.76
1:O:153:ASP:OD1	1:O:155:THR:HG23	1.86	0.76
1:O:281:PHE:HB2	1:O:285:HIS:ND1	1.99	0.76
1:G:156:GLY:HA2	1:G:179:ARG:O	1.85	0.76
1:G:197:HIS:CD2	1:G:222:ASN:HD22	2.04	0.76
1:G:415:ASN:O	1:G:417:LYS:HG2	1.84	0.76
1:E:176:ARG:CZ	1:E:419:GLN:H	1.98	0.76
1:G:131:LYS:HD3	2:H:130:LYS:O	1.85	0.76
1:E:215:SER:CA	1:E:237:THR:HG21	2.14	0.76
1:K:415:ASN:O	1:K:417:LYS:HG2	1.84	0.76
1:M:281:PHE:HB2	1:M:285:HIS:CG	2.21	0.76
1:E:220:LEU:H	1:E:241:ASN:HD22	1.34	0.76
1:O:109:TRP:HZ3	2:P:104:ILE:HD12	1.49	0.76
1:C:281:PHE:HB2	1:C:285:HIS:ND1	2.01	0.76
1:K:155:THR:HG21	1:K:176:ARG:NH1	2.00	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:161:PRO:HB3	1:K:209:THR:HG23	1.65	0.75
1:M:416:LYS:CB	1:M:417:LYS:HE2	2.17	0.75
1:C:416:LYS:NZ	1:G:414:GLY:HA2	2.00	0.75
1:M:148:LEU:HA	1:M:427:ARG:NH2	2.01	0.75
1:G:189:HIS:O	1:G:190:PHE:O	1.68	0.75
1:O:156:GLY:HA2	1:O:179:ARG:O	1.85	0.75
2:H:5:LYS:HG2	2:H:15:GLU:HG2	1.67	0.75
1:I:350:GLN:HA	1:I:374:LEU:HA	1.69	0.75
1:C:389:GLN:NE2	1:C:392:LYS:HD2	2.02	0.75
2:L:43:ASP:HB2	2:L:44:PRO:CD	2.11	0.75
1:M:215:SER:CA	1:M:237:THR:HG21	2.17	0.75
2:N:133:GLU:O	2:N:137:LYS:HG3	1.88	0.74
1:C:197:HIS:HD2	1:C:222:ASN:HD22	1.33	0.74
1:G:220:LEU:HB2	1:G:241:ASN:ND2	2.03	0.74
1:I:281:PHE:HB2	1:I:285:HIS:CG	2.22	0.74
1:M:399:GLN:HE21	1:M:399:GLN:HA	1.51	0.74
1:A:153:ASP:OD1	1:A:155:THR:HG23	1.88	0.74
1:C:175:PHE:HD2	1:C:198:MET:HE2	1.52	0.74
2:P:22:LYS:CG	2:P:28:LYS:HZ3	2.01	0.74
2:B:86:PRO:O	2:B:90:GLN:HG3	1.86	0.74
1:A:215:SER:HA	1:A:237:THR:CG2	2.15	0.74
1:M:159:LEU:HD23	1:M:164:THR:HG22	1.69	0.74
1:O:281:PHE:HB2	1:O:285:HIS:CG	2.23	0.74
1:C:274:ASN:C	1:C:274:ASN:HD22	1.94	0.74
1:C:281:PHE:HB2	1:C:285:HIS:CG	2.22	0.74
1:I:175:PHE:HD2	1:I:198:MET:HE2	1.53	0.74
1:E:289:ALA:O	1:E:293:VAL:HG23	1.88	0.73
1:E:399:GLN:HE21	1:E:399:GLN:HA	1.51	0.73
1:G:176:ARG:HD3	1:G:419:GLN:HA	1.67	0.73
2:H:107:ALA:HB2	2:H:115:LEU:CD1	2.18	0.73
1:A:417:LYS:CE	1:E:417:LYS:HG3	2.19	0.73
1:I:367:GLU:O	1:I:370:GLU:HG2	1.86	0.73
1:M:155:THR:HG21	1:M:176:ARG:NH1	2.02	0.73
2:B:88:TRP:O	2:B:91:GLU:HG3	1.89	0.73
1:E:148:LEU:HA	1:E:427:ARG:NH2	2.03	0.73
1:I:417:LYS:HD2	1:M:417:LYS:CG	2.15	0.73
2:J:49:ASN:HB2	2:J:109:TYR:CE2	2.23	0.73
2:J:88:TRP:O	2:J:91:GLU:HG3	1.87	0.73
1:K:289:ALA:O	1:K:293:VAL:HG23	1.88	0.73
1:O:281:PHE:CE1	1:O:305:TYR:CE2	2.76	0.73
1:A:176:ARG:CZ	1:A:419:GLN:H	2.01	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:148:LEU:HA	1:A:427:ARG:HH22	1.54	0.73
1:C:416:LYS:HD2	1:C:417:LYS:NZ	2.03	0.73
1:C:399:GLN:HA	1:C:399:GLN:HE21	1.53	0.73
1:G:220:LEU:H	1:G:241:ASN:HD22	1.35	0.73
1:G:350:GLN:HA	1:G:374:LEU:HA	1.70	0.73
1:M:191:SER:HB3	1:M:192:PRO:HD2	1.70	0.73
1:A:155:THR:HG21	1:A:176:ARG:NH1	2.04	0.73
1:C:215:SER:CA	1:C:237:THR:HG21	2.17	0.73
1:E:415:ASN:O	1:E:417:LYS:HG2	1.89	0.73
1:C:176:ARG:CZ	1:C:419:GLN:H	2.01	0.72
1:I:176:ARG:CZ	1:I:419:GLN:H	2.01	0.72
1:K:417:LYS:CG	1:O:417:LYS:HD2	2.18	0.72
1:O:367:GLU:O	1:O:370:GLU:HG2	1.89	0.72
1:C:159:LEU:HD23	1:C:164:THR:HG22	1.70	0.72
1:C:351:HIS:HD2	1:C:376:THR:OG1	1.72	0.72
1:E:168:LEU:HD22	1:E:195:VAL:CG2	2.19	0.72
1:C:376:THR:HG22	1:C:399:GLN:HB3	1.70	0.72
2:J:117:ASP:OD1	2:J:121:LYS:HE3	1.89	0.72
2:N:130:LYS:HE3	2:N:138:THR:HG21	1.71	0.72
1:O:137:LYS:HE3	2:P:144:ASP:O	1.88	0.72
1:O:176:ARG:CD	1:O:419:GLN:HA	2.19	0.72
1:M:128:GLU:HG2	2:N:128:LYS:HG2	1.71	0.72
1:K:220:LEU:HB2	1:K:241:ASN:ND2	2.05	0.72
1:E:351:HIS:HD2	1:E:376:THR:OG1	1.72	0.72
1:I:220:LEU:H	1:I:241:ASN:HD22	1.33	0.72
2:D:146:THR:HB	2:D:149:GLU:HG3	1.72	0.72
1:G:389:GLN:O	1:G:393:GLU:HG2	1.88	0.72
1:M:197:HIS:CD2	1:M:222:ASN:HD22	2.05	0.72
1:O:376:THR:HG22	1:O:399:GLN:HB3	1.72	0.72
2:F:49:ASN:HB2	2:F:109:TYR:CE2	2.25	0.72
2:J:26:THR:O	2:J:29:THR:HG22	1.89	0.72
2:B:49:ASN:HB2	2:B:109:TYR:CE2	2.25	0.71
1:E:397:HIS:CE1	1:E:398:LEU:HG	2.25	0.71
2:P:26:THR:O	2:P:29:THR:HG22	1.89	0.71
2:L:5:LYS:HG2	2:L:15:GLU:HG2	1.72	0.71
1:O:161:PRO:HB3	1:O:209:THR:HG23	1.70	0.71
1:O:215:SER:CA	1:O:237:THR:HG21	2.20	0.71
1:O:283:GLU:O	1:O:287:GLN:HG3	1.91	0.71
1:A:356:ARG:HG2	1:A:358:TYR:OH	1.91	0.71
1:A:399:GLN:HA	1:A:399:GLN:HE21	1.54	0.71
1:E:197:HIS:CD2	1:E:222:ASN:HD22	2.08	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:107:ALA:HB2	2:L:115:LEU:HD12	1.72	0.71
1:C:350:GLN:HA	1:C:374:LEU:HA	1.73	0.71
1:E:220:LEU:HB2	1:E:241:ASN:ND2	2.06	0.71
1:M:116:LEU:HD21	2:N:108:ASN:CG	2.15	0.71
1:E:220:LEU:HB2	1:E:241:ASN:HD22	1.56	0.71
1:C:109:TRP:CD2	2:D:101:PHE:HD2	2.09	0.71
1:M:176:ARG:CZ	1:M:419:GLN:H	2.03	0.71
1:O:220:LEU:HB2	1:O:241:ASN:ND2	2.06	0.71
2:N:107:ALA:HB2	2:N:115:LEU:CD1	2.20	0.71
2:J:107:ALA:HB2	2:J:115:LEU:HD12	1.73	0.71
1:O:350:GLN:HA	1:O:374:LEU:HA	1.71	0.71
1:C:197:HIS:CD2	1:C:222:ASN:HD22	2.08	0.71
1:I:181:PHE:CD1	1:K:384:PRO:HG3	2.26	0.71
1:M:416:LYS:HD2	1:M:417:LYS:NZ	2.05	0.71
1:E:376:THR:HG22	1:E:399:GLN:HB3	1.73	0.70
1:M:164:THR:HG21	1:M:182:MET:SD	2.32	0.70
1:I:283:GLU:O	1:I:287:GLN:HG3	1.92	0.70
1:K:164:THR:HG21	1:K:182:MET:SD	2.31	0.70
1:M:176:ARG:HD3	1:M:419:GLN:HA	1.73	0.70
1:O:109:TRP:HA	2:P:101:PHE:CE2	2.24	0.70
2:F:98:GLY:HA3	1:O:284:LYS:NZ	2.06	0.70
1:K:416:LYS:HD2	1:K:417:LYS:NZ	2.06	0.70
1:A:417:LYS:HD2	1:E:417:LYS:CG	2.18	0.70
2:L:54:ILE:HG23	2:L:103:LEU:HD23	1.73	0.70
2:H:130:LYS:HE3	2:H:138:THR:HG21	1.73	0.70
1:M:176:ARG:CD	1:M:419:GLN:HA	2.22	0.70
1:C:164:THR:HG21	1:C:182:MET:SD	2.32	0.70
1:G:367:GLU:O	1:G:370:GLU:HG2	1.92	0.70
1:K:109:TRP:HA	2:L:101:PHE:HE2	1.57	0.70
1:K:156:GLY:HA2	1:K:179:ARG:O	1.92	0.70
1:C:161:PRO:HB3	1:C:209:THR:HG23	1.72	0.70
1:K:367:GLU:O	1:K:370:GLU:HG2	1.90	0.70
1:I:274:ASN:HD22	1:I:274:ASN:C	2.00	0.70
1:A:175:PHE:HD2	1:A:198:MET:HE2	1.56	0.69
1:A:351:HIS:HD2	1:A:376:THR:OG1	1.75	0.69
1:A:389:GLN:NE2	1:A:392:LYS:HD2	2.06	0.69
1:K:215:SER:CA	1:K:237:THR:HG21	2.22	0.69
1:M:389:GLN:NE2	1:M:392:LYS:HD2	2.07	0.69
1:C:417:LYS:HG3	1:G:417:LYS:CE	2.22	0.69
2:D:107:ALA:HB2	2:D:115:LEU:HD12	1.74	0.69
1:G:137:LYS:HE3	2:H:144:ASP:O	1.92	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:281:PHE:HB2	1:M:285:HIS:ND1	2.07	0.69
1:C:168:LEU:HD22	1:C:195:VAL:HG21	1.73	0.69
1:E:168:LEU:HD22	1:E:195:VAL:HG21	1.71	0.69
1:A:281:PHE:CE1	1:A:305:TYR:CE2	2.80	0.69
2:D:130:LYS:HE3	2:D:138:THR:HG21	1.75	0.69
1:G:215:SER:CA	1:G:237:THR:HG21	2.21	0.69
1:O:416:LYS:C	1:O:417:LYS:HD3	2.18	0.69
2:B:130:LYS:HE3	2:B:138:THR:HG21	1.73	0.69
1:C:417:LYS:HG3	1:G:417:LYS:CD	2.22	0.69
1:K:397:HIS:CE1	1:K:398:LEU:HG	2.28	0.69
2:F:43:ASP:HB2	2:F:44:PRO:CD	2.21	0.69
1:I:215:SER:CA	1:I:237:THR:HG21	2.22	0.69
2:L:146:THR:HB	2:L:149:GLU:HG3	1.74	0.68
1:C:367:GLU:O	1:C:370:GLU:HG2	1.92	0.68
1:C:109:TRP:CZ2	2:D:101:PHE:HB2	2.29	0.68
1:O:274:ASN:HD22	1:O:274:ASN:C	2.01	0.68
1:A:295:GLU:HG2	1:A:320:ARG:O	1.94	0.68
1:A:416:LYS:C	1:A:417:LYS:HD3	2.18	0.68
1:E:188:GLU:O	1:E:212:GLY:O	2.12	0.68
1:E:376:THR:HG22	1:E:399:GLN:CB	2.24	0.68
1:K:109:TRP:CD2	2:L:101:PHE:HD2	2.11	0.68
1:A:414:GLY:HA2	1:E:416:LYS:HZ2	1.57	0.68
1:E:350:GLN:HA	1:E:374:LEU:HA	1.76	0.68
1:M:295:GLU:HG2	1:M:320:ARG:O	1.93	0.68
1:O:197:HIS:HD2	1:O:222:ASN:HD22	1.42	0.68
1:A:148:LEU:HA	1:A:427:ARG:NH2	2.09	0.68
1:G:416:LYS:C	1:G:417:LYS:HD3	2.19	0.68
2:H:146:THR:HG22	2:H:148:GLU:H	1.59	0.68
1:I:281:PHE:HB2	1:I:285:HIS:ND1	2.09	0.68
1:C:281:PHE:CE1	1:C:305:TYR:CE2	2.81	0.68
1:E:416:LYS:HD2	1:E:417:LYS:HZ1	1.57	0.68
1:I:295:GLU:HG2	1:I:320:ARG:O	1.93	0.68
2:L:86:PRO:O	2:L:90:GLN:HG3	1.94	0.68
1:O:164:THR:HG21	1:O:182:MET:SD	2.33	0.68
1:G:186:LEU:C	1:G:187:ALA:C	2.62	0.68
1:G:383:VAL:HG12	1:G:388:LEU:HB2	1.76	0.68
1:M:331:ASP:OD2	1:M:356:ARG:NH1	2.27	0.68
1:A:378:GLN:NE2	1:A:401:ASN:HA	2.09	0.67
1:K:140:TYR:CE2	2:L:153:VAL:HG22	2.29	0.67
1:G:155:THR:HG21	1:G:176:ARG:NH1	2.09	0.67
1:M:175:PHE:CD2	1:M:198:MET:HE2	2.30	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:98:GLY:HA3	1:O:284:LYS:HZ1	1.57	0.67
1:G:416:LYS:O	1:G:417:LYS:HD3	1.94	0.67
1:K:314:LEU:O	1:K:318:VAL:HG23	1.94	0.67
1:A:176:ARG:HD3	1:A:419:GLN:HA	1.76	0.67
1:E:367:GLU:O	1:E:370:GLU:HG2	1.95	0.67
1:M:367:GLU:O	1:M:370:GLU:HG2	1.94	0.67
2:N:86:PRO:O	2:N:90:GLN:HG3	1.95	0.67
1:C:415:ASN:O	1:C:416:LYS:HB3	1.95	0.67
1:E:231:SER:OG	1:E:233:PRO:HD2	1.94	0.67
1:K:351:HIS:HD2	1:K:376:THR:OG1	1.76	0.67
2:H:88:TRP:O	2:H:91:GLU:HG3	1.93	0.67
2:H:133:GLU:O	2:H:137:LYS:HG3	1.94	0.67
1:E:153:ASP:OD1	1:E:155:THR:HG23	1.95	0.67
1:I:389:GLN:NE2	1:I:392:LYS:HD2	2.09	0.67
2:L:130:LYS:HE3	2:L:138:THR:HG21	1.76	0.67
1:O:193:PHE:HB3	1:O:195:VAL:HG23	1.76	0.67
1:O:399:GLN:HA	1:O:399:GLN:HE21	1.59	0.67
2:P:43:ASP:HB2	2:P:44:PRO:CD	2.22	0.67
1:A:176:ARG:CD	1:A:419:GLN:HA	2.25	0.67
1:G:220:LEU:HB2	1:G:241:ASN:HD22	1.60	0.67
1:G:384:PRO:HD2	1:G:387:THR:OG1	1.95	0.67
1:G:397:HIS:CE1	1:G:398:LEU:HG	2.30	0.67
2:P:143:ASN:OD1	2:P:145:PHE:HB2	1.95	0.66
2:J:130:LYS:HE3	2:J:138:THR:HG21	1.76	0.66
1:M:376:THR:HG22	1:M:399:GLN:HB3	1.76	0.66
2:P:117:ASP:OD1	2:P:121:LYS:HE3	1.95	0.66
1:A:344:PHE:CE2	1:A:370:GLU:HB2	2.31	0.66
1:E:137:LYS:HE3	2:F:145:PHE:HA	1.76	0.66
1:M:153:ASP:OD1	1:M:155:THR:HG23	1.95	0.66
1:O:415:ASN:O	1:O:416:LYS:HB3	1.95	0.66
1:C:148:LEU:CA	1:C:427:ARG:HH22	2.04	0.66
1:G:168:LEU:HD22	1:G:195:VAL:HG21	1.76	0.66
1:O:376:THR:HG22	1:O:399:GLN:CB	2.24	0.66
1:C:295:GLU:HG2	1:C:320:ARG:O	1.95	0.66
2:H:43:ASP:CB	2:H:44:PRO:HD3	2.21	0.66
1:O:175:PHE:HD2	1:O:198:MET:HE2	1.61	0.66
1:E:215:SER:O	1:E:240:LYS:NZ	2.29	0.66
1:K:415:ASN:O	1:K:416:LYS:HB3	1.95	0.66
2:F:4:ILE:CG1	2:F:18:VAL:HG12	2.22	0.66
1:K:350:GLN:HA	1:K:374:LEU:HA	1.76	0.66
1:O:197:HIS:CD2	1:O:222:ASN:HD22	2.13	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:220:LEU:HB2	1:A:241:ASN:ND2	2.11	0.66
2:B:107:ALA:HB2	2:B:115:LEU:CD1	2.25	0.66
1:O:344:PHE:CE2	1:O:370:GLU:HB2	2.31	0.66
1:A:162:ASP:HA	1:A:190:PHE:HZ	1.60	0.66
1:O:137:LYS:HG3	2:P:144:ASP:HB2	1.77	0.66
1:A:159:LEU:HD23	1:A:164:THR:HG22	1.77	0.66
1:A:383:VAL:HG12	1:A:388:LEU:HB2	1.77	0.66
1:C:376:THR:HG22	1:C:399:GLN:CB	2.25	0.66
1:A:225:LEU:O	1:A:252:CYS:SG	2.54	0.65
1:K:158:ASN:HA	1:K:181:PHE:HB2	1.77	0.65
1:A:281:PHE:HB2	1:A:285:HIS:ND1	2.12	0.65
1:G:351:HIS:HD2	1:G:376:THR:OG1	1.79	0.65
1:M:182:MET:O	1:M:205:ILE:HA	1.95	0.65
2:N:104:ILE:HG12	2:N:119:THR:HB	1.79	0.65
1:O:158:ASN:HA	1:O:181:PHE:HB2	1.78	0.65
1:O:397:HIS:CE1	1:O:398:LEU:HG	2.30	0.65
2:B:43:ASP:HB2	2:B:44:PRO:CD	2.23	0.65
1:M:161:PRO:HB3	1:M:209:THR:HG23	1.78	0.65
1:E:175:PHE:HD2	1:E:198:MET:HE2	1.60	0.65
1:O:148:LEU:HA	1:O:427:ARG:NH2	2.10	0.65
1:O:176:ARG:HD3	1:O:419:GLN:HA	1.77	0.65
2:L:101:PHE:CE1	2:L:105:LEU:HD11	2.31	0.65
2:B:146:THR:HB	2:B:149:GLU:HG3	1.78	0.65
2:F:26:THR:O	2:F:29:THR:HG22	1.97	0.65
1:G:274:ASN:HD22	1:G:274:ASN:C	2.05	0.65
1:I:415:ASN:O	1:I:416:LYS:HB3	1.95	0.65
1:A:367:GLU:O	1:A:370:GLU:HG2	1.97	0.65
2:F:107:ALA:HB2	2:F:115:LEU:CD1	2.26	0.65
1:I:416:LYS:C	1:I:417:LYS:HD3	2.22	0.65
1:M:350:GLN:HA	1:M:374:LEU:HA	1.77	0.65
1:K:176:ARG:CZ	1:K:419:GLN:H	2.10	0.65
1:K:220:LEU:H	1:K:241:ASN:ND2	1.95	0.65
1:A:137:LYS:HG3	2:B:144:ASP:HB2	1.78	0.65
1:E:295:GLU:HG2	1:E:320:ARG:O	1.96	0.65
1:M:155:THR:CG2	1:M:178:PRO:HD2	2.25	0.65
1:A:215:SER:CA	1:A:237:THR:HG21	2.21	0.64
1:K:182:MET:O	1:K:205:ILE:HA	1.97	0.64
1:C:253:SER:HB2	1:C:280:ASP:CB	2.25	0.64
1:I:159:LEU:HD23	1:I:164:THR:HG22	1.78	0.64
1:I:362:PRO:HG3	1:I:382:ILE:O	1.97	0.64
1:O:220:LEU:HB2	1:O:241:ASN:HD22	1.60	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:350:GLN:HA	1:A:374:LEU:HA	1.79	0.64
1:C:131:LYS:NZ	2:D:127:ILE:O	2.29	0.64
1:G:186:LEU:HA	1:G:187:ALA:N	2.13	0.64
1:E:383:VAL:HG12	1:E:388:LEU:HB2	1.78	0.64
1:G:419:GLN:NE2	1:G:426:CYS:O	2.31	0.64
1:I:220:LEU:CB	1:I:241:ASN:HD22	2.11	0.64
1:K:253:SER:HB2	1:K:280:ASP:CB	2.26	0.64
1:K:344:PHE:CE2	1:K:370:GLU:HB2	2.33	0.64
1:K:356:ARG:HG2	1:K:358:TYR:OH	1.98	0.64
2:N:43:ASP:CB	2:N:44:PRO:HD3	2.22	0.64
1:K:109:TRP:CE2	2:L:101:PHE:HD2	2.16	0.64
2:P:133:GLU:O	2:P:137:LYS:HG3	1.98	0.64
2:J:101:PHE:CE1	2:J:105:LEU:HD11	2.33	0.64
1:A:397:HIS:CE1	1:A:398:LEU:HG	2.33	0.64
1:E:389:GLN:NE2	1:E:392:LYS:HD2	2.13	0.64
1:I:148:LEU:HA	1:I:427:ARG:HH22	1.62	0.64
2:N:43:ASP:HB2	2:N:44:PRO:CD	2.22	0.64
2:P:61:TRP:CD2	2:P:115:LEU:HD23	2.33	0.63
1:C:186:LEU:O	1:C:188:GLU:N	2.30	0.63
1:I:414:GLY:CA	1:M:416:LYS:NZ	2.60	0.63
1:O:159:LEU:HD23	1:O:164:THR:HG22	1.80	0.63
2:P:61:TRP:CG	2:P:115:LEU:HD23	2.32	0.63
2:P:107:ALA:HB2	2:P:115:LEU:CD1	2.28	0.63
2:B:16:VAL:CG2	2:B:21:ALA:HB2	2.29	0.63
1:C:416:LYS:HD2	1:C:417:LYS:HZ1	1.63	0.63
2:D:146:THR:HG22	2:D:148:GLU:H	1.61	0.63
2:F:133:GLU:O	2:F:137:LYS:HG3	1.99	0.63
1:C:187:ALA:C	1:C:212:GLY:HA3	2.23	0.63
2:D:117:ASP:OD1	2:D:121:LYS:HE3	1.99	0.63
1:E:176:ARG:CD	1:E:419:GLN:HA	2.29	0.63
1:K:220:LEU:HB2	1:K:241:ASN:HD22	1.63	0.63
1:M:151:THR:HB	1:M:426:CYS:HA	1.79	0.63
1:A:415:ASN:O	1:A:416:LYS:HB3	1.98	0.63
1:K:162:ASP:HA	1:K:190:PHE:CZ	2.30	0.63
2:L:51:ASN:ND2	2:L:53:ALA:H	1.95	0.63
1:G:128:GLU:HG2	2:H:128:LYS:HG2	1.80	0.63
1:G:415:ASN:O	1:G:416:LYS:HB3	1.97	0.63
1:K:384:PRO:HD2	1:K:387:THR:OG1	1.98	0.63
1:O:295:GLU:HG2	1:O:320:ARG:O	1.97	0.63
2:J:104:ILE:HG12	2:J:119:THR:HB	1.79	0.63
1:A:231:SER:HA	1:A:254:GLY:O	1.99	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:362:PRO:HG3	1:C:382:ILE:O	1.99	0.63
1:E:415:ASN:O	1:E:416:LYS:HB3	1.99	0.63
2:D:43:ASP:HB2	2:D:44:PRO:CD	2.24	0.62
1:O:220:LEU:H	1:O:241:ASN:ND2	1.96	0.62
1:C:188:GLU:O	1:C:212:GLY:C	2.42	0.62
1:C:417:LYS:CG	1:G:417:LYS:HD2	2.29	0.62
1:I:155:THR:HG21	1:I:176:ARG:NH1	2.13	0.62
1:M:377:LEU:O	1:M:400:ILE:HA	1.99	0.62
2:P:61:TRP:CE3	2:P:115:LEU:HB2	2.34	0.62
1:C:128:GLU:HG2	2:D:128:LYS:HG2	1.81	0.62
1:C:156:GLY:HA2	1:C:179:ARG:O	2.00	0.62
1:E:176:ARG:HD3	1:E:419:GLN:HA	1.81	0.62
1:G:168:LEU:HD22	1:G:195:VAL:CG2	2.28	0.62
2:P:43:ASP:CB	2:P:44:PRO:HD3	2.23	0.62
1:A:253:SER:HB2	1:A:280:ASP:CB	2.29	0.62
2:B:26:THR:O	2:B:29:THR:HG22	1.99	0.62
1:O:362:PRO:HG3	1:O:382:ILE:O	1.99	0.62
1:A:184:GLN:CB	1:A:185:PRO:HD2	2.29	0.62
1:E:155:THR:HG21	1:E:176:ARG:NH1	2.14	0.62
1:I:376:THR:HG22	1:I:399:GLN:HB3	1.81	0.62
1:O:257:GLU:HB2	1:O:288:VAL:HG21	1.81	0.62
1:O:325:VAL:HG13	1:O:350:GLN:HG2	1.81	0.62
1:O:411:PRO:HG3	1:O:422:TRP:CH2	2.34	0.62
2:P:22:LYS:HG2	2:P:28:LYS:HZ3	1.62	0.62
1:C:378:GLN:NE2	1:C:401:ASN:HA	2.14	0.62
1:M:220:LEU:HB2	1:M:241:ASN:ND2	2.15	0.62
1:C:220:LEU:H	1:C:241:ASN:HD22	1.48	0.62
1:I:162:ASP:HA	1:I:190:PHE:HZ	1.65	0.62
1:K:176:ARG:CD	1:K:419:GLN:HA	2.30	0.62
2:L:16:VAL:CG2	2:L:21:ALA:HB2	2.30	0.62
1:I:253:SER:HB2	1:I:280:ASP:CB	2.30	0.62
1:M:397:HIS:CE1	1:M:398:LEU:HG	2.35	0.62
1:A:419:GLN:O	1:A:419:GLN:HG2	2.00	0.62
1:E:158:ASN:HA	1:E:181:PHE:HB2	1.81	0.62
1:I:376:THR:HG22	1:I:399:GLN:CB	2.30	0.62
2:J:4:ILE:CG1	2:J:18:VAL:HG12	2.23	0.62
1:O:155:THR:CG2	1:O:178:PRO:HD2	2.29	0.62
1:I:309:LEU:HD23	1:I:332:SER:OG	1.99	0.61
2:J:49:ASN:HB2	2:J:109:TYR:CZ	2.35	0.61
1:K:168:LEU:HD22	1:K:195:VAL:HG21	1.82	0.61
1:O:155:THR:HG21	1:O:176:ARG:NH1	2.14	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:158:ASN:HA	1:C:181:PHE:HB2	1.82	0.61
1:C:411:PRO:HG3	1:C:422:TRP:CH2	2.35	0.61
2:B:4:ILE:CG1	2:B:18:VAL:HG12	2.24	0.61
1:C:137:LYS:H	2:D:144:ASP:HB2	1.65	0.61
2:J:22:LYS:O	2:J:22:LYS:HD3	2.01	0.61
1:M:356:ARG:HG2	1:M:358:TYR:OH	1.99	0.61
1:O:253:SER:HB2	1:O:280:ASP:CB	2.30	0.61
1:E:253:SER:HB2	1:E:280:ASP:CB	2.29	0.61
2:H:26:THR:O	2:H:29:THR:HG22	2.00	0.61
2:H:51:ASN:ND2	2:H:53:ALA:H	1.98	0.61
1:I:153:ASP:OD1	1:I:155:THR:HG23	2.00	0.61
2:J:86:PRO:O	2:J:90:GLN:HG3	1.99	0.61
1:E:362:PRO:HG3	1:E:382:ILE:O	2.00	0.61
2:F:43:ASP:CB	2:F:44:PRO:HD3	2.23	0.61
2:D:4:ILE:CG1	2:D:18:VAL:HG12	2.22	0.61
2:F:54:ILE:HG23	2:F:103:LEU:HD23	1.83	0.61
2:J:107:ALA:HB2	2:J:115:LEU:CD1	2.31	0.61
2:J:127:ILE:HG23	2:J:135:ILE:CD1	2.30	0.61
2:L:4:ILE:CG1	2:L:18:VAL:HG12	2.23	0.61
1:C:283:GLU:O	1:C:287:GLN:HG3	2.01	0.61
1:E:128:GLU:HG2	2:F:128:LYS:HG2	1.83	0.61
2:J:43:ASP:CB	2:J:44:PRO:HD3	2.22	0.61
1:K:191:SER:HB3	1:K:192:PRO:HD2	1.82	0.61
1:M:314:LEU:O	1:M:318:VAL:HG23	2.01	0.61
1:O:416:LYS:CB	1:O:417:LYS:HE2	2.24	0.61
2:H:43:ASP:HB2	2:H:44:PRO:CD	2.21	0.61
1:I:158:ASN:HA	1:I:181:PHE:HB2	1.81	0.61
2:L:107:ALA:HB2	2:L:115:LEU:CD1	2.31	0.61
1:A:116:LEU:HD13	2:B:104:ILE:HG22	1.81	0.61
1:A:164:THR:HG21	1:A:182:MET:SD	2.40	0.61
1:A:168:LEU:HD22	1:A:195:VAL:HG21	1.83	0.61
1:M:253:SER:HB2	1:M:280:ASP:CB	2.27	0.61
1:M:415:ASN:O	1:M:416:LYS:HB3	2.00	0.61
2:N:16:VAL:CG2	2:N:21:ALA:HB2	2.31	0.61
2:N:146:THR:HG22	2:N:148:GLU:H	1.66	0.61
2:B:101:PHE:CE1	2:B:105:LEU:HD11	2.36	0.61
2:D:16:VAL:CG2	2:D:21:ALA:HB2	2.31	0.61
1:K:109:TRP:CZ2	2:L:101:PHE:HB2	2.36	0.61
2:F:16:VAL:CG2	2:F:21:ALA:HB2	2.31	0.60
1:G:344:PHE:CE2	1:G:370:GLU:HB2	2.36	0.60
1:I:231:SER:HA	1:I:254:GLY:O	2.01	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:O:168:LEU:HD22	1:O:195:VAL:HG21	1.83	0.60
1:C:175:PHE:CD2	1:C:198:MET:HE2	2.35	0.60
1:K:399:GLN:HE21	1:K:399:GLN:CA	2.11	0.60
1:C:281:PHE:CB	1:C:285:HIS:CG	2.84	0.60
1:K:281:PHE:CB	1:K:285:HIS:CG	2.84	0.60
2:N:12:GLU:OE2	2:N:52:ALA:HB1	2.02	0.60
1:A:382:ILE:HG13	1:A:383:VAL:HG23	1.83	0.60
1:I:399:GLN:HA	1:I:399:GLN:HE21	1.66	0.60
1:A:124:LEU:HB3	1:A:128:GLU:HB3	1.83	0.60
1:A:416:LYS:O	1:A:417:LYS:HD3	2.02	0.60
1:G:416:LYS:CB	1:G:417:LYS:HE2	2.27	0.60
2:J:16:VAL:CG2	2:J:21:ALA:HB2	2.30	0.60
1:M:383:VAL:HG12	1:M:388:LEU:HB2	1.82	0.60
1:O:109:TRP:CZ3	2:P:104:ILE:HD12	2.33	0.60
1:G:295:GLU:HG2	1:G:320:ARG:O	2.02	0.60
1:K:276:SER:HB3	1:K:301:ASN:OD1	2.01	0.60
1:K:325:VAL:HG13	1:K:350:GLN:HG2	1.84	0.60
1:M:351:HIS:HD2	1:M:376:THR:OG1	1.84	0.60
1:O:389:GLN:NE2	1:O:392:LYS:HD2	2.17	0.60
1:A:281:PHE:CB	1:A:285:HIS:CG	2.85	0.60
1:C:344:PHE:CE2	1:C:370:GLU:HB2	2.36	0.60
1:C:417:LYS:HG3	1:G:417:LYS:HZ2	1.67	0.60
1:I:281:PHE:CB	1:I:285:HIS:CG	2.85	0.60
2:J:16:VAL:HG21	2:J:21:ALA:HB2	1.84	0.60
1:C:215:SER:O	1:C:240:LYS:NZ	2.35	0.60
1:E:325:VAL:HG13	1:E:350:GLN:HG2	1.82	0.60
2:H:49:ASN:HB2	2:H:109:TYR:CE2	2.37	0.60
1:M:137:LYS:HE3	2:N:145:PHE:HA	1.84	0.60
1:M:159:LEU:CD2	1:M:164:THR:HG22	2.32	0.60
1:M:181:PHE:CD1	1:O:384:PRO:HG3	2.37	0.60
2:D:16:VAL:HG21	2:D:21:ALA:HB2	1.84	0.60
1:E:274:ASN:C	1:E:274:ASN:HD22	2.10	0.59
1:O:132:VAL:CG1	2:P:127:ILE:HG21	2.32	0.59
1:C:168:LEU:HD22	1:C:195:VAL:CG2	2.32	0.59
2:H:101:PHE:CE1	2:H:105:LEU:HD11	2.37	0.59
1:K:175:PHE:HD2	1:K:198:MET:HE2	1.66	0.59
1:O:419:GLN:OE1	1:O:429:THR:HG21	2.02	0.59
2:P:54:ILE:HG23	2:P:103:LEU:HD23	1.83	0.59
1:A:175:PHE:CD2	1:A:198:MET:HE2	2.36	0.59
1:E:281:PHE:CB	1:E:285:HIS:CG	2.85	0.59
1:G:175:PHE:HD2	1:G:198:MET:HE2	1.67	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:281:PHE:CB	1:G:285:HIS:CG	2.85	0.59
2:H:146:THR:HB	2:H:149:GLU:HG3	1.83	0.59
1:I:149:TRP:O	1:I:172:VAL:HA	2.02	0.59
1:I:181:PHE:CE1	1:K:384:PRO:HA	2.37	0.59
2:P:101:PHE:CE1	2:P:105:LEU:HD11	2.37	0.59
2:B:127:ILE:HG23	2:B:135:ILE:CD1	2.32	0.59
1:C:151:THR:HB	1:C:426:CYS:HA	1.84	0.59
2:D:43:ASP:CB	2:D:44:PRO:HD3	2.25	0.59
1:E:115:GLU:OE1	1:E:115:GLU:N	2.36	0.59
2:F:88:TRP:O	2:F:91:GLU:HG3	2.01	0.59
1:K:168:LEU:HD22	1:K:195:VAL:CG2	2.32	0.59
1:K:281:PHE:CE1	1:K:305:TYR:CE2	2.91	0.59
2:B:43:ASP:CB	2:B:44:PRO:HD3	2.25	0.59
1:C:176:ARG:CD	1:C:419:GLN:HA	2.32	0.59
1:C:383:VAL:HG12	1:C:388:LEU:HB2	1.85	0.59
2:F:146:THR:HG22	2:F:148:GLU:H	1.67	0.59
1:G:162:ASP:HA	1:G:190:PHE:CZ	2.31	0.59
1:I:184:GLN:CB	1:I:185:PRO:HD2	2.32	0.59
2:J:115:LEU:O	2:J:115:LEU:HD13	2.01	0.59
2:L:127:ILE:HG23	2:L:135:ILE:CD1	2.33	0.59
2:P:16:VAL:CG1	2:P:59:ILE:HD13	2.33	0.59
2:P:130:LYS:HE3	2:P:138:THR:HG21	1.85	0.59
1:C:187:ALA:O	1:C:188:GLU:C	2.46	0.59
2:D:107:ALA:HB2	2:D:115:LEU:CD1	2.32	0.59
1:G:399:GLN:HA	1:G:399:GLN:HE21	1.68	0.59
1:K:416:LYS:HZ3	1:O:414:GLY:HA2	1.68	0.59
2:L:26:THR:O	2:L:29:THR:HG22	2.02	0.59
2:H:117:ASP:OD1	2:H:121:LYS:HE3	2.03	0.59
1:I:225:LEU:O	1:I:252:CYS:SG	2.61	0.59
1:A:184:GLN:HB3	1:A:185:PRO:HD2	1.84	0.59
1:C:416:LYS:HZ2	1:G:414:GLY:HA2	1.68	0.59
1:G:325:VAL:HG13	1:G:350:GLN:HG2	1.84	0.59
1:G:376:THR:HG22	1:G:399:GLN:CB	2.33	0.59
1:E:419:GLN:OE1	1:E:429:THR:HG21	2.02	0.59
1:G:153:ASP:HB3	1:G:429:THR:HG22	1.85	0.59
1:K:417:LYS:CE	1:O:417:LYS:CG	2.65	0.59
2:N:149:GLU:O	2:N:153:VAL:HG23	2.02	0.59
2:P:16:VAL:CG2	2:P:21:ALA:HB2	2.33	0.59
2:B:87:VAL:HG21	1:E:296:THR:HA	1.83	0.58
1:C:136:CYS:HB2	2:D:144:ASP:OD2	2.03	0.58
1:C:419:GLN:NE2	1:C:426:CYS:O	2.36	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:220:LEU:H	1:I:241:ASN:ND2	2.00	0.58
1:I:419:GLN:NE2	1:I:426:CYS:O	2.36	0.58
2:L:43:ASP:CB	2:L:44:PRO:HD3	2.20	0.58
2:N:26:THR:O	2:N:29:THR:HG22	2.02	0.58
1:A:374:LEU:HG	1:A:398:LEU:HD21	1.86	0.58
2:H:148:GLU:O	2:H:151:ALA:HB3	2.04	0.58
1:O:383:VAL:HG12	1:O:388:LEU:HB2	1.85	0.58
1:E:137:LYS:HG2	2:F:145:PHE:CE2	2.39	0.58
2:P:146:THR:HB	2:P:149:GLU:HG3	1.84	0.58
2:J:43:ASP:HB2	2:J:44:PRO:CD	2.24	0.58
1:K:416:LYS:HZ2	1:O:414:GLY:HA2	1.68	0.58
1:K:424:ILE:HD12	1:K:424:ILE:N	2.18	0.58
1:O:377:LEU:O	1:O:400:ILE:HA	2.04	0.58
1:E:155:THR:CG2	1:E:178:PRO:HD2	2.26	0.58
1:K:383:VAL:HG12	1:K:388:LEU:HB2	1.86	0.58
1:K:419:GLN:OE1	1:K:429:THR:HG21	2.04	0.58
1:A:220:LEU:N	1:A:241:ASN:HD22	2.01	0.58
1:I:149:TRP:O	1:I:172:VAL:HG22	2.03	0.58
1:K:151:THR:CB	1:K:426:CYS:HA	2.32	0.58
1:K:331:ASP:OD2	1:K:356:ARG:NH1	2.37	0.58
1:K:411:PRO:HG3	1:K:422:TRP:CH2	2.39	0.58
2:P:33:ASP:O	2:P:34:LEU:HD23	2.02	0.58
1:A:120:ILE:HD13	2:B:123:VAL:HG11	1.85	0.58
1:A:237:THR:HG23	1:A:240:LYS:HE3	1.86	0.58
1:C:176:ARG:HD3	1:C:419:GLN:HA	1.85	0.58
2:J:146:THR:HB	2:J:149:GLU:HG3	1.85	0.58
1:A:325:VAL:HG13	1:A:350:GLN:HG2	1.85	0.58
2:D:88:TRP:O	2:D:91:GLU:HG3	2.04	0.58
2:D:133:GLU:O	2:D:137:LYS:HG3	2.04	0.58
1:M:124:LEU:HB3	1:M:128:GLU:HB3	1.86	0.58
1:M:419:GLN:OE1	1:M:429:THR:HG21	2.03	0.58
2:N:4:ILE:CG1	2:N:18:VAL:HG12	2.25	0.58
2:P:88:TRP:O	2:P:91:GLU:HG3	2.02	0.58
1:E:220:LEU:H	1:E:241:ASN:ND2	2.01	0.58
1:E:309:LEU:HD23	1:E:332:SER:OG	2.04	0.58
1:G:389:GLN:NE2	1:G:392:LYS:HD2	2.18	0.58
1:I:155:THR:CG2	1:I:178:PRO:HD2	2.28	0.58
1:K:378:GLN:NE2	1:K:401:ASN:HA	2.18	0.58
2:P:127:ILE:HG23	2:P:135:ILE:CD1	2.34	0.58
1:C:184:GLN:CB	1:C:185:PRO:HD2	2.34	0.58
1:E:153:ASP:HB3	1:E:429:THR:HG22	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:383:VAL:HG12	1:I:388:LEU:HB2	1.85	0.58
1:K:225:LEU:O	1:K:252:CYS:SG	2.62	0.58
1:K:362:PRO:HG3	1:K:382:ILE:O	2.04	0.58
1:M:225:LEU:O	1:M:252:CYS:SG	2.62	0.58
1:A:115:GLU:N	1:A:115:GLU:OE1	2.37	0.57
1:A:283:GLU:O	1:A:287:GLN:HG3	2.04	0.57
1:E:184:GLN:CB	1:E:185:PRO:HD2	2.34	0.57
1:E:418:ASN:ND2	1:E:418:ASN:C	2.61	0.57
2:F:86:PRO:O	2:F:90:GLN:HG3	2.03	0.57
1:I:389:GLN:HE22	1:I:392:LYS:NZ	2.02	0.57
1:M:168:LEU:HD22	1:M:195:VAL:HG21	1.85	0.57
1:C:155:THR:CG2	1:C:178:PRO:HD2	2.31	0.57
2:D:51:ASN:ND2	2:D:53:ALA:H	2.03	0.57
1:M:300:LEU:HD23	1:M:317:LEU:HD11	1.85	0.57
1:O:419:GLN:NE2	1:O:426:CYS:O	2.37	0.57
2:B:127:ILE:HG23	2:B:135:ILE:HD13	1.86	0.57
1:K:419:GLN:NE2	1:K:426:CYS:O	2.38	0.57
1:A:411:PRO:HG3	1:A:422:TRP:CH2	2.40	0.57
2:B:64:HIS:CE1	1:E:269:ARG:NH2	2.73	0.57
1:K:116:LEU:HD11	2:L:108:ASN:ND2	2.19	0.57
1:K:274:ASN:C	1:K:274:ASN:HD22	2.11	0.57
1:M:418:ASN:C	1:M:418:ASN:HD22	2.11	0.57
1:E:281:PHE:CE1	1:E:305:TYR:CE2	2.92	0.57
1:I:416:LYS:CD	1:I:417:LYS:HZ1	2.01	0.57
1:O:309:LEU:HD21	1:O:329:LEU:HD22	1.86	0.57
1:E:231:SER:HA	1:E:254:GLY:O	2.05	0.57
1:I:176:ARG:HD3	1:I:419:GLN:HA	1.85	0.57
1:M:384:PRO:HD2	1:M:387:THR:OG1	2.05	0.57
2:N:89:ASP:O	2:N:92:PHE:HB3	2.04	0.57
1:O:123:CYS:SG	2:P:121:LYS:HG2	2.45	0.57
1:A:161:PRO:HB3	1:A:209:THR:HG23	1.86	0.57
2:F:146:THR:HB	2:F:149:GLU:HG3	1.86	0.57
1:G:161:PRO:HB3	1:G:209:THR:HG23	1.86	0.57
1:I:148:LEU:HA	1:I:427:ARG:NH2	2.20	0.57
2:J:133:GLU:O	2:J:137:LYS:HG3	2.05	0.57
2:N:65:HIS:CB	2:N:68:ASP:HB2	2.21	0.57
2:P:101:PHE:HE1	2:P:105:LEU:HD11	1.70	0.57
2:B:49:ASN:HB2	2:B:109:TYR:CZ	2.39	0.57
2:J:101:PHE:HE1	2:J:105:LEU:HD11	1.67	0.57
2:L:104:ILE:HG12	2:L:119:THR:HB	1.86	0.57
1:G:157:LYS:HD2	1:G:430:LEU:HD12	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:378:GLN:NE2	1:I:401:ASN:HA	2.20	0.57
1:M:151:THR:HG22	1:M:174:ALA:HB3	1.86	0.57
1:O:161:PRO:HB3	1:O:209:THR:CG2	2.34	0.57
1:A:220:LEU:H	1:A:241:ASN:ND2	2.01	0.56
1:A:220:LEU:HB2	1:A:241:ASN:HD22	1.69	0.56
1:I:220:LEU:N	1:I:241:ASN:HD22	2.03	0.56
1:K:215:SER:O	1:K:240:LYS:NZ	2.38	0.56
1:A:274:ASN:HD22	1:A:274:ASN:C	2.13	0.56
1:A:419:GLN:OE1	1:A:429:THR:HG21	2.05	0.56
1:C:182:MET:O	1:C:205:ILE:HA	2.05	0.56
1:C:309:LEU:HD21	1:C:329:LEU:HD22	1.87	0.56
1:G:378:GLN:NE2	1:G:401:ASN:HA	2.20	0.56
1:K:159:LEU:HD23	1:K:164:THR:HG22	1.87	0.56
1:O:220:LEU:N	1:O:241:ASN:HD22	1.99	0.56
1:A:181:PHE:CD1	1:C:384:PRO:HG3	2.40	0.56
1:A:370:GLU:O	1:A:372:PRO:HD3	2.05	0.56
1:C:416:LYS:NZ	1:G:414:GLY:CA	2.68	0.56
1:E:418:ASN:C	1:E:418:ASN:HD22	2.12	0.56
2:F:16:VAL:HG21	2:F:21:ALA:HB2	1.88	0.56
1:G:199:ASP:OD1	1:G:411:PRO:HG2	2.06	0.56
1:G:376:THR:HG22	1:G:399:GLN:HB3	1.86	0.56
2:J:146:THR:HG22	2:J:148:GLU:H	1.70	0.56
1:K:320:ARG:HB2	1:K:320:ARG:HH11	1.71	0.56
1:M:418:ASN:C	1:M:418:ASN:ND2	2.62	0.56
1:M:421:ILE:HG23	1:M:424:ILE:HB	1.87	0.56
1:O:124:LEU:HB3	1:O:128:GLU:HB3	1.86	0.56
1:O:378:GLN:NE2	1:O:401:ASN:HA	2.19	0.56
2:P:146:THR:HG22	2:P:148:GLU:H	1.70	0.56
2:B:16:VAL:HG21	2:B:21:ALA:HB2	1.88	0.56
1:C:159:LEU:CD2	1:C:164:THR:HG22	2.34	0.56
1:G:418:ASN:C	1:G:418:ASN:HD22	2.12	0.56
2:H:130:LYS:HB3	2:H:134:GLU:HB2	1.86	0.56
1:K:161:PRO:HB3	1:K:209:THR:CG2	2.36	0.56
1:M:283:GLU:O	1:M:287:GLN:HG3	2.04	0.56
1:M:374:LEU:HG	1:M:398:LEU:HD21	1.87	0.56
1:A:331:ASP:OD2	1:A:356:ARG:NH1	2.39	0.56
1:E:419:GLN:HG2	1:E:419:GLN:O	2.06	0.56
1:G:124:LEU:HB3	1:G:128:GLU:HB3	1.86	0.56
1:K:148:LEU:CA	1:K:427:ARG:HH22	2.04	0.56
1:K:193:PHE:HB3	1:K:195:VAL:HG23	1.88	0.56
2:L:49:ASN:HB2	2:L:109:TYR:CE2	2.41	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:420:GLU:O	1:M:423:GLY:N	2.37	0.56
2:N:49:ASN:HB2	2:N:109:TYR:CE2	2.41	0.56
1:O:299:GLN:HG2	1:O:326:HIS:HB2	1.88	0.56
1:O:300:LEU:HD23	1:O:317:LEU:HD11	1.87	0.56
1:A:125:CYS:SG	1:A:127:PRO:HG2	2.45	0.56
1:A:418:ASN:HD22	1:A:418:ASN:C	2.13	0.56
1:G:159:LEU:CD2	1:G:164:THR:HG22	2.35	0.56
1:G:309:LEU:HD21	1:G:329:LEU:HD22	1.88	0.56
1:M:416:LYS:HD2	1:M:417:LYS:HZ1	1.67	0.56
1:A:418:ASN:C	1:A:418:ASN:ND2	2.63	0.56
1:C:140:TYR:CE2	2:D:153:VAL:HG22	2.40	0.56
1:C:309:LEU:HD23	1:C:332:SER:OG	2.06	0.56
1:I:397:HIS:CE1	1:I:398:LEU:HG	2.41	0.56
1:M:124:LEU:HD23	2:N:124:ALA:HB1	1.88	0.56
1:E:159:LEU:CD2	1:E:164:THR:HG22	2.35	0.56
1:K:418:ASN:C	1:K:418:ASN:HD22	2.14	0.56
2:L:127:ILE:HG23	2:L:135:ILE:HD13	1.88	0.56
1:M:376:THR:HG22	1:M:399:GLN:CB	2.35	0.56
2:N:65:HIS:HA	2:N:68:ASP:OD1	2.05	0.56
1:E:300:LEU:HD23	1:E:317:LEU:HD11	1.88	0.56
1:G:253:SER:HB2	1:G:280:ASP:CB	2.35	0.56
1:G:283:GLU:O	1:G:287:GLN:HG3	2.06	0.56
1:G:419:GLN:OE1	1:G:429:THR:HG21	2.05	0.56
1:M:281:PHE:CE1	1:M:305:TYR:CE2	2.94	0.56
1:A:108:SER:OG	2:B:140:ASN:ND2	2.34	0.56
2:F:30:MET:HB3	2:F:36:MET:HE2	1.88	0.56
1:G:418:ASN:C	1:G:418:ASN:ND2	2.62	0.56
1:O:215:SER:O	1:O:240:LYS:NZ	2.39	0.56
1:A:149:TRP:O	1:A:172:VAL:HG22	2.06	0.55
1:E:220:LEU:CB	1:E:241:ASN:HD22	2.18	0.55
1:M:362:PRO:HG3	1:M:382:ILE:O	2.06	0.55
2:N:47:LEU:HD12	2:N:55:LEU:HD11	1.88	0.55
1:O:418:ASN:C	1:O:418:ASN:ND2	2.64	0.55
1:A:137:LYS:H	2:B:144:ASP:HB2	1.71	0.55
1:K:141:ARG:HB3	1:K:141:ARG:HH11	1.72	0.55
1:C:421:ILE:O	1:C:421:ILE:HG23	2.05	0.55
1:E:374:LEU:HG	1:E:398:LEU:HD21	1.88	0.55
1:M:128:GLU:CG	2:N:128:LYS:HG2	2.35	0.55
1:O:314:LEU:O	1:O:318:VAL:HG23	2.05	0.55
1:G:158:ASN:HA	1:G:181:PHE:HB2	1.89	0.55
1:G:377:LEU:O	1:G:400:ILE:HA	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:12:GLU:OE2	2:H:52:ALA:HB1	2.06	0.55
2:J:98:GLY:O	2:J:102:GLU:HG2	2.06	0.55
1:E:344:PHE:CE2	1:E:370:GLU:HB2	2.41	0.55
2:H:30:MET:HB3	2:H:36:MET:HE2	1.88	0.55
1:C:418:ASN:C	1:C:418:ASN:HD22	2.15	0.55
2:D:122:THR:O	2:D:126:MET:HG3	2.07	0.55
1:E:182:MET:O	1:E:205:ILE:HA	2.06	0.55
1:G:309:LEU:HD23	1:G:332:SER:OG	2.07	0.55
2:H:16:VAL:CG2	2:H:21:ALA:HB2	2.36	0.55
1:K:389:GLN:NE2	1:K:392:LYS:HD2	2.21	0.55
1:O:335:LEU:HD22	1:O:339:CYS:SG	2.47	0.55
1:A:182:MET:O	1:A:205:ILE:HA	2.06	0.55
1:E:141:ARG:HB3	1:E:141:ARG:HH11	1.71	0.55
2:H:43:ASP:O	2:H:45:VAL:N	2.39	0.55
2:H:98:GLY:O	2:H:102:GLU:HG2	2.07	0.55
1:I:274:ASN:C	1:I:274:ASN:ND2	2.65	0.55
2:N:16:VAL:HG21	2:N:21:ALA:HB2	1.88	0.55
1:O:416:LYS:O	1:O:417:LYS:HD3	2.06	0.55
1:A:151:THR:HB	1:A:426:CYS:HA	1.87	0.55
1:C:424:ILE:N	1:C:424:ILE:HD12	2.22	0.55
1:I:220:LEU:CB	1:I:241:ASN:ND2	2.68	0.55
1:M:257:GLU:HB2	1:M:288:VAL:HG21	1.88	0.55
1:M:281:PHE:CB	1:M:285:HIS:CG	2.89	0.55
1:A:155:THR:CG2	1:A:178:PRO:HD2	2.31	0.55
2:B:98:GLY:O	2:B:102:GLU:HG2	2.07	0.55
1:E:149:TRP:O	1:E:172:VAL:HG22	2.07	0.55
1:G:378:GLN:HE21	1:G:380:PHE:HE1	1.54	0.55
1:K:176:ARG:HD3	1:K:419:GLN:HA	1.87	0.55
2:D:101:PHE:CE1	2:D:105:LEU:HD11	2.42	0.55
1:G:151:THR:HB	1:G:426:CYS:HA	1.87	0.55
2:H:47:LEU:HD12	2:H:55:LEU:HD11	1.89	0.55
1:I:417:LYS:CD	1:M:417:LYS:CG	2.80	0.55
2:B:101:PHE:HE1	2:B:105:LEU:HD11	1.71	0.54
1:C:220:LEU:CB	1:C:241:ASN:HD22	2.18	0.54
1:C:418:ASN:C	1:C:418:ASN:ND2	2.64	0.54
2:D:54:ILE:HG13	2:D:102:GLU:HB3	1.87	0.54
1:E:157:LYS:NZ	1:E:430:LEU:HD13	2.22	0.54
1:E:257:GLU:HB2	1:E:288:VAL:HG21	1.88	0.54
2:F:12:GLU:OE2	2:F:52:ALA:HB1	2.07	0.54
2:H:61:TRP:HB2	2:H:92:PHE:CZ	2.43	0.54
1:I:182:MET:O	1:I:205:ILE:HA	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:158:ASN:HA	1:M:181:PHE:HB2	1.89	0.54
1:M:331:ASP:CG	1:M:356:ARG:HH11	2.13	0.54
2:N:88:TRP:O	2:N:91:GLU:HG3	2.07	0.54
1:O:309:LEU:HD23	1:O:332:SER:OG	2.07	0.54
1:G:374:LEU:HG	1:G:398:LEU:HD21	1.89	0.54
1:I:176:ARG:CD	1:I:419:GLN:HA	2.37	0.54
1:K:419:GLN:O	1:K:419:GLN:HG2	2.07	0.54
1:M:196:GLN:HB3	1:M:221:GLN:HE21	1.73	0.54
1:M:220:LEU:HB2	1:M:241:ASN:HD22	1.72	0.54
1:C:155:THR:HG21	1:C:176:ARG:NH1	2.23	0.54
1:G:141:ARG:HB3	1:G:141:ARG:HH11	1.71	0.54
1:K:420:GLU:O	1:K:420:GLU:CG	2.55	0.54
2:L:20:ILE:HA	2:L:23:GLN:HE21	1.71	0.54
1:C:397:HIS:CE1	1:C:398:LEU:HG	2.43	0.54
1:G:257:GLU:HB2	1:G:288:VAL:HG21	1.89	0.54
1:G:362:PRO:HG3	1:G:382:ILE:O	2.08	0.54
1:E:187:ALA:O	1:E:188:GLU:C	2.49	0.54
1:M:131:LYS:HD3	2:N:130:LYS:O	2.08	0.54
1:M:215:SER:O	1:M:240:LYS:NZ	2.39	0.54
1:A:362:PRO:HG3	1:A:382:ILE:O	2.08	0.54
1:C:340:PHE:N	1:C:340:PHE:CD1	2.74	0.54
1:C:415:ASN:C	1:C:417:LYS:H	2.14	0.54
1:G:182:MET:O	1:G:205:ILE:HA	2.08	0.54
1:K:418:ASN:C	1:K:418:ASN:ND2	2.65	0.54
1:M:415:ASN:C	1:M:417:LYS:H	2.14	0.54
1:I:414:GLY:CA	1:M:416:LYS:HZ2	2.14	0.54
1:M:370:GLU:O	1:M:372:PRO:HD3	2.08	0.54
1:M:420:GLU:O	1:M:420:GLU:CG	2.56	0.54
2:N:22:LYS:O	2:N:22:LYS:HD3	2.06	0.54
1:O:279:PHE:C	1:O:279:PHE:CD1	2.86	0.54
2:P:65:HIS:CB	2:P:68:ASP:HB2	2.25	0.54
1:A:277:TRP:HZ2	1:A:331:ASP:OD2	1.90	0.54
2:F:130:LYS:HE3	2:F:138:THR:HG21	1.88	0.54
1:K:109:TRP:CD2	2:L:101:PHE:CD2	2.95	0.54
1:K:320:ARG:NH1	1:K:320:ARG:CB	2.71	0.54
2:F:133:GLU:HG3	2:F:137:LYS:HE2	1.90	0.54
2:H:16:VAL:HG21	2:H:21:ALA:HB2	1.89	0.54
1:O:281:PHE:HE1	1:O:305:TYR:CE2	2.24	0.54
1:A:377:LEU:O	1:A:400:ILE:HA	2.08	0.54
1:C:205:ILE:O	1:C:206:GLU:C	2.51	0.54
1:I:424:ILE:N	1:I:424:ILE:HD12	2.23	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:220:LEU:N	1:K:241:ASN:HD22	2.01	0.54
2:L:24:SER:OG	2:L:110:LEU:HB3	2.07	0.54
1:C:184:GLN:HB3	1:C:185:PRO:HD2	1.89	0.53
1:K:126:LEU:HD23	1:K:430:LEU:CD2	2.38	0.53
1:A:158:ASN:HA	1:A:181:PHE:HB2	1.88	0.53
1:A:187:ALA:O	1:A:188:GLU:C	2.51	0.53
1:A:380:PHE:CE2	1:A:401:ASN:HB3	2.43	0.53
2:D:61:TRP:CD2	2:D:115:LEU:HD23	2.44	0.53
1:I:418:ASN:HD22	1:I:418:ASN:C	2.16	0.53
1:K:137:LYS:HE3	2:L:145:PHE:HA	1.91	0.53
1:O:141:ARG:HB3	1:O:141:ARG:HH11	1.73	0.53
1:C:115:GLU:OE1	1:C:115:GLU:N	2.41	0.53
1:E:164:THR:HG21	1:E:182:MET:SD	2.49	0.53
1:E:419:GLN:NE2	1:E:426:CYS:O	2.40	0.53
1:G:420:GLU:CG	1:G:420:GLU:O	2.56	0.53
2:H:54:ILE:HG23	2:H:103:LEU:HD23	1.89	0.53
1:I:115:GLU:N	1:I:115:GLU:OE1	2.40	0.53
1:I:417:LYS:CG	1:M:417:LYS:CE	2.67	0.53
1:M:420:GLU:O	1:M:420:GLU:HG3	2.08	0.53
1:A:197:HIS:CD2	1:A:222:ASN:ND2	2.72	0.53
2:B:47:LEU:HD12	2:B:55:LEU:HD11	1.89	0.53
1:C:276:SER:HB3	1:C:301:ASN:OD1	2.08	0.53
1:C:325:VAL:HG13	1:C:350:GLN:HG2	1.89	0.53
1:G:186:LEU:N	1:G:187:ALA:N	2.55	0.53
1:M:153:ASP:HB3	1:M:429:THR:HG22	1.90	0.53
1:O:281:PHE:CB	1:O:285:HIS:CG	2.91	0.53
2:B:146:THR:HG22	2:B:148:GLU:H	1.73	0.53
2:B:149:GLU:O	2:B:153:VAL:HG23	2.08	0.53
1:G:300:LEU:HD23	1:G:317:LEU:HD11	1.91	0.53
1:K:124:LEU:HB3	1:K:128:GLU:HB3	1.91	0.53
1:K:133:SER:O	1:K:140:TYR:HB2	2.08	0.53
1:O:231:SER:OG	1:O:233:PRO:HD2	2.09	0.53
1:E:157:LYS:HD2	1:E:430:LEU:HD12	1.89	0.53
2:F:101:PHE:CE1	2:F:105:LEU:HD11	2.44	0.53
1:G:419:GLN:HG2	1:G:419:GLN:O	2.08	0.53
1:I:231:SER:OG	1:I:233:PRO:HD2	2.08	0.53
2:L:16:VAL:HG21	2:L:21:ALA:HB2	1.90	0.53
1:O:220:LEU:CB	1:O:241:ASN:HD22	2.22	0.53
1:C:384:PRO:HD2	1:C:387:THR:OG1	2.08	0.53
1:E:157:LYS:HD2	1:E:430:LEU:CD1	2.38	0.53
1:E:189:HIS:HA	1:E:216:GLN:NE2	2.24	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:331:ASP:CG	1:E:356:ARG:HH11	2.17	0.53
1:E:378:GLN:NE2	1:E:401:ASN:HA	2.24	0.53
1:G:157:LYS:HD2	1:G:430:LEU:CD1	2.39	0.53
1:G:370:GLU:O	1:G:372:PRO:HD3	2.09	0.53
2:J:54:ILE:O	2:J:57:LYS:HB2	2.09	0.53
1:K:415:ASN:C	1:K:417:LYS:H	2.15	0.53
1:M:151:THR:CB	1:M:426:CYS:HA	2.39	0.53
1:A:246:ARG:NE	1:A:272:GLU:OE2	2.41	0.53
1:A:300:LEU:HD23	1:A:317:LEU:HD11	1.89	0.53
1:C:124:LEU:HB3	1:C:128:GLU:HB3	1.90	0.53
1:C:131:LYS:HD3	2:D:130:LYS:O	2.09	0.53
1:I:416:LYS:O	1:I:417:LYS:HD3	2.08	0.53
1:O:420:GLU:CG	1:O:420:GLU:O	2.57	0.53
2:P:115:LEU:O	2:P:115:LEU:HD13	2.09	0.53
1:I:390:LEU:HA	1:I:393:GLU:HG2	1.89	0.53
1:M:149:TRP:O	1:M:172:VAL:HG22	2.09	0.53
2:B:61:TRP:CE3	2:B:115:LEU:HB2	2.44	0.53
1:E:420:GLU:CG	1:E:420:GLU:O	2.56	0.53
1:G:184:GLN:CB	1:G:185:PRO:HD2	2.39	0.53
1:M:109:TRP:CD2	2:N:101:PHE:HD2	2.27	0.53
1:M:367:GLU:C	1:M:369:GLY:H	2.17	0.53
2:P:133:GLU:HA	2:P:136:ARG:NH1	2.24	0.53
1:A:416:LYS:CB	1:A:417:LYS:HE2	2.26	0.52
2:F:61:TRP:CG	2:F:115:LEU:HD23	2.44	0.52
1:G:157:LYS:HG2	1:G:159:LEU:HD13	1.90	0.52
1:I:331:ASP:OD1	1:I:356:ARG:NH1	2.36	0.52
1:I:416:LYS:CB	1:I:417:LYS:HE2	2.29	0.52
2:N:16:VAL:CG1	2:N:59:ILE:HD13	2.39	0.52
1:E:184:GLN:HB3	1:E:185:PRO:HD2	1.91	0.52
2:F:22:LYS:O	2:F:22:LYS:HD3	2.09	0.52
1:I:137:LYS:HE3	2:J:144:ASP:O	2.08	0.52
1:I:300:LEU:HD23	1:I:317:LEU:HD11	1.91	0.52
1:I:418:ASN:C	1:I:418:ASN:ND2	2.65	0.52
1:K:186:LEU:O	1:K:188:GLU:N	2.42	0.52
1:M:235:VAL:HG21	1:M:255:PHE:CG	2.45	0.52
1:A:420:GLU:CG	1:A:420:GLU:O	2.58	0.52
1:C:417:LYS:HG3	1:G:417:LYS:NZ	2.24	0.52
1:E:124:LEU:HB3	1:E:128:GLU:HB3	1.91	0.52
1:I:157:LYS:HG2	1:I:159:LEU:HD13	1.91	0.52
1:I:351:HIS:CD2	1:I:376:THR:OG1	2.56	0.52
2:J:85:ILE:HG22	2:J:90:GLN:HG3	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:J:127:ILE:HG23	2:J:135:ILE:HD11	1.90	0.52
1:K:417:LYS:HG3	1:O:417:LYS:NZ	2.25	0.52
1:M:125:CYS:SG	1:M:127:PRO:HG2	2.49	0.52
1:M:274:ASN:C	1:M:274:ASN:HD22	2.17	0.52
1:O:314:LEU:O	1:O:314:LEU:HD23	2.09	0.52
1:A:253:SER:HB2	1:A:280:ASP:CG	2.34	0.52
1:A:414:GLY:HA2	1:E:416:LYS:HZ3	1.74	0.52
2:F:117:ASP:OD1	2:F:121:LYS:HE3	2.09	0.52
1:I:190:PHE:N	1:I:190:PHE:CD1	2.78	0.52
1:I:196:GLN:HB3	1:I:221:GLN:HE21	1.74	0.52
1:O:418:ASN:C	1:O:418:ASN:HD22	2.17	0.52
1:C:417:LYS:CD	1:G:417:LYS:HD2	2.40	0.52
1:G:155:THR:CG2	1:G:178:PRO:HD2	2.27	0.52
1:G:420:GLU:O	1:G:423:GLY:N	2.39	0.52
2:N:26:THR:HB	2:N:110:LEU:HA	1.92	0.52
1:O:168:LEU:HD22	1:O:195:VAL:CG2	2.40	0.52
1:C:257:GLU:OE2	1:C:257:GLU:N	2.36	0.52
1:I:420:GLU:O	1:I:423:GLY:N	2.40	0.52
2:B:54:ILE:HG21	2:B:106:ALA:HB2	1.92	0.52
1:C:141:ARG:HB3	1:C:141:ARG:HH11	1.74	0.52
1:G:420:GLU:O	1:G:420:GLU:HG3	2.08	0.52
1:M:378:GLN:NE2	1:M:401:ASN:HA	2.25	0.52
1:C:374:LEU:HG	1:C:398:LEU:HD21	1.91	0.52
1:E:130:LEU:HD12	2:F:160:CYS:SG	2.50	0.52
1:E:235:VAL:HG21	1:E:255:PHE:CG	2.45	0.52
1:E:415:ASN:C	1:E:417:LYS:H	2.18	0.52
1:I:325:VAL:HG13	1:I:350:GLN:HG2	1.92	0.52
1:K:149:TRP:O	1:K:172:VAL:HG22	2.09	0.52
2:L:133:GLU:O	2:L:137:LYS:HG3	2.10	0.52
1:M:237:THR:HG23	1:M:240:LYS:HE3	1.92	0.52
1:C:420:GLU:O	1:C:421:ILE:C	2.53	0.52
1:I:272:GLU:HG2	1:I:299:GLN:HB2	1.92	0.52
1:M:411:PRO:HG3	1:M:422:TRP:CH2	2.45	0.52
1:A:314:LEU:O	1:A:318:VAL:HG23	2.10	0.52
1:C:247:LEU:HD21	1:C:263:LEU:HD21	1.92	0.52
1:C:377:LEU:O	1:C:400:ILE:HA	2.10	0.52
1:G:193:PHE:HB3	1:G:195:VAL:HG23	1.92	0.52
2:H:143:ASN:OD1	2:H:145:PHE:HB2	2.09	0.52
1:I:159:LEU:HD23	1:I:164:THR:CG2	2.39	0.52
1:K:399:GLN:HA	1:K:399:GLN:NE2	2.13	0.52
1:C:419:GLN:O	1:C:419:GLN:HG2	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:115:GLU:OE1	1:G:115:GLU:N	2.43	0.51
1:I:161:PRO:HB3	1:I:209:THR:HG23	1.91	0.51
1:K:235:VAL:HG21	1:K:255:PHE:CG	2.45	0.51
1:K:416:LYS:HZ3	1:O:414:GLY:CA	2.23	0.51
2:L:26:THR:HB	2:L:110:LEU:HA	1.92	0.51
2:L:136:ARG:CZ	2:L:143:ASN:ND2	2.73	0.51
1:I:124:LEU:HB3	1:I:128:GLU:HB3	1.92	0.51
1:K:109:TRP:HZ3	2:L:104:ILE:HD12	1.74	0.51
2:L:101:PHE:HE1	2:L:105:LEU:HD11	1.75	0.51
1:M:109:TRP:CZ2	2:N:101:PHE:HB2	2.46	0.51
1:M:161:PRO:HB3	1:M:209:THR:CG2	2.40	0.51
2:P:127:ILE:HG23	2:P:135:ILE:HD13	1.93	0.51
1:A:340:PHE:N	1:A:340:PHE:CD1	2.78	0.51
2:B:12:GLU:OE2	2:B:52:ALA:HB1	2.11	0.51
1:E:159:LEU:HD23	1:E:164:THR:CG2	2.38	0.51
1:G:367:GLU:C	1:G:369:GLY:H	2.18	0.51
2:H:124:ALA:O	2:H:128:LYS:HG3	2.11	0.51
1:M:220:LEU:H	1:M:241:ASN:ND2	2.05	0.51
2:P:49:ASN:HB2	2:P:109:TYR:CE2	2.45	0.51
1:C:191:SER:HB3	1:C:192:PRO:CD	2.35	0.51
1:C:299:GLN:HG2	1:C:326:HIS:HB2	1.91	0.51
1:C:416:LYS:HZ3	1:G:414:GLY:CA	2.23	0.51
2:F:104:ILE:HG12	2:F:119:THR:HB	1.93	0.51
1:G:419:GLN:HG3	1:G:426:CYS:HB2	1.91	0.51
1:I:419:GLN:OE1	1:I:429:THR:HG21	2.11	0.51
1:O:420:GLU:O	1:O:423:GLY:N	2.42	0.51
1:C:151:THR:CB	1:C:426:CYS:HA	2.40	0.51
2:D:26:THR:O	2:D:29:THR:HG22	2.11	0.51
1:E:133:SER:O	1:E:140:TYR:HB2	2.10	0.51
1:E:151:THR:HB	1:E:426:CYS:HA	1.92	0.51
1:E:420:GLU:O	1:E:423:GLY:N	2.42	0.51
2:F:61:TRP:CD2	2:F:115:LEU:HD23	2.46	0.51
1:G:159:LEU:HD23	1:G:164:THR:CG2	2.36	0.51
1:G:383:VAL:CG1	1:G:388:LEU:HB2	2.40	0.51
1:M:378:GLN:HE21	1:M:380:PHE:HE1	1.58	0.51
1:A:254:GLY:H	1:A:280:ASP:CG	2.19	0.51
1:G:149:TRP:O	1:G:172:VAL:HA	2.11	0.51
2:N:24:SER:OG	2:N:110:LEU:HB3	2.09	0.51
1:A:181:PHE:CE1	1:C:384:PRO:HA	2.45	0.51
1:A:237:THR:CG2	1:A:240:LYS:HE3	2.40	0.51
1:A:307:LYS:HD2	1:A:307:LYS:C	2.36	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:351:HIS:CD2	1:C:376:THR:OG1	2.60	0.51
1:G:157:LYS:NZ	1:G:430:LEU:HD13	2.26	0.51
1:O:351:HIS:CD2	1:O:376:THR:OG1	2.55	0.51
1:O:419:GLN:O	1:O:419:GLN:HG2	2.10	0.51
1:C:281:PHE:HE1	1:C:305:TYR:CE2	2.28	0.51
2:D:61:TRP:CE3	2:D:115:LEU:HB2	2.45	0.51
1:G:351:HIS:CD2	1:G:376:THR:OG1	2.62	0.51
2:H:16:VAL:CG1	2:H:59:ILE:HD13	2.40	0.51
2:L:132:PRO:O	2:L:136:ARG:HG3	2.10	0.51
1:O:149:TRP:O	1:O:172:VAL:HA	2.10	0.51
1:O:153:ASP:HB3	1:O:429:THR:HG22	1.91	0.51
1:A:417:LYS:NZ	1:E:417:LYS:HG3	2.25	0.51
2:B:61:TRP:CG	2:B:115:LEU:HD23	2.45	0.51
1:O:394:ALA:O	1:O:396:PRO:HD2	2.11	0.51
2:F:98:GLY:O	2:F:102:GLU:HG2	2.11	0.51
1:G:189:HIS:H	1:G:190:PHE:HD1	1.58	0.51
1:I:133:SER:O	1:I:140:TYR:HB2	2.11	0.51
1:I:420:GLU:O	1:I:420:GLU:CG	2.58	0.51
2:J:87:VAL:HG21	1:M:296:THR:HA	1.92	0.51
1:K:320:ARG:CZ	1:K:320:ARG:HB3	2.41	0.51
1:K:376:THR:HG22	1:K:399:GLN:CB	2.41	0.51
1:K:417:LYS:CD	1:O:417:LYS:HD2	2.41	0.51
1:M:130:LEU:HD11	2:N:159:TRP:CZ2	2.46	0.51
1:M:231:SER:HA	1:M:254:GLY:O	2.11	0.51
1:O:115:GLU:N	1:O:115:GLU:OE1	2.44	0.51
1:O:420:GLU:O	1:O:421:ILE:C	2.52	0.51
1:A:181:PHE:CG	1:C:384:PRO:HG3	2.45	0.50
1:A:419:GLN:HG3	1:A:426:CYS:HB2	1.94	0.50
1:C:149:TRP:O	1:C:172:VAL:HG22	2.11	0.50
1:E:424:ILE:HD12	1:E:424:ILE:N	2.26	0.50
1:I:162:ASP:OD2	1:I:186:LEU:HD12	2.11	0.50
1:I:215:SER:O	1:I:240:LYS:NZ	2.44	0.50
2:J:132:PRO:O	2:J:136:ARG:HG3	2.10	0.50
1:K:184:GLN:CB	1:K:185:PRO:HD2	2.41	0.50
1:K:382:ILE:HG13	1:K:383:VAL:HG23	1.93	0.50
1:M:116:LEU:HD21	2:N:108:ASN:OD1	2.10	0.50
1:M:331:ASP:CG	1:M:356:ARG:NH1	2.69	0.50
2:P:86:PRO:O	2:P:90:GLN:HG3	2.10	0.50
2:B:22:LYS:O	2:B:22:LYS:HD3	2.11	0.50
1:C:137:LYS:HE3	2:D:145:PHE:HA	1.91	0.50
2:D:24:SER:OG	2:D:110:LEU:HB3	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:54:ILE:HG23	2:D:103:LEU:HD23	1.93	0.50
1:K:189:HIS:H	1:K:190:PHE:HD1	1.59	0.50
1:O:220:LEU:CB	1:O:241:ASN:ND2	2.74	0.50
1:A:107:VAL:HG12	1:A:108:SER:N	2.26	0.50
1:A:424:ILE:N	1:A:424:ILE:HD12	2.26	0.50
1:C:126:LEU:HD23	1:C:430:LEU:HG	1.93	0.50
1:E:220:LEU:N	1:E:241:ASN:HD22	2.05	0.50
2:F:85:ILE:HG22	2:F:90:GLN:HG3	1.93	0.50
1:M:184:GLN:CB	1:M:185:PRO:HD2	2.41	0.50
1:M:325:VAL:HG13	1:M:350:GLN:HG2	1.93	0.50
1:O:284:LYS:O	1:O:288:VAL:HG23	2.12	0.50
1:A:376:THR:HG22	1:A:399:GLN:HB3	1.93	0.50
2:D:98:GLY:O	2:D:102:GLU:HG2	2.11	0.50
1:E:253:SER:HB3	1:E:280:ASP:HB2	1.92	0.50
1:I:253:SER:HB2	1:I:280:ASP:CG	2.37	0.50
2:J:51:ASN:ND2	2:J:53:ALA:H	2.09	0.50
2:J:127:ILE:HG23	2:J:135:ILE:HD13	1.92	0.50
1:K:370:GLU:O	1:K:372:PRO:HD3	2.11	0.50
1:C:417:LYS:CE	1:G:417:LYS:CG	2.71	0.50
1:K:419:GLN:O	1:K:425:LYS:HA	2.10	0.50
1:A:220:LEU:CB	1:A:241:ASN:ND2	2.75	0.50
1:A:278:CYS:HB2	1:A:281:PHE:CE2	2.47	0.50
1:A:417:LYS:HD2	1:E:417:LYS:CD	2.40	0.50
2:D:65:HIS:CB	2:D:68:ASP:HB2	2.32	0.50
1:E:196:GLN:HB3	1:E:221:GLN:HE21	1.76	0.50
1:E:377:LEU:O	1:E:400:ILE:HA	2.11	0.50
1:G:314:LEU:CD2	1:G:314:LEU:O	2.59	0.50
1:K:237:THR:HG23	1:K:240:LYS:HE3	1.92	0.50
2:P:16:VAL:HG23	2:P:21:ALA:HB2	1.93	0.50
1:A:157:LYS:HG2	1:A:159:LEU:HD13	1.93	0.50
1:A:414:GLY:CA	1:E:416:LYS:NZ	2.72	0.50
1:A:421:ILE:HG23	1:A:421:ILE:O	2.10	0.50
1:E:220:LEU:CB	1:E:241:ASN:ND2	2.75	0.50
1:G:220:LEU:N	1:G:241:ASN:HD22	2.07	0.50
1:K:197:HIS:CD2	1:K:222:ASN:ND2	2.78	0.50
1:M:419:GLN:NE2	1:M:426:CYS:O	2.44	0.50
1:A:384:PRO:HG3	1:C:181:PHE:CD1	2.47	0.50
2:D:61:TRP:CG	2:D:115:LEU:HD23	2.47	0.50
1:K:416:LYS:NZ	1:O:414:GLY:CA	2.73	0.50
1:O:196:GLN:HB3	1:O:221:GLN:HE21	1.77	0.50
1:O:274:ASN:C	1:O:274:ASN:ND2	2.69	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:378:GLN:HE22	1:A:401:ASN:HA	1.74	0.50
2:D:130:LYS:HB3	2:D:134:GLU:HB2	1.93	0.50
2:F:64:HIS:HD2	2:F:65:HIS:CE1	2.29	0.50
2:F:127:ILE:HG23	2:F:135:ILE:CD1	2.42	0.50
2:H:122:THR:O	2:H:126:MET:HG3	2.12	0.50
1:I:159:LEU:CD2	1:I:164:THR:HG22	2.42	0.50
1:I:340:PHE:N	1:I:340:PHE:CD1	2.78	0.50
1:I:415:ASN:C	1:I:417:LYS:H	2.20	0.50
1:K:115:GLU:OE1	1:K:115:GLU:N	2.45	0.50
1:M:164:THR:O	1:M:168:LEU:HG	2.11	0.50
1:O:314:LEU:O	1:O:314:LEU:CD2	2.59	0.50
2:P:12:GLU:OE2	2:P:52:ALA:HB1	2.12	0.50
1:A:133:SER:O	1:A:140:TYR:HB2	2.12	0.49
2:B:25:VAL:HB	2:B:111:ASP:OD2	2.12	0.49
2:B:130:LYS:CE	2:B:138:THR:HG21	2.40	0.49
1:C:314:LEU:O	1:C:314:LEU:CD2	2.60	0.49
1:E:331:ASP:OD2	1:E:356:ARG:NH1	2.45	0.49
1:G:281:PHE:CE1	1:G:305:TYR:CE2	2.99	0.49
1:I:164:THR:HG21	1:I:182:MET:SD	2.52	0.49
2:L:22:LYS:O	2:L:22:LYS:HD3	2.12	0.49
1:M:166:ARG:NH1	2:N:156:GLU:HA	2.27	0.49
1:A:188:GLU:O	1:A:212:GLY:O	2.28	0.49
1:A:367:GLU:C	1:A:369:GLY:H	2.20	0.49
1:A:383:VAL:CG1	1:A:388:LEU:HB2	2.41	0.49
2:B:61:TRP:CD1	2:B:115:LEU:HD23	2.47	0.49
2:B:133:GLU:O	2:B:137:LYS:HG3	2.12	0.49
1:E:244:LEU:HD12	1:E:267:CYS:SG	2.53	0.49
2:F:47:LEU:HD12	2:F:55:LEU:HD11	1.94	0.49
2:H:32:GLU:C	2:H:34:LEU:H	2.19	0.49
1:I:175:PHE:CD2	1:I:198:MET:HE2	2.41	0.49
2:J:61:TRP:CE3	2:J:115:LEU:HB2	2.48	0.49
2:N:30:MET:HB3	2:N:36:MET:HE2	1.94	0.49
1:A:193:PHE:HB3	1:A:195:VAL:HG23	1.93	0.49
1:A:417:LYS:CD	1:E:417:LYS:CG	2.86	0.49
2:B:133:GLU:HG3	2:B:137:LYS:HE2	1.93	0.49
1:C:420:GLU:O	1:C:420:GLU:CG	2.59	0.49
1:G:153:ASP:HB2	1:G:426:CYS:HB2	1.92	0.49
2:H:107:ALA:CB	2:H:115:LEU:HD12	2.33	0.49
1:K:247:LEU:HD21	1:K:263:LEU:HD21	1.95	0.49
1:M:344:PHE:CE2	1:M:370:GLU:HB2	2.48	0.49
1:M:419:GLN:HG2	1:M:419:GLN:O	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:136:CYS:SG	1:A:139:TRP:CD1	3.05	0.49
2:H:8:SER:C	2:H:10:ASP:H	2.20	0.49
2:J:16:VAL:CG1	2:J:59:ILE:HD13	2.42	0.49
1:A:184:GLN:CB	1:A:185:PRO:CD	2.90	0.49
2:B:6:LEU:O	2:B:13:ILE:HA	2.13	0.49
2:D:49:ASN:HB2	2:D:109:TYR:CE2	2.47	0.49
1:E:107:VAL:HG12	1:E:108:SER:N	2.27	0.49
1:G:320:ARG:HB2	1:G:320:ARG:HH11	1.76	0.49
2:H:6:LEU:O	2:H:13:ILE:HA	2.12	0.49
2:H:127:ILE:HG23	2:H:135:ILE:CD1	2.42	0.49
1:I:184:GLN:CG	1:I:185:PRO:HD2	2.42	0.49
2:J:61:TRP:CG	2:J:115:LEU:HD23	2.47	0.49
1:K:220:LEU:CB	1:K:241:ASN:ND2	2.75	0.49
2:N:133:GLU:HG3	2:N:137:LYS:HE2	1.93	0.49
1:A:153:ASP:HB3	1:A:429:THR:HG22	1.94	0.49
1:C:420:GLU:O	1:C:423:GLY:N	2.44	0.49
2:D:30:MET:C	2:D:32:GLU:H	2.20	0.49
2:D:32:GLU:C	2:D:34:LEU:H	2.21	0.49
1:E:181:PHE:CD1	1:G:384:PRO:HG3	2.48	0.49
1:E:253:SER:HB2	1:E:280:ASP:CG	2.37	0.49
1:E:331:ASP:CG	1:E:356:ARG:NH1	2.71	0.49
2:F:127:ILE:HG23	2:F:135:ILE:HD13	1.93	0.49
2:J:54:ILE:HG21	2:J:106:ALA:HB2	1.93	0.49
1:O:109:TRP:C	1:O:111:SER:H	2.20	0.49
1:G:107:VAL:HG12	1:G:108:SER:N	2.28	0.49
1:G:162:ASP:OD2	1:G:186:LEU:HD12	2.12	0.49
1:G:184:GLN:CG	1:G:185:PRO:HD2	2.43	0.49
1:I:189:HIS:H	1:I:190:PHE:HD1	1.60	0.49
1:I:370:GLU:O	1:I:372:PRO:HD3	2.12	0.49
1:K:197:HIS:HD2	1:K:222:ASN:ND2	2.00	0.49
1:K:294:SER:OG	1:K:296:THR:OG1	2.24	0.49
1:O:232:ASP:HB2	1:O:233:PRO:HD3	1.93	0.49
1:O:415:ASN:C	1:O:417:LYS:H	2.19	0.49
2:P:32:GLU:C	2:P:34:LEU:H	2.20	0.49
1:A:220:LEU:CB	1:A:241:ASN:HD22	2.25	0.49
2:D:54:ILE:O	2:D:57:LYS:HB2	2.13	0.49
1:E:420:GLU:O	1:E:420:GLU:HG3	2.13	0.49
2:H:61:TRP:CG	2:H:115:LEU:HD23	2.48	0.49
1:I:116:LEU:HD21	2:J:108:ASN:CG	2.38	0.49
1:I:168:LEU:HD22	1:I:195:VAL:HG21	1.93	0.49
1:K:186:LEU:C	1:K:188:GLU:N	2.69	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:311:LYS:HE3	1:A:338:ASP:O	2.13	0.49
2:B:61:TRP:HB2	2:B:92:PHE:CZ	2.47	0.49
1:E:370:GLU:O	1:E:372:PRO:HD3	2.12	0.49
1:I:344:PHE:CE2	1:I:370:GLU:HB2	2.47	0.49
1:A:331:ASP:CG	1:A:356:ARG:NH1	2.71	0.49
2:B:22:LYS:HD3	2:B:22:LYS:C	2.38	0.49
1:E:225:LEU:O	1:E:252:CYS:SG	2.71	0.49
1:G:220:LEU:H	1:G:241:ASN:ND2	2.06	0.49
2:H:104:ILE:HG12	2:H:119:THR:HB	1.94	0.49
1:I:389:GLN:HE21	1:I:392:LYS:HD2	1.77	0.49
2:P:61:TRP:CZ3	2:P:115:LEU:HB2	2.47	0.49
2:P:133:GLU:HG3	2:P:137:LYS:HE2	1.94	0.49
1:C:278:CYS:HB2	1:C:281:PHE:CZ	2.48	0.48
1:G:389:GLN:HE22	1:G:392:LYS:NZ	2.11	0.48
2:H:2:PRO:C	2:H:18:VAL:HG13	2.38	0.48
2:L:117:ASP:OD1	2:L:121:LYS:HE3	2.13	0.48
2:N:22:LYS:HD3	2:N:22:LYS:C	2.38	0.48
1:O:163:VAL:O	1:O:167:LEU:HG	2.13	0.48
1:O:418:ASN:HD22	1:O:418:ASN:N	2.10	0.48
2:P:22:LYS:C	2:P:22:LYS:HD3	2.38	0.48
1:A:161:PRO:HB3	1:A:209:THR:CG2	2.43	0.48
2:B:55:LEU:HD22	2:B:59:ILE:HD11	1.94	0.48
1:E:137:LYS:CE	2:F:149:GLU:OE1	2.61	0.48
1:G:153:ASP:HB2	1:G:426:CYS:CB	2.42	0.48
2:H:101:PHE:HE1	2:H:105:LEU:HD11	1.77	0.48
2:J:2:PRO:C	2:J:18:VAL:HG13	2.38	0.48
2:N:2:PRO:C	2:N:18:VAL:HG13	2.37	0.48
1:O:190:PHE:N	1:O:190:PHE:CD1	2.81	0.48
1:C:193:PHE:HB3	1:C:195:VAL:HG23	1.93	0.48
1:E:130:LEU:HD12	2:F:160:CYS:CB	2.44	0.48
2:F:51:ASN:ND2	2:F:53:ALA:H	2.11	0.48
1:G:149:TRP:O	1:G:172:VAL:HG22	2.13	0.48
1:G:176:ARG:NE	1:G:419:GLN:H	2.10	0.48
1:K:175:PHE:CD2	1:K:198:MET:HE2	2.48	0.48
1:M:129:LEU:O	1:M:132:VAL:HG22	2.14	0.48
2:N:130:LYS:CE	2:N:138:THR:HG21	2.42	0.48
2:P:133:GLU:HA	2:P:136:ARG:HH11	1.76	0.48
1:A:376:THR:HG22	1:A:399:GLN:CB	2.43	0.48
1:C:390:LEU:HA	1:C:393:GLU:HG2	1.95	0.48
1:E:301:ASN:HB2	1:E:408:ILE:HD12	1.94	0.48
2:F:24:SER:OG	2:F:110:LEU:HB3	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:22:LYS:O	2:H:22:LYS:HD3	2.13	0.48
1:K:417:LYS:CG	1:O:417:LYS:CD	2.81	0.48
1:M:284:LYS:O	1:M:288:VAL:HG23	2.13	0.48
1:M:415:ASN:C	1:M:417:LYS:N	2.71	0.48
1:O:184:GLN:CB	1:O:185:PRO:HD2	2.43	0.48
1:A:417:LYS:CG	1:E:417:LYS:CE	2.70	0.48
1:G:314:LEU:O	1:G:318:VAL:HG23	2.13	0.48
2:J:95:VAL:CG2	2:J:99:THR:HB	2.43	0.48
1:K:416:LYS:CG	1:K:417:LYS:HE2	2.43	0.48
1:M:168:LEU:HD22	1:M:195:VAL:CG2	2.44	0.48
1:A:159:LEU:CD2	1:A:164:THR:HG22	2.42	0.48
2:F:30:MET:C	2:F:32:GLU:H	2.22	0.48
2:F:133:GLU:HA	2:F:136:ARG:NH1	2.28	0.48
1:G:390:LEU:HA	1:G:393:GLU:HG2	1.95	0.48
1:K:309:LEU:HD21	1:K:329:LEU:HD22	1.96	0.48
1:K:417:LYS:CG	1:O:417:LYS:CE	2.87	0.48
1:M:115:GLU:OE1	1:M:115:GLU:N	2.46	0.48
2:B:51:ASN:ND2	2:B:53:ALA:H	2.11	0.48
1:C:135:VAL:HG13	2:D:136:ARG:HG2	1.94	0.48
1:C:378:GLN:HE22	1:C:401:ASN:HA	1.78	0.48
1:C:416:LYS:HZ3	1:G:414:GLY:HA2	1.77	0.48
2:F:101:PHE:HE1	2:F:105:LEU:HD11	1.78	0.48
1:G:225:LEU:O	1:G:252:CYS:SG	2.72	0.48
1:I:320:ARG:HB3	1:I:320:ARG:CZ	2.43	0.48
2:J:149:GLU:O	2:J:153:VAL:HG23	2.14	0.48
1:K:161:PRO:HG3	1:K:184:GLN:O	2.13	0.48
1:K:231:SER:HA	1:K:254:GLY:O	2.14	0.48
1:M:133:SER:O	1:M:140:TYR:HB2	2.13	0.48
1:M:220:LEU:N	1:M:241:ASN:HD22	2.06	0.48
1:M:382:ILE:HG13	1:M:383:VAL:HG23	1.94	0.48
1:A:417:LYS:HZ2	1:E:417:LYS:HG3	1.78	0.48
2:B:16:VAL:CG1	2:B:59:ILE:HD13	2.44	0.48
1:C:133:SER:O	1:C:140:TYR:HB2	2.13	0.48
1:C:194:ARG:O	1:C:195:VAL:C	2.56	0.48
1:C:419:GLN:HG3	1:C:426:CYS:HB2	1.95	0.48
1:E:421:ILE:HG23	1:E:421:ILE:O	2.13	0.48
2:F:16:VAL:CG1	2:F:59:ILE:HD13	2.44	0.48
1:G:231:SER:OG	1:G:233:PRO:HD2	2.13	0.48
1:G:254:GLY:H	1:G:280:ASP:CG	2.21	0.48
2:J:122:THR:O	2:J:126:MET:HG3	2.14	0.48
2:N:6:LEU:O	2:N:13:ILE:HA	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:P:61:TRP:HB2	2:P:92:PHE:CZ	2.48	0.48
2:P:102:GLU:HA	2:P:102:GLU:OE2	2.14	0.48
1:A:395:LEU:HD13	1:A:398:LEU:HD12	1.96	0.48
2:D:127:ILE:HG23	2:D:135:ILE:CD1	2.44	0.48
1:E:186:LEU:O	1:E:188:GLU:N	2.47	0.48
1:E:187:ALA:C	1:E:212:GLY:HA3	2.39	0.48
1:G:125:CYS:SG	1:G:127:PRO:HG2	2.54	0.48
1:G:137:LYS:HE3	2:H:145:PHE:HA	1.96	0.48
2:L:47:LEU:HD12	2:L:55:LEU:HD11	1.96	0.48
1:M:420:GLU:O	1:M:421:ILE:C	2.56	0.48
1:O:159:LEU:CD2	1:O:164:THR:HG22	2.44	0.48
1:O:374:LEU:HG	1:O:398:LEU:HD21	1.96	0.48
1:O:384:PRO:HD2	1:O:387:THR:OG1	2.14	0.48
1:A:241:ASN:O	1:A:244:LEU:HG	2.13	0.48
1:E:309:LEU:HD21	1:E:329:LEU:HD22	1.95	0.48
2:F:32:GLU:C	2:F:34:LEU:H	2.22	0.48
1:G:231:SER:HA	1:G:254:GLY:O	2.14	0.48
1:K:231:SER:HB2	1:K:233:PRO:HD2	1.96	0.48
2:N:146:THR:HB	2:N:149:GLU:HG3	1.94	0.48
2:B:117:ASP:OD1	2:B:121:LYS:HE3	2.14	0.47
2:B:124:ALA:O	2:B:128:LYS:HG3	2.13	0.47
1:E:367:GLU:C	1:E:369:GLY:H	2.22	0.47
1:E:418:ASN:HD22	1:E:418:ASN:N	2.12	0.47
2:F:22:LYS:HD3	2:F:22:LYS:C	2.39	0.47
1:I:137:LYS:HG3	2:J:144:ASP:HB2	1.96	0.47
1:O:340:PHE:CD1	1:O:340:PHE:N	2.81	0.47
1:A:419:GLN:NE2	1:A:426:CYS:O	2.47	0.47
1:C:314:LEU:O	1:C:314:LEU:HD23	2.15	0.47
1:C:340:PHE:N	1:C:340:PHE:HD1	2.13	0.47
2:D:131:THR:OG1	2:D:134:GLU:HG3	2.14	0.47
1:E:109:TRP:C	1:E:111:SER:H	2.21	0.47
1:E:370:GLU:OE1	1:E:370:GLU:HA	2.14	0.47
1:K:126:LEU:HD23	1:K:430:LEU:HG	1.96	0.47
1:K:307:LYS:HD2	1:K:307:LYS:C	2.39	0.47
1:E:148:LEU:CA	1:E:427:ARG:HH22	2.21	0.47
1:K:190:PHE:CD1	1:K:190:PHE:N	2.81	0.47
1:K:416:LYS:HD2	1:K:417:LYS:HZ1	1.79	0.47
2:L:25:VAL:HB	2:L:111:ASP:OD2	2.13	0.47
2:L:65:HIS:HA	2:L:68:ASP:OD1	2.13	0.47
1:O:205:ILE:O	1:O:206:GLU:C	2.57	0.47
1:O:331:ASP:CG	1:O:356:ARG:HH11	2.21	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:148:LEU:CA	1:C:427:ARG:NH2	2.65	0.47
1:G:235:VAL:HG21	1:G:255:PHE:CG	2.48	0.47
1:G:419:GLN:O	1:G:425:LYS:HA	2.15	0.47
1:I:164:THR:HG21	1:I:182:MET:CE	2.44	0.47
1:I:301:ASN:HB2	1:I:408:ILE:HD12	1.96	0.47
1:I:418:ASN:HD22	1:I:418:ASN:N	2.11	0.47
1:I:419:GLN:O	1:I:419:GLN:HG2	2.14	0.47
2:J:22:LYS:HD3	2:J:22:LYS:C	2.40	0.47
2:J:32:GLU:C	2:J:34:LEU:H	2.23	0.47
2:J:50:VAL:HG23	2:J:109:TYR:CE2	2.50	0.47
1:K:107:VAL:HG12	1:K:108:SER:N	2.30	0.47
1:K:109:TRP:C	1:K:111:SER:H	2.23	0.47
1:K:331:ASP:CG	1:K:356:ARG:HH11	2.22	0.47
2:P:24:SER:HB3	2:P:27:ILE:HB	1.96	0.47
1:A:196:GLN:OE1	1:A:221:GLN:NE2	2.48	0.47
2:B:32:GLU:C	2:B:34:LEU:H	2.22	0.47
1:K:278:CYS:HB2	1:K:281:PHE:CE2	2.49	0.47
2:N:132:PRO:O	2:N:136:ARG:HG3	2.14	0.47
1:C:109:TRP:C	1:C:111:SER:H	2.22	0.47
2:D:61:TRP:HB2	2:D:92:PHE:CZ	2.49	0.47
1:E:137:LYS:HG2	2:F:145:PHE:CD2	2.50	0.47
1:E:420:GLU:O	1:E:421:ILE:C	2.58	0.47
1:G:340:PHE:CD1	1:G:340:PHE:N	2.82	0.47
2:H:127:ILE:HG23	2:H:135:ILE:HD13	1.97	0.47
2:L:130:LYS:CE	2:L:138:THR:HG21	2.44	0.47
2:P:16:VAL:HG21	2:P:21:ALA:HB2	1.95	0.47
1:A:235:VAL:HG21	1:A:255:PHE:CG	2.49	0.47
1:A:257:GLU:HB2	1:A:288:VAL:HG21	1.96	0.47
2:B:16:VAL:HG23	2:B:21:ALA:HB2	1.96	0.47
2:B:115:LEU:O	2:B:115:LEU:HD13	2.13	0.47
2:B:148:GLU:O	2:B:151:ALA:HB3	2.15	0.47
1:C:107:VAL:HG12	1:C:108:SER:N	2.29	0.47
1:C:415:ASN:C	1:C:417:LYS:N	2.72	0.47
1:C:418:ASN:HD22	1:C:418:ASN:N	2.12	0.47
1:E:382:ILE:HG13	1:E:383:VAL:HG23	1.96	0.47
1:G:137:LYS:CE	2:H:144:ASP:O	2.62	0.47
1:G:190:PHE:N	1:G:190:PHE:CD1	2.83	0.47
1:G:220:LEU:CB	1:G:241:ASN:HD22	2.27	0.47
1:G:277:TRP:HZ2	1:G:331:ASP:OD2	1.97	0.47
1:G:278:CYS:HB2	1:G:281:PHE:CE2	2.50	0.47
1:I:193:PHE:HB3	1:I:195:VAL:HG23	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:J:6:LEU:O	2:J:13:ILE:HA	2.15	0.47
1:K:220:LEU:CB	1:K:241:ASN:HD22	2.27	0.47
1:K:231:SER:CB	1:K:233:PRO:HD2	2.44	0.47
1:K:377:LEU:O	1:K:400:ILE:HA	2.15	0.47
1:M:159:LEU:N	1:M:181:PHE:O	2.42	0.47
1:M:176:ARG:NE	1:M:419:GLN:H	2.12	0.47
1:O:276:SER:HB3	1:O:301:ASN:OD1	2.14	0.47
1:O:301:ASN:HB2	1:O:408:ILE:HD12	1.97	0.47
2:P:54:ILE:O	2:P:57:LYS:HB2	2.15	0.47
2:D:85:ILE:HG22	2:D:90:GLN:HG3	1.95	0.47
1:I:107:VAL:HG12	1:I:108:SER:N	2.29	0.47
1:I:153:ASP:HB3	1:I:429:THR:HG22	1.96	0.47
2:N:32:GLU:C	2:N:34:LEU:H	2.23	0.47
2:N:101:PHE:CE1	2:N:105:LEU:HD11	2.49	0.47
2:P:122:THR:O	2:P:126:MET:HG3	2.15	0.47
1:A:176:ARG:NE	1:A:419:GLN:H	2.12	0.47
1:A:271:ASP:C	1:A:297:ILE:HD12	2.40	0.47
2:B:131:THR:O	2:B:135:ILE:HG13	2.15	0.47
2:B:155:LYS:HA	2:B:158:GLN:HG3	1.96	0.47
1:C:161:PRO:HB3	1:C:209:THR:CG2	2.41	0.47
1:E:279:PHE:CD1	1:E:279:PHE:C	2.93	0.47
1:E:340:PHE:N	1:E:340:PHE:CD1	2.81	0.47
1:K:155:THR:CG2	1:K:178:PRO:HD2	2.39	0.47
1:M:132:VAL:CG1	2:N:127:ILE:HG21	2.45	0.47
1:A:232:ASP:HB2	1:A:233:PRO:HD3	1.96	0.47
1:A:417:LYS:CE	1:E:417:LYS:CG	2.93	0.47
1:C:116:LEU:HD21	2:D:108:ASN:CG	2.40	0.47
1:C:164:THR:O	1:C:168:LEU:HG	2.14	0.47
1:C:419:GLN:O	1:C:425:LYS:HA	2.15	0.47
1:G:420:GLU:HG3	1:G:423:GLY:HA2	1.96	0.47
1:I:389:GLN:HE22	1:I:392:LYS:HZ1	1.63	0.47
1:K:162:ASP:OD2	1:K:186:LEU:HD12	2.15	0.47
2:L:133:GLU:HG3	2:L:137:LYS:HE2	1.96	0.47
2:P:139:PHE:O	2:P:140:ASN:O	2.33	0.47
1:A:420:GLU:HG3	1:A:423:GLY:HA2	1.96	0.46
1:C:331:ASP:CG	1:C:356:ARG:HH11	2.23	0.46
1:E:278:CYS:HB2	1:E:281:PHE:CZ	2.49	0.46
1:E:271:ASP:C	1:E:297:ILE:HD12	2.40	0.46
2:F:96:ASP:OD1	1:O:284:LYS:CE	2.48	0.46
2:F:133:GLU:HA	2:F:136:ARG:HH11	1.80	0.46
1:G:416:LYS:CD	1:G:417:LYS:NZ	2.74	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:420:GLU:O	1:G:421:ILE:C	2.59	0.46
1:M:185:PRO:O	1:M:186:LEU:HB2	2.15	0.46
1:O:162:ASP:OD2	1:O:186:LEU:HD12	2.16	0.46
1:A:163:VAL:HG11	2:B:159:TRP:CG	2.51	0.46
1:C:159:LEU:HD23	1:C:164:THR:CG2	2.41	0.46
1:C:389:GLN:HE21	1:C:392:LYS:HD2	1.76	0.46
1:G:320:ARG:CB	1:G:320:ARG:NH1	2.78	0.46
1:I:235:VAL:HG21	1:I:255:PHE:CG	2.50	0.46
1:I:278:CYS:HB2	1:I:281:PHE:CZ	2.50	0.46
1:I:335:LEU:HD22	1:I:339:CYS:SG	2.55	0.46
1:I:347:ASN:O	1:I:373:THR:HG21	2.15	0.46
1:K:415:ASN:C	1:K:417:LYS:N	2.72	0.46
2:L:22:LYS:HG2	2:L:28:LYS:HZ2	1.80	0.46
1:M:367:GLU:C	1:M:369:GLY:N	2.74	0.46
1:O:132:VAL:O	1:O:135:VAL:HG23	2.15	0.46
1:C:136:CYS:HG	1:C:139:TRP:CD1	2.32	0.46
1:E:356:ARG:HG2	1:E:358:TYR:OH	2.15	0.46
2:F:30:MET:HB3	2:F:36:MET:CE	2.46	0.46
2:F:148:GLU:O	2:F:151:ALA:HB3	2.14	0.46
1:G:164:THR:HG21	1:G:182:MET:SD	2.55	0.46
1:G:220:LEU:CB	1:G:241:ASN:ND2	2.77	0.46
1:G:274:ASN:C	1:G:274:ASN:ND2	2.71	0.46
1:G:356:ARG:HG2	1:G:358:TYR:OH	2.15	0.46
2:L:130:LYS:HB3	2:L:134:GLU:HB2	1.96	0.46
1:M:390:LEU:HA	1:M:393:GLU:HG2	1.97	0.46
1:O:109:TRP:C	1:O:111:SER:N	2.74	0.46
1:A:141:ARG:HB3	1:A:141:ARG:HH11	1.79	0.46
1:G:424:ILE:HD12	1:G:424:ILE:N	2.31	0.46
1:I:417:LYS:CE	1:M:417:LYS:CG	2.91	0.46
2:L:22:LYS:HD3	2:L:22:LYS:C	2.41	0.46
1:O:185:PRO:O	1:O:186:LEU:HB2	2.15	0.46
1:O:189:HIS:HA	1:O:216:GLN:NE2	2.30	0.46
1:O:246:ARG:HG2	1:O:272:GLU:HB2	1.98	0.46
2:B:49:ASN:HB2	2:B:109:TYR:CD2	2.51	0.46
1:C:341:GLN:HA	1:C:344:PHE:CD1	2.51	0.46
2:D:26:THR:HB	2:D:110:LEU:HA	1.97	0.46
2:D:132:PRO:O	2:D:136:ARG:HG3	2.16	0.46
1:G:215:SER:O	1:G:240:LYS:NZ	2.49	0.46
1:K:367:GLU:C	1:K:369:GLY:H	2.22	0.46
2:N:65:HIS:HB3	2:N:68:ASP:CB	2.22	0.46
1:O:189:HIS:H	1:O:190:PHE:HD1	1.64	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:O:278:CYS:C	1:O:280:ASP:N	2.74	0.46
1:A:215:SER:O	1:A:240:LYS:NZ	2.49	0.46
2:B:64:HIS:HD2	2:B:65:HIS:CE1	2.34	0.46
2:B:104:ILE:HG12	2:B:119:THR:HB	1.96	0.46
1:E:278:CYS:HB2	1:E:281:PHE:CE2	2.51	0.46
2:F:65:HIS:CB	2:F:68:ASP:HB2	2.29	0.46
1:G:184:GLN:HG2	1:G:185:PRO:HD2	1.98	0.46
1:I:199:ASP:OD1	1:I:411:PRO:HG2	2.14	0.46
1:I:374:LEU:HG	1:I:398:LEU:HD21	1.97	0.46
1:I:417:LYS:HD2	1:M:417:LYS:CD	2.46	0.46
2:J:61:TRP:CD2	2:J:115:LEU:HD23	2.51	0.46
2:L:16:VAL:CG1	2:L:59:ILE:HD13	2.45	0.46
1:O:107:VAL:HG12	1:O:108:SER:N	2.30	0.46
1:O:390:LEU:HA	1:O:393:GLU:HG2	1.97	0.46
2:P:30:MET:HB3	2:P:36:MET:HE2	1.97	0.46
1:C:328:ASP:HA	1:C:353:SER:HB2	1.98	0.46
1:E:229:ARG:CZ	1:E:253:SER:OG	2.64	0.46
2:F:6:LEU:O	2:F:13:ILE:HA	2.15	0.46
2:F:132:PRO:O	2:F:136:ARG:HG3	2.16	0.46
1:G:186:LEU:C	1:G:188:GLU:N	2.74	0.46
1:G:367:GLU:C	1:G:369:GLY:N	2.73	0.46
1:I:278:CYS:HB2	1:I:281:PHE:CE2	2.51	0.46
2:J:30:MET:HB3	2:J:36:MET:HE2	1.97	0.46
1:K:272:GLU:HG2	1:K:299:GLN:HB2	1.98	0.46
2:L:32:GLU:C	2:L:34:LEU:H	2.24	0.46
1:O:246:ARG:NE	1:O:272:GLU:OE2	2.48	0.46
1:O:331:ASP:OD2	1:O:356:ARG:NH1	2.49	0.46
1:O:367:GLU:C	1:O:369:GLY:H	2.24	0.46
2:B:33:ASP:O	2:B:34:LEU:HD23	2.16	0.46
1:C:184:GLN:CB	1:C:185:PRO:CD	2.94	0.46
2:D:2:PRO:C	2:D:18:VAL:HG13	2.40	0.46
1:E:131:LYS:NZ	2:F:128:LYS:HA	2.30	0.46
1:E:141:ARG:CB	1:E:141:ARG:NH1	2.79	0.46
1:E:314:LEU:O	1:E:318:VAL:HG23	2.16	0.46
1:G:378:GLN:NE2	1:G:380:PHE:CE1	2.83	0.46
1:I:411:PRO:HG3	1:I:422:TRP:CH2	2.50	0.46
1:I:414:GLY:CA	1:M:416:LYS:HZ3	2.29	0.46
1:I:415:ASN:O	1:I:417:LYS:HE2	2.15	0.46
1:K:244:LEU:HD12	1:K:267:CYS:SG	2.56	0.46
1:K:320:ARG:NH1	1:K:320:ARG:HB3	2.31	0.46
2:N:122:THR:O	2:N:126:MET:HG3	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:151:THR:HG22	1:A:174:ALA:HB3	1.98	0.46
1:A:157:LYS:NZ	1:A:430:LEU:HD13	2.31	0.46
1:A:196:GLN:HB3	1:A:221:GLN:HE21	1.81	0.46
1:A:415:ASN:C	1:A:417:LYS:H	2.24	0.46
2:F:122:THR:O	2:F:126:MET:HG3	2.15	0.46
2:H:24:SER:OG	2:H:110:LEU:HB3	2.15	0.46
1:I:205:ILE:O	1:I:206:GLU:C	2.59	0.46
1:I:367:GLU:C	1:I:369:GLY:H	2.24	0.46
1:K:420:GLU:O	1:K:421:ILE:C	2.58	0.46
1:M:109:TRP:C	1:M:111:SER:H	2.23	0.46
2:N:27:ILE:HG12	2:N:47:LEU:HD21	1.98	0.46
1:A:129:LEU:O	1:A:132:VAL:HG22	2.17	0.45
1:A:279:PHE:C	1:A:279:PHE:CD1	2.94	0.45
1:G:253:SER:HB2	1:G:280:ASP:CG	2.41	0.45
1:G:286:VAL:O	1:G:289:ALA:HB3	2.16	0.45
1:I:187:ALA:C	1:I:212:GLY:HA3	2.40	0.45
2:J:136:ARG:CZ	2:J:143:ASN:ND2	2.79	0.45
1:M:137:LYS:HE3	2:N:144:ASP:O	2.16	0.45
1:O:277:TRP:HZ2	1:O:331:ASP:OD2	2.00	0.45
1:O:278:CYS:C	1:O:280:ASP:H	2.23	0.45
2:P:22:LYS:C	2:P:24:SER:H	2.24	0.45
2:P:127:ILE:HG23	2:P:135:ILE:HD11	1.97	0.45
2:B:65:HIS:HB3	2:B:68:ASP:CB	2.26	0.45
1:C:237:THR:HG23	1:C:240:LYS:HE3	1.99	0.45
2:H:19:GLU:HA	2:H:19:GLU:OE2	2.16	0.45
1:I:159:LEU:N	1:I:181:PHE:O	2.42	0.45
1:I:276:SER:HB3	1:I:301:ASN:OD1	2.16	0.45
1:M:353:SER:O	1:M:354:LEU:HD23	2.17	0.45
1:O:278:CYS:HB2	1:O:281:PHE:CE2	2.51	0.45
1:A:185:PRO:O	1:A:186:LEU:HB2	2.17	0.45
1:E:109:TRP:C	1:E:111:SER:N	2.74	0.45
1:E:155:THR:HG22	1:E:178:PRO:CD	2.32	0.45
1:E:205:ILE:O	1:E:206:GLU:C	2.60	0.45
1:E:320:ARG:HB3	1:E:320:ARG:CZ	2.46	0.45
1:E:395:LEU:HD13	1:E:398:LEU:HD12	1.99	0.45
2:H:33:ASP:O	2:H:34:LEU:HD23	2.16	0.45
1:M:153:ASP:OD2	1:M:419:GLN:HB2	2.17	0.45
1:M:245:VAL:HG12	1:M:246:ARG:HG3	1.99	0.45
1:O:415:ASN:O	1:O:417:LYS:HE2	2.16	0.45
1:O:424:ILE:HD12	1:O:424:ILE:N	2.32	0.45
2:H:61:TRP:CD2	2:H:115:LEU:HD23	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:109:TRP:C	1:I:111:SER:H	2.24	0.45
1:I:184:GLN:HB3	1:I:185:PRO:HD2	1.96	0.45
1:K:149:TRP:O	1:K:172:VAL:HA	2.17	0.45
1:K:253:SER:HB2	1:K:280:ASP:CG	2.41	0.45
1:K:309:LEU:HD23	1:K:332:SER:OG	2.16	0.45
1:K:380:PHE:CE2	1:K:401:ASN:HB3	2.51	0.45
1:M:141:ARG:HB3	1:M:141:ARG:HH11	1.81	0.45
1:O:157:LYS:HG2	1:O:159:LEU:HD13	1.97	0.45
1:O:356:ARG:HG2	1:O:358:TYR:OH	2.16	0.45
2:P:55:LEU:HD22	2:P:59:ILE:HD11	1.96	0.45
1:E:416:LYS:CG	1:E:417:LYS:HE2	2.47	0.45
2:H:32:GLU:O	2:H:34:LEU:N	2.50	0.45
1:I:303:SER:HB2	1:I:330:SER:O	2.17	0.45
1:K:141:ARG:NH1	1:K:141:ARG:CB	2.79	0.45
2:L:61:TRP:CG	2:L:115:LEU:HD23	2.52	0.45
2:L:136:ARG:NH2	2:L:143:ASN:ND2	2.64	0.45
1:C:189:HIS:O	1:C:190:PHE:C	2.59	0.45
1:C:220:LEU:CB	1:C:241:ASN:ND2	2.77	0.45
1:G:153:ASP:OD2	1:G:419:GLN:HB2	2.16	0.45
1:G:411:PRO:HG3	1:G:422:TRP:CH2	2.52	0.45
1:I:420:GLU:O	1:I:421:ILE:C	2.60	0.45
1:C:367:GLU:C	1:C:369:GLY:H	2.24	0.45
1:C:417:LYS:HA	1:C:417:LYS:HD3	1.87	0.45
2:F:130:LYS:HB3	2:F:134:GLU:HB2	1.99	0.45
1:I:181:PHE:CG	1:K:384:PRO:HG3	2.50	0.45
1:K:149:TRP:CZ3	1:K:167:LEU:HD22	2.52	0.45
1:M:159:LEU:HD23	1:M:164:THR:CG2	2.42	0.45
1:M:419:GLN:O	1:M:425:LYS:HA	2.17	0.45
1:O:162:ASP:HA	1:O:190:PHE:CZ	2.37	0.45
2:P:98:GLY:O	2:P:102:GLU:HG2	2.17	0.45
2:P:130:LYS:HB3	2:P:134:GLU:HB2	1.98	0.45
2:B:30:MET:HB3	2:B:36:MET:HE2	1.98	0.45
1:E:161:PRO:HB3	1:E:209:THR:HG23	1.97	0.45
1:E:191:SER:HB3	1:E:192:PRO:HD2	1.99	0.45
1:G:378:GLN:NE2	1:G:380:PHE:HE1	2.14	0.45
1:G:415:ASN:C	1:G:417:LYS:H	2.25	0.45
1:K:109:TRP:C	1:K:111:SER:N	2.75	0.45
1:K:232:ASP:HB2	1:K:233:PRO:HD3	1.99	0.45
1:K:254:GLY:H	1:K:280:ASP:CG	2.24	0.45
2:L:55:LEU:HD22	2:L:59:ILE:HD11	1.98	0.45
1:M:109:TRP:HA	2:N:101:PHE:HE2	1.82	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:O:415:ASN:C	1:O:417:LYS:N	2.75	0.45
1:A:415:ASN:O	1:A:417:LYS:HE2	2.17	0.45
1:A:421:ILE:HG23	1:A:424:ILE:HB	1.99	0.45
1:C:232:ASP:HB2	1:C:233:PRO:HD3	1.98	0.45
1:C:272:GLU:HG2	1:C:299:GLN:HB2	1.98	0.45
2:D:136:ARG:CZ	2:D:143:ASN:ND2	2.80	0.45
1:G:378:GLN:HE22	1:G:401:ASN:HA	1.81	0.45
2:J:130:LYS:CE	2:J:138:THR:HG21	2.45	0.45
2:J:143:ASN:OD1	2:J:145:PHE:HB2	2.17	0.45
1:K:340:PHE:N	1:K:340:PHE:CD1	2.83	0.45
1:A:109:TRP:C	1:A:111:SER:H	2.23	0.45
1:A:159:LEU:N	1:A:181:PHE:O	2.47	0.45
1:A:274:ASN:C	1:A:274:ASN:ND2	2.75	0.45
1:A:278:CYS:HB2	1:A:281:PHE:CZ	2.52	0.45
1:A:320:ARG:HB2	1:A:320:ARG:HH11	1.82	0.45
2:D:133:GLU:HA	2:D:136:ARG:NH1	2.32	0.45
1:E:168:LEU:HD22	1:E:195:VAL:HG22	1.97	0.45
2:P:131:THR:OG1	2:P:134:GLU:HG3	2.17	0.45
1:A:132:VAL:O	1:A:135:VAL:HG23	2.16	0.44
1:A:331:ASP:CG	1:A:356:ARG:HH11	2.21	0.44
2:B:50:VAL:HG23	2:B:109:TYR:CD2	2.53	0.44
1:C:253:SER:HB2	1:C:280:ASP:CG	2.42	0.44
1:C:331:ASP:OD2	1:C:356:ARG:NH1	2.50	0.44
1:E:140:TYR:CE2	2:F:153:VAL:HG22	2.52	0.44
1:E:193:PHE:HB3	1:E:195:VAL:HG23	1.98	0.44
1:E:314:LEU:O	1:E:314:LEU:CD2	2.65	0.44
1:G:151:THR:HG22	1:G:174:ALA:HB3	1.99	0.44
1:G:164:THR:O	1:G:168:LEU:HG	2.17	0.44
1:K:159:LEU:CD2	1:K:164:THR:HG22	2.46	0.44
2:L:25:VAL:HG21	2:L:111:ASP:OD2	2.17	0.44
1:A:191:SER:O	1:A:192:PRO:C	2.61	0.44
1:A:253:SER:HB3	1:A:280:ASP:HB2	1.95	0.44
1:A:337:ASN:C	1:A:339:CYS:H	2.25	0.44
1:C:257:GLU:OE1	1:C:258:PHE:CE1	2.71	0.44
1:C:279:PHE:CD1	1:C:279:PHE:C	2.95	0.44
2:D:88:TRP:HE3	2:D:89:ASP:OD2	1.99	0.44
2:D:127:ILE:HG23	2:D:135:ILE:HD11	1.98	0.44
1:E:184:GLN:CB	1:E:185:PRO:CD	2.96	0.44
1:E:299:GLN:HA	1:E:326:HIS:HB2	1.99	0.44
1:E:389:GLN:HE21	1:E:392:LYS:HD2	1.82	0.44
1:G:241:ASN:O	1:G:244:LEU:HG	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:136:CYS:HB2	2:J:144:ASP:OD2	2.16	0.44
2:J:22:LYS:HG2	2:J:28:LYS:HZ2	1.82	0.44
1:M:424:ILE:HD12	1:M:424:ILE:N	2.32	0.44
1:O:244:LEU:HD12	1:O:267:CYS:SG	2.57	0.44
1:O:417:LYS:HD3	1:O:417:LYS:N	2.31	0.44
1:C:235:VAL:HG21	1:C:255:PHE:CG	2.52	0.44
1:C:335:LEU:HD22	1:C:339:CYS:SG	2.57	0.44
1:E:376:THR:HG22	1:E:399:GLN:HB2	1.98	0.44
1:G:205:ILE:O	1:G:206:GLU:C	2.60	0.44
1:I:185:PRO:O	1:I:186:LEU:HB2	2.17	0.44
1:K:400:ILE:O	1:K:401:ASN:C	2.61	0.44
1:A:210:LEU:HD12	1:A:210:LEU:HA	1.86	0.44
1:C:109:TRP:C	1:C:111:SER:N	2.74	0.44
1:C:417:LYS:NZ	1:G:417:LYS:HG3	2.29	0.44
2:D:16:VAL:CG1	2:D:59:ILE:HD13	2.47	0.44
1:E:136:CYS:C	2:F:145:PHE:CE1	2.95	0.44
1:K:281:PHE:HB2	1:K:285:HIS:CE1	2.51	0.44
1:K:390:LEU:HA	1:K:393:GLU:HG2	1.99	0.44
2:L:18:VAL:HG23	2:L:19:GLU:N	2.32	0.44
1:M:107:VAL:HG12	1:M:108:SER:N	2.31	0.44
1:M:421:ILE:HG23	1:M:421:ILE:O	2.17	0.44
2:N:20:ILE:O	2:N:23:GLN:HB2	2.17	0.44
1:O:420:GLU:O	1:O:422:TRP:N	2.50	0.44
2:P:58:VAL:HG11	2:P:110:LEU:HD12	1.98	0.44
1:A:320:ARG:HB3	1:A:320:ARG:CZ	2.47	0.44
1:A:367:GLU:C	1:A:369:GLY:N	2.76	0.44
1:C:109:TRP:CH2	2:D:101:PHE:HA	2.52	0.44
1:I:314:LEU:O	1:I:314:LEU:HD23	2.17	0.44
1:K:109:TRP:CH2	2:L:101:PHE:HA	2.51	0.44
1:K:160:HIS:CE1	1:K:161:PRO:HG2	2.52	0.44
2:N:16:VAL:HG23	2:N:21:ALA:HB2	2.00	0.44
1:O:153:ASP:HB2	1:O:426:CYS:HB2	1.99	0.44
1:O:157:LYS:HD2	1:O:430:LEU:HD12	1.99	0.44
1:O:184:GLN:CG	1:O:185:PRO:HD2	2.47	0.44
2:P:124:ALA:O	2:P:128:LYS:HG3	2.16	0.44
1:A:160:HIS:CE1	1:A:161:PRO:HG2	2.52	0.44
1:A:271:ASP:O	1:A:297:ILE:HD12	2.17	0.44
1:G:148:LEU:CA	1:G:427:ARG:HH22	2.23	0.44
2:J:12:GLU:OE2	2:J:52:ALA:HB1	2.17	0.44
1:K:176:ARG:NE	1:K:419:GLN:H	2.16	0.44
1:K:424:ILE:HD12	1:K:424:ILE:H	1.83	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:166:ARG:HH12	2:N:156:GLU:HA	1.83	0.44
1:M:340:PHE:N	1:M:340:PHE:CD1	2.83	0.44
1:A:182:MET:HB2	1:A:205:ILE:HG12	1.99	0.44
1:A:420:GLU:O	1:A:421:ILE:C	2.60	0.44
1:C:185:PRO:O	1:C:186:LEU:HB2	2.16	0.44
2:D:30:MET:C	2:D:32:GLU:N	2.76	0.44
1:E:419:GLN:HG3	1:E:426:CYS:HB2	2.00	0.44
1:G:320:ARG:HB3	1:G:320:ARG:CZ	2.47	0.44
2:H:133:GLU:HA	2:H:136:ARG:NH1	2.33	0.44
1:I:244:LEU:HD12	1:I:267:CYS:SG	2.58	0.44
1:I:314:LEU:O	1:I:318:VAL:HG23	2.18	0.44
1:I:415:ASN:C	1:I:417:LYS:N	2.74	0.44
1:M:126:LEU:HD23	1:M:430:LEU:HG	1.98	0.44
1:M:162:ASP:HA	1:M:190:PHE:CZ	2.53	0.44
1:M:215:SER:C	1:M:240:LYS:NZ	2.76	0.44
1:M:418:ASN:HD22	1:M:418:ASN:N	2.14	0.44
1:O:370:GLU:O	1:O:372:PRO:HD3	2.18	0.44
2:B:2:PRO:C	2:B:18:VAL:HG13	2.43	0.44
2:B:136:ARG:CZ	2:B:143:ASN:ND2	2.80	0.44
2:D:22:LYS:C	2:D:22:LYS:HD3	2.43	0.44
1:E:351:HIS:CD2	1:E:376:THR:OG1	2.61	0.44
1:E:390:LEU:HA	1:E:393:GLU:HG2	1.99	0.44
2:F:33:ASP:O	2:F:34:LEU:HD23	2.17	0.44
1:G:132:VAL:HA	1:G:135:VAL:HG23	2.00	0.44
1:G:186:LEU:O	1:G:188:GLU:N	2.51	0.44
1:I:129:LEU:O	1:I:132:VAL:HG22	2.17	0.44
1:K:126:LEU:HD23	1:K:430:LEU:HD21	1.99	0.44
2:L:136:ARG:CZ	2:L:143:ASN:HD22	2.30	0.44
1:C:274:ASN:C	1:C:274:ASN:ND2	2.65	0.44
1:I:109:TRP:C	1:I:111:SER:N	2.76	0.44
1:I:377:LEU:O	1:I:400:ILE:HA	2.18	0.44
2:J:54:ILE:HG23	2:J:103:LEU:HD23	1.99	0.44
1:K:140:TYR:HB2	2:L:145:PHE:HZ	1.82	0.44
1:K:299:GLN:HG2	1:K:326:HIS:HB2	2.00	0.44
1:M:419:GLN:HG3	1:M:426:CYS:HB2	2.00	0.44
1:O:237:THR:HG23	1:O:240:LYS:HE3	1.99	0.44
1:A:172:VAL:O	1:A:195:VAL:HA	2.18	0.43
1:A:184:GLN:CG	1:A:185:PRO:HD2	2.47	0.43
2:B:54:ILE:HG23	2:B:103:LEU:HD23	2.00	0.43
1:C:307:LYS:HD2	1:C:307:LYS:C	2.43	0.43
1:E:415:ASN:C	1:E:417:LYS:N	2.75	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:16:VAL:HG23	2:F:21:ALA:HB2	1.99	0.43
1:I:370:GLU:HA	1:I:370:GLU:OE1	2.18	0.43
2:J:49:ASN:HB2	2:J:109:TYR:CD2	2.53	0.43
2:J:85:ILE:HA	2:J:86:PRO:HD3	1.87	0.43
1:K:241:ASN:O	1:K:244:LEU:HG	2.18	0.43
1:M:320:ARG:CZ	1:M:320:ARG:HB3	2.47	0.43
1:O:151:THR:HB	1:O:426:CYS:HA	1.99	0.43
1:O:237:THR:CG2	1:O:240:LYS:HE3	2.48	0.43
1:A:414:GLY:CA	1:E:416:LYS:HZ3	2.30	0.43
1:C:246:ARG:HH11	1:C:246:ARG:HG3	1.82	0.43
2:D:64:HIS:HD2	2:D:65:HIS:CE1	2.37	0.43
1:E:128:GLU:CG	2:F:128:LYS:HG2	2.48	0.43
1:E:383:VAL:HA	1:E:384:PRO:HD3	1.90	0.43
2:F:49:ASN:HB2	2:F:109:TYR:CD2	2.53	0.43
1:G:311:LYS:HE3	1:G:338:ASP:O	2.17	0.43
1:G:418:ASN:HD22	1:G:418:ASN:N	2.16	0.43
1:I:419:GLN:HG3	1:I:426:CYS:HB2	2.00	0.43
2:L:6:LEU:O	2:L:13:ILE:HA	2.18	0.43
1:M:241:ASN:O	1:M:244:LEU:HG	2.18	0.43
2:N:25:VAL:HB	2:N:111:ASP:OD2	2.19	0.43
1:O:153:ASP:OD2	1:O:419:GLN:HB2	2.18	0.43
1:O:175:PHE:CD2	1:O:198:MET:HE2	2.48	0.43
1:O:415:ASN:O	1:O:416:LYS:CB	2.64	0.43
1:A:109:TRP:NE1	2:B:139:PHE:HD2	2.16	0.43
2:B:133:GLU:HA	2:B:136:ARG:NH1	2.33	0.43
1:E:417:LYS:HA	1:E:417:LYS:HD3	1.88	0.43
1:G:141:ARG:CB	1:G:141:ARG:NH1	2.81	0.43
1:G:382:ILE:HG13	1:G:383:VAL:HG23	2.00	0.43
1:I:137:LYS:HE2	2:J:149:GLU:OE2	2.18	0.43
1:I:161:PRO:HB3	1:I:209:THR:CG2	2.48	0.43
1:K:331:ASP:CG	1:K:356:ARG:NH1	2.75	0.43
1:K:417:LYS:HG3	1:O:417:LYS:HZ2	1.82	0.43
2:L:49:ASN:HB2	2:L:109:TYR:CD2	2.53	0.43
1:O:420:GLU:O	1:O:420:GLU:HG3	2.18	0.43
2:B:55:LEU:HD22	2:B:59:ILE:CD1	2.48	0.43
2:D:133:GLU:HA	2:D:136:ARG:HH11	1.83	0.43
1:I:416:LYS:HD2	1:I:417:LYS:CE	2.47	0.43
1:K:367:GLU:C	1:K:369:GLY:N	2.77	0.43
2:L:24:SER:HB3	2:L:27:ILE:HB	2.00	0.43
1:M:271:ASP:C	1:M:297:ILE:HD12	2.44	0.43
1:M:378:GLN:NE2	1:M:380:PHE:CE1	2.86	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:N:8:SER:C	2:N:10:ASP:N	2.76	0.43
2:P:18:VAL:HG23	2:P:19:GLU:N	2.34	0.43
2:P:130:LYS:CE	2:P:138:THR:HG21	2.47	0.43
2:B:50:VAL:HG23	2:B:109:TYR:CE2	2.54	0.43
1:G:129:LEU:O	1:G:132:VAL:HG22	2.18	0.43
1:G:278:CYS:HB2	1:G:281:PHE:CZ	2.54	0.43
2:H:30:MET:HB3	2:H:36:MET:CE	2.49	0.43
2:H:43:ASP:CB	2:H:44:PRO:CD	2.89	0.43
1:I:331:ASP:CG	1:I:356:ARG:NH1	2.76	0.43
1:I:416:LYS:CD	1:I:417:LYS:NZ	2.69	0.43
1:K:417:LYS:NZ	1:O:417:LYS:HG3	2.31	0.43
2:L:25:VAL:HB	2:L:111:ASP:HB3	2.01	0.43
1:M:278:CYS:HB2	1:M:281:PHE:CE2	2.54	0.43
2:P:8:SER:C	2:P:10:ASP:H	2.27	0.43
1:A:379:VAL:N	1:A:401:ASN:OD1	2.41	0.43
1:A:420:GLU:O	1:A:423:GLY:N	2.51	0.43
1:C:399:GLN:HE21	1:C:399:GLN:CA	2.26	0.43
1:E:411:PRO:HG3	1:E:422:TRP:CH2	2.53	0.43
2:F:61:TRP:HB2	2:F:92:PHE:CZ	2.54	0.43
1:K:420:GLU:O	1:K:423:GLY:N	2.50	0.43
1:M:149:TRP:CZ3	1:M:167:LEU:HD22	2.53	0.43
1:M:220:LEU:CB	1:M:241:ASN:HD22	2.31	0.43
1:O:178:PRO:O	1:O:179:ARG:HB2	2.18	0.43
2:P:133:GLU:HA	2:P:133:GLU:OE2	2.19	0.43
1:A:153:ASP:OD2	1:A:419:GLN:HB2	2.19	0.43
1:A:299:GLN:HG2	1:A:326:HIS:HB2	2.01	0.43
1:A:418:ASN:HD22	1:A:418:ASN:N	2.16	0.43
1:C:128:GLU:OE2	2:D:128:LYS:HA	2.19	0.43
1:C:215:SER:C	1:C:240:LYS:NZ	2.77	0.43
1:E:132:VAL:O	1:E:135:VAL:HG23	2.19	0.43
1:E:278:CYS:C	1:E:280:ASP:N	2.76	0.43
1:G:314:LEU:O	1:G:314:LEU:HD23	2.18	0.43
1:G:320:ARG:HB2	1:G:320:ARG:NH1	2.33	0.43
1:G:394:ALA:O	1:G:396:PRO:HD2	2.18	0.43
1:I:380:PHE:CE2	1:I:401:ASN:HB3	2.54	0.43
1:O:416:LYS:CD	1:O:417:LYS:NZ	2.72	0.43
2:P:24:SER:OG	2:P:110:LEU:HB3	2.18	0.43
2:P:54:ILE:HG13	2:P:102:GLU:HB3	2.01	0.43
1:A:416:LYS:CD	1:A:417:LYS:HZ1	2.08	0.43
1:C:380:PHE:CE2	1:C:401:ASN:HB3	2.54	0.43
1:G:415:ASN:C	1:G:417:LYS:N	2.77	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:22:LYS:HG2	2:H:28:LYS:HZ2	1.83	0.43
1:I:125:CYS:SG	1:I:127:PRO:HG2	2.59	0.43
2:L:104:ILE:HD11	2:L:123:VAL:HG21	2.00	0.43
2:L:127:ILE:HG23	2:L:135:ILE:HD11	2.00	0.43
1:M:307:LYS:HD2	1:M:307:LYS:C	2.44	0.43
2:N:43:ASP:CB	2:N:44:PRO:CD	2.91	0.43
1:O:299:GLN:HG2	1:O:326:HIS:ND1	2.34	0.43
2:P:55:LEU:HD22	2:P:59:ILE:CD1	2.48	0.43
1:A:244:LEU:HD12	1:A:267:CYS:SG	2.58	0.43
1:A:389:GLN:HE22	1:A:392:LYS:NZ	2.16	0.43
2:D:133:GLU:HG3	2:D:137:LYS:HE2	2.01	0.43
1:E:157:LYS:HZ2	1:E:430:LEU:HD13	1.82	0.43
1:E:419:GLN:O	1:E:425:LYS:HA	2.19	0.43
1:I:420:GLU:O	1:I:422:TRP:N	2.52	0.43
1:K:141:ARG:HH11	1:K:141:ARG:CB	2.31	0.43
1:K:300:LEU:HD23	1:K:317:LEU:HD11	2.01	0.43
2:L:16:VAL:HG23	2:L:21:ALA:HB2	2.00	0.43
1:M:184:GLN:CG	1:M:185:PRO:HD2	2.49	0.43
1:M:370:GLU:OE1	1:M:370:GLU:HA	2.19	0.43
2:N:127:ILE:HG23	2:N:135:ILE:CD1	2.49	0.43
1:O:176:ARG:NE	1:O:419:GLN:H	2.16	0.43
2:P:26:THR:HB	2:P:110:LEU:HA	2.00	0.43
1:A:151:THR:CB	1:A:426:CYS:HA	2.49	0.43
1:C:274:ASN:ND2	1:C:276:SER:H	2.16	0.43
1:E:130:LEU:CD1	2:F:160:CYS:HB2	2.48	0.43
1:E:232:ASP:O	1:E:236:ASN:ND2	2.52	0.43
2:F:30:MET:C	2:F:32:GLU:N	2.76	0.43
1:K:294:SER:C	1:K:296:THR:H	2.26	0.43
2:L:43:ASP:CB	2:L:44:PRO:CD	2.87	0.43
1:M:109:TRP:C	1:M:111:SER:N	2.76	0.43
1:M:420:GLU:HG3	1:M:423:GLY:HA2	2.00	0.43
2:N:117:ASP:OD1	2:N:121:LYS:HE3	2.19	0.43
1:O:131:LYS:NZ	2:P:127:ILE:O	2.51	0.43
1:O:307:LYS:HD2	1:O:307:LYS:C	2.44	0.43
1:O:419:GLN:HG3	1:O:426:CYS:HB2	2.01	0.43
1:A:137:LYS:HE2	2:B:149:GLU:OE2	2.18	0.42
1:A:384:PRO:HD2	1:A:387:THR:OG1	2.19	0.42
2:B:64:HIS:CD2	2:B:88:TRP:CE3	3.07	0.42
2:D:6:LEU:O	2:D:13:ILE:HA	2.19	0.42
1:E:383:VAL:HG11	1:E:388:LEU:HD13	2.01	0.42
2:F:155:LYS:HA	2:F:158:GLN:HG3	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:274:ASN:HB2	1:G:408:ILE:HG22	2.01	0.42
2:J:131:THR:HG23	2:J:134:GLU:OE2	2.18	0.42
1:M:271:ASP:O	1:M:297:ILE:HD12	2.19	0.42
2:N:124:ALA:O	2:N:128:LYS:HG3	2.19	0.42
1:O:141:ARG:HB3	1:O:141:ARG:NH1	2.34	0.42
1:A:272:GLU:HG2	1:A:299:GLN:HB2	2.00	0.42
1:A:320:ARG:NH1	1:A:320:ARG:CB	2.82	0.42
2:B:143:ASN:OD1	2:B:145:PHE:HB2	2.19	0.42
1:C:124:LEU:HD23	2:D:124:ALA:HB1	2.01	0.42
1:C:278:CYS:HB2	1:C:281:PHE:CE2	2.54	0.42
1:E:159:LEU:N	1:E:181:PHE:O	2.44	0.42
1:E:384:PRO:HG3	1:G:181:PHE:CD2	2.54	0.42
1:E:420:GLU:O	1:E:422:TRP:N	2.52	0.42
1:G:130:LEU:HD12	2:H:160:CYS:SG	2.59	0.42
1:G:416:LYS:C	1:G:417:LYS:CD	2.90	0.42
1:K:370:GLU:OE1	1:K:370:GLU:HA	2.18	0.42
1:K:378:GLN:HE22	1:K:401:ASN:HA	1.81	0.42
1:K:416:LYS:HD2	1:K:417:LYS:HZ3	1.83	0.42
2:L:30:MET:HB3	2:L:36:MET:HE2	2.01	0.42
1:M:399:GLN:HE21	1:M:399:GLN:CA	2.25	0.42
1:C:420:GLU:O	1:C:422:TRP:N	2.52	0.42
1:K:215:SER:C	1:K:240:LYS:NZ	2.78	0.42
1:O:307:LYS:HE3	1:O:308:ASN:OD1	2.19	0.42
1:A:109:TRP:C	1:A:111:SER:N	2.76	0.42
2:B:24:SER:HB3	2:B:27:ILE:HB	2.01	0.42
1:E:185:PRO:O	1:E:186:LEU:HB2	2.18	0.42
2:F:18:VAL:HG23	2:F:19:GLU:N	2.34	0.42
2:F:131:THR:O	2:F:135:ILE:HG13	2.18	0.42
1:G:245:VAL:HG12	1:G:246:ARG:HG3	2.01	0.42
1:G:415:ASN:O	1:G:417:LYS:HE2	2.19	0.42
2:H:24:SER:HB3	2:H:27:ILE:HB	2.02	0.42
1:I:281:PHE:CE1	1:I:305:TYR:CE2	3.08	0.42
2:J:115:LEU:HD13	2:J:115:LEU:C	2.44	0.42
1:K:271:ASP:C	1:K:297:ILE:HD12	2.45	0.42
1:K:379:VAL:HG23	1:K:379:VAL:O	2.19	0.42
1:M:162:ASP:OD2	1:M:186:LEU:HD12	2.20	0.42
1:M:272:GLU:HG2	1:M:299:GLN:HB2	2.00	0.42
2:N:33:ASP:O	2:N:34:LEU:HD23	2.20	0.42
1:O:416:LYS:CD	1:O:417:LYS:HZ1	2.05	0.42
2:P:45:VAL:HA	2:P:46:PRO:HD3	1.88	0.42
2:P:149:GLU:O	2:P:153:VAL:HG23	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:164:THR:O	1:A:168:LEU:HG	2.19	0.42
1:A:187:ALA:O	1:A:188:GLU:O	2.38	0.42
1:A:415:ASN:O	1:A:416:LYS:CB	2.66	0.42
2:B:132:PRO:O	2:B:136:ARG:HG3	2.19	0.42
1:C:416:LYS:CG	1:C:417:LYS:HE2	2.49	0.42
2:D:24:SER:HB3	2:D:27:ILE:HB	2.01	0.42
2:D:47:LEU:HD12	2:D:55:LEU:HD11	2.00	0.42
2:F:130:LYS:CE	2:F:138:THR:HG21	2.50	0.42
1:G:153:ASP:OD2	1:G:419:GLN:HG3	2.19	0.42
1:I:151:THR:HB	1:I:426:CYS:HA	2.01	0.42
1:I:164:THR:HG21	1:I:182:MET:HE2	2.01	0.42
1:K:131:LYS:NZ	2:L:127:ILE:O	2.52	0.42
1:K:277:TRP:HZ2	1:K:331:ASP:OD2	2.03	0.42
1:K:394:ALA:O	1:K:396:PRO:HD2	2.20	0.42
1:M:191:SER:HB3	1:M:192:PRO:CD	2.45	0.42
1:M:231:SER:OG	1:M:233:PRO:HD2	2.19	0.42
1:M:278:CYS:HB2	1:M:281:PHE:CZ	2.55	0.42
2:N:8:SER:C	2:N:10:ASP:H	2.26	0.42
2:N:102:GLU:HA	2:N:102:GLU:OE2	2.20	0.42
1:O:378:GLN:HE21	1:O:380:PHE:HE1	1.68	0.42
2:B:30:MET:C	2:B:32:GLU:H	2.28	0.42
1:C:314:LEU:O	1:C:318:VAL:HG23	2.19	0.42
1:E:272:GLU:HG2	1:E:299:GLN:HB2	2.00	0.42
1:E:284:LYS:O	1:E:288:VAL:HG23	2.19	0.42
2:F:2:PRO:C	2:F:18:VAL:HG13	2.45	0.42
1:I:279:PHE:CD1	1:I:279:PHE:C	2.98	0.42
1:I:378:GLN:HE21	1:I:380:PHE:HE1	1.67	0.42
1:K:418:ASN:HD22	1:K:418:ASN:N	2.17	0.42
1:K:419:GLN:HG3	1:K:426:CYS:HB2	2.00	0.42
2:L:25:VAL:CB	2:L:111:ASP:OD2	2.67	0.42
1:O:376:THR:HG22	1:O:399:GLN:HB2	2.00	0.42
1:A:199:ASP:OD1	1:A:411:PRO:HG2	2.20	0.42
1:A:232:ASP:O	1:A:233:PRO:C	2.59	0.42
1:C:389:GLN:HE22	1:C:392:LYS:NZ	2.16	0.42
1:C:421:ILE:HG23	1:C:424:ILE:HB	2.02	0.42
1:E:314:LEU:O	1:E:314:LEU:HD23	2.20	0.42
1:G:132:VAL:CG1	2:H:127:ILE:HG21	2.50	0.42
2:H:18:VAL:HG23	2:H:19:GLU:N	2.34	0.42
1:O:383:VAL:CG1	1:O:388:LEU:HB2	2.48	0.42
2:P:148:GLU:O	2:P:151:ALA:HB3	2.18	0.42
1:C:141:ARG:NH1	1:C:141:ARG:CB	2.83	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:420:GLU:O	1:C:420:GLU:HG3	2.20	0.42
2:F:45:VAL:HA	2:F:46:PRO:HD3	1.81	0.42
1:G:329:LEU:O	1:G:332:SER:OG	2.32	0.42
1:G:421:ILE:HG23	1:G:424:ILE:HB	2.02	0.42
1:I:421:ILE:HG23	1:I:421:ILE:O	2.18	0.42
2:J:101:PHE:C	2:J:101:PHE:CD1	2.98	0.42
1:K:274:ASN:C	1:K:274:ASN:ND2	2.77	0.42
1:K:376:THR:HG22	1:K:399:GLN:HB3	2.02	0.42
2:N:100:LEU:O	2:N:104:ILE:HG13	2.20	0.42
1:O:109:TRP:CZ2	2:P:101:PHE:HB2	2.55	0.42
1:O:125:CYS:SG	1:O:127:PRO:HG2	2.60	0.42
1:O:132:VAL:HG12	2:P:127:ILE:HG21	2.01	0.42
2:P:30:MET:C	2:P:32:GLU:H	2.28	0.42
1:A:340:PHE:N	1:A:340:PHE:HD1	2.18	0.42
1:C:177:CYS:HB3	1:C:200:LEU:HD23	2.02	0.42
2:D:130:LYS:CE	2:D:138:THR:HG21	2.47	0.42
1:E:231:SER:CB	1:E:233:PRO:HD2	2.49	0.42
2:F:8:SER:C	2:F:10:ASP:H	2.27	0.42
1:I:299:GLN:HG2	1:I:326:HIS:HB2	2.02	0.42
1:I:330:SER:O	1:I:331:ASP:HB2	2.20	0.42
1:I:382:ILE:HG13	1:I:383:VAL:HG23	2.01	0.42
2:J:49:ASN:ND2	2:J:109:TYR:CE1	2.86	0.42
1:M:254:GLY:H	1:M:280:ASP:CG	2.28	0.42
1:M:299:GLN:HG2	1:M:326:HIS:HB2	2.01	0.42
1:O:253:SER:HB2	1:O:280:ASP:CG	2.45	0.42
1:O:320:ARG:HB3	1:O:320:ARG:CZ	2.50	0.42
1:O:389:GLN:HE22	1:O:392:LYS:NZ	2.18	0.42
1:C:374:LEU:HG	1:C:398:LEU:CD2	2.50	0.42
2:H:22:LYS:HD3	2:H:22:LYS:C	2.45	0.42
2:J:16:VAL:HG12	2:J:59:ILE:HG21	2.02	0.42
2:J:148:GLU:O	2:J:151:ALA:HB3	2.19	0.42
2:L:64:HIS:HD2	2:L:65:HIS:CE1	2.38	0.42
1:C:186:LEU:HD23	1:C:186:LEU:HA	1.91	0.41
2:D:30:MET:HB3	2:D:36:MET:HE2	2.02	0.41
2:D:36:MET:HE3	2:D:36:MET:HB2	1.88	0.41
2:D:146:THR:HB	2:D:149:GLU:CG	2.48	0.41
1:E:151:THR:HG22	1:E:174:ALA:HB3	2.02	0.41
1:E:243:ASN:HA	1:E:269:ARG:HD2	2.02	0.41
1:G:130:LEU:HD11	2:H:159:TRP:CZ2	2.55	0.41
2:H:8:SER:C	2:H:10:ASP:N	2.77	0.41
2:H:65:HIS:CB	2:H:68:ASP:HB2	2.37	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:220:LEU:HD13	1:K:223:LEU:HD22	2.02	0.41
1:K:351:HIS:CD2	1:K:376:THR:OG1	2.65	0.41
1:K:420:GLU:O	1:K:420:GLU:HG3	2.20	0.41
1:O:197:HIS:CD2	1:O:222:ASN:ND2	2.85	0.41
1:O:215:SER:C	1:O:240:LYS:NZ	2.77	0.41
2:P:33:ASP:O	2:P:34:LEU:CD2	2.68	0.41
2:P:55:LEU:O	2:P:59:ILE:HG13	2.20	0.41
2:P:85:ILE:HG22	2:P:90:GLN:HG3	2.01	0.41
1:A:168:LEU:HD22	1:A:195:VAL:CG2	2.48	0.41
1:A:351:HIS:CD2	1:A:376:THR:OG1	2.65	0.41
2:B:136:ARG:O	2:B:140:ASN:HA	2.20	0.41
1:E:137:LYS:HE2	2:F:149:GLU:CD	2.45	0.41
1:K:185:PRO:O	1:K:186:LEU:HB2	2.20	0.41
1:K:196:GLN:HB3	1:K:221:GLN:HE21	1.86	0.41
1:M:205:ILE:O	1:M:206:GLU:C	2.62	0.41
1:O:141:ARG:NH1	1:O:141:ARG:CB	2.83	0.41
1:O:184:GLN:HG2	1:O:185:PRO:HD2	2.02	0.41
1:O:419:GLN:O	1:O:425:LYS:HA	2.20	0.41
1:A:116:LEU:HD21	2:B:108:ASN:HB2	2.01	0.41
1:A:390:LEU:HA	1:A:393:GLU:HG2	2.01	0.41
1:A:420:GLU:O	1:A:420:GLU:HG3	2.20	0.41
2:B:127:ILE:HG23	2:B:135:ILE:HD11	2.01	0.41
1:C:394:ALA:O	1:C:396:PRO:HD2	2.21	0.41
1:E:416:LYS:CD	1:E:417:LYS:NZ	2.75	0.41
1:I:378:GLN:HE22	1:I:401:ASN:HA	1.85	0.41
1:I:420:GLU:O	1:I:420:GLU:HG3	2.20	0.41
2:J:30:MET:O	2:J:34:LEU:HB2	2.21	0.41
1:K:235:VAL:HG21	1:K:255:PHE:CB	2.50	0.41
1:K:327:LEU:O	1:K:352:LEU:HD12	2.21	0.41
2:L:55:LEU:HD22	2:L:59:ILE:CD1	2.49	0.41
1:M:157:LYS:HD2	1:M:430:LEU:HD12	2.02	0.41
1:M:164:THR:HG21	1:M:182:MET:CE	2.51	0.41
2:N:158:GLN:O	2:N:160:CYS:N	2.53	0.41
1:O:410:ARG:NH1	1:O:410:ARG:HG3	2.35	0.41
1:A:207:VAL:O	1:A:234:ILE:HD11	2.20	0.41
1:C:211:HIS:HB2	1:C:234:ILE:HG12	2.01	0.41
1:C:390:LEU:HA	1:C:393:GLU:CG	2.50	0.41
1:E:199:ASP:OD1	1:E:411:PRO:HG2	2.20	0.41
2:F:136:ARG:O	2:F:140:ASN:HA	2.20	0.41
1:G:109:TRP:C	1:G:111:SER:H	2.27	0.41
2:J:64:HIS:HD2	2:J:65:HIS:CE1	2.38	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:229:ARG:CZ	1:K:253:SER:OG	2.68	0.41
1:K:294:SER:C	1:K:296:THR:N	2.79	0.41
2:L:45:VAL:HA	2:L:46:PRO:HD3	1.86	0.41
1:M:281:PHE:HB2	1:M:285:HIS:CE1	2.56	0.41
2:N:25:VAL:HG21	2:N:111:ASP:OD2	2.20	0.41
1:O:149:TRP:O	1:O:172:VAL:HG22	2.20	0.41
1:O:390:LEU:HA	1:O:393:GLU:CG	2.51	0.41
2:P:2:PRO:C	2:P:18:VAL:HG13	2.46	0.41
1:A:415:ASN:C	1:A:417:LYS:N	2.78	0.41
2:D:61:TRP:CZ3	2:D:115:LEU:HB2	2.56	0.41
2:D:65:HIS:HD2	2:D:68:ASP:OD2	2.03	0.41
2:F:54:ILE:O	2:F:58:VAL:HG23	2.20	0.41
2:F:149:GLU:O	2:F:153:VAL:HG23	2.19	0.41
2:H:28:LYS:HG2	2:H:32:GLU:OE2	2.21	0.41
2:H:29:THR:HG23	2:H:30:MET:N	2.36	0.41
2:L:8:SER:C	2:L:10:ASP:N	2.79	0.41
2:L:146:THR:HB	2:L:149:GLU:CG	2.47	0.41
1:M:246:ARG:HG2	1:M:272:GLU:HB2	2.01	0.41
2:N:30:MET:C	2:N:32:GLU:H	2.27	0.41
1:O:191:SER:HB3	1:O:192:PRO:HD2	2.01	0.41
1:A:137:LYS:CG	2:B:144:ASP:HB2	2.47	0.41
2:B:61:TRP:CD2	2:B:115:LEU:HD23	2.55	0.41
1:G:196:GLN:HB3	1:G:221:GLN:HE21	1.86	0.41
1:I:417:LYS:NZ	1:M:417:LYS:HG3	2.35	0.41
2:J:133:GLU:HG3	2:J:137:LYS:HE2	2.01	0.41
1:K:390:LEU:O	1:K:393:GLU:HB2	2.21	0.41
1:M:309:LEU:HD23	1:M:332:SER:OG	2.20	0.41
2:N:133:GLU:HA	2:N:136:ARG:NH1	2.35	0.41
1:O:132:VAL:C	1:O:134:GLY:N	2.77	0.41
1:O:378:GLN:HE22	1:O:402:CYS:N	2.19	0.41
2:P:19:GLU:HA	2:P:19:GLU:OE2	2.21	0.41
2:D:112:ILE:O	2:D:112:ILE:HG22	2.20	0.41
2:F:116:LEU:O	2:F:120:CYS:SG	2.71	0.41
2:H:30:MET:C	2:H:32:GLU:H	2.29	0.41
1:I:176:ARG:NH2	1:I:418:ASN:HB2	2.36	0.41
1:I:253:SER:HB3	1:I:280:ASP:HB2	1.96	0.41
1:K:205:ILE:O	1:K:206:GLU:C	2.62	0.41
1:M:193:PHE:HB3	1:M:195:VAL:HG23	2.03	0.41
1:O:159:LEU:HD23	1:O:164:THR:CG2	2.47	0.41
2:P:88:TRP:HA	2:P:91:GLU:HG3	2.02	0.41
1:A:159:LEU:HD23	1:A:164:THR:CG2	2.48	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:89:ASP:O	2:B:92:PHE:HB3	2.20	0.41
1:C:300:LEU:HD23	1:C:317:LEU:HD11	2.02	0.41
1:E:341:GLN:HA	1:E:344:PHE:CD1	2.55	0.41
1:E:384:PRO:HA	1:G:181:PHE:CE1	2.56	0.41
2:F:49:ASN:HB2	2:F:109:TYR:CZ	2.55	0.41
1:G:415:ASN:O	1:G:416:LYS:CB	2.66	0.41
2:J:61:TRP:CD1	2:J:115:LEU:HD23	2.56	0.41
1:K:124:LEU:HD23	2:L:124:ALA:HB1	2.02	0.41
1:K:182:MET:HB2	1:K:205:ILE:HG23	2.02	0.41
1:K:341:GLN:HA	1:K:344:PHE:CD1	2.56	0.41
1:K:421:ILE:O	1:K:421:ILE:HG23	2.21	0.41
2:L:136:ARG:NE	2:L:143:ASN:ND2	2.69	0.41
1:M:354:LEU:O	1:M:357:CYS:SG	2.79	0.41
2:N:54:ILE:O	2:N:57:LYS:HB2	2.21	0.41
1:O:159:LEU:N	1:O:181:PHE:O	2.42	0.41
1:O:367:GLU:C	1:O:369:GLY:N	2.78	0.41
2:P:22:LYS:C	2:P:24:SER:N	2.78	0.41
2:B:101:PHE:CD1	2:B:101:PHE:C	2.98	0.41
1:C:254:GLY:H	1:C:280:ASP:CG	2.29	0.41
1:C:356:ARG:HG2	1:C:358:TYR:OH	2.21	0.41
1:C:417:LYS:HE3	1:G:417:LYS:CD	2.49	0.41
2:F:85:ILE:HA	2:F:86:PRO:HD3	1.94	0.41
1:G:135:VAL:O	2:H:136:ARG:NE	2.54	0.41
1:G:185:PRO:O	1:G:186:LEU:HB2	2.21	0.41
1:G:279:PHE:CD1	1:G:279:PHE:C	2.99	0.41
1:I:194:ARG:O	1:I:195:VAL:C	2.63	0.41
2:J:33:ASP:O	2:J:34:LEU:HD23	2.20	0.41
2:L:54:ILE:HG13	2:L:102:GLU:HB3	2.01	0.41
1:M:378:GLN:NE2	1:M:380:PHE:HE1	2.19	0.41
1:O:164:THR:HG21	1:O:182:MET:CE	2.51	0.41
2:P:6:LEU:O	2:P:13:ILE:HA	2.21	0.41
1:A:210:LEU:HB3	1:A:234:ILE:CD1	2.51	0.41
1:A:297:ILE:HD12	1:A:297:ILE:HA	1.97	0.41
2:B:45:VAL:HA	2:B:46:PRO:HD3	1.81	0.41
1:C:140:TYR:HB2	2:D:145:PHE:HZ	1.85	0.41
1:C:162:ASP:OD2	1:C:186:LEU:HD12	2.21	0.41
1:C:370:GLU:HA	1:C:370:GLU:OE1	2.21	0.41
1:E:157:LYS:HG2	1:E:159:LEU:HD13	2.02	0.41
1:E:203:SER:O	1:E:228:LEU:CD2	2.69	0.41
1:E:237:THR:HG23	1:E:240:LYS:HE3	2.03	0.41
1:E:421:ILE:HG23	1:E:424:ILE:HB	2.04	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:235:VAL:HG21	1:G:255:PHE:CB	2.51	0.41
1:G:278:CYS:CB	1:G:281:PHE:CE2	3.04	0.41
2:H:4:ILE:HG21	2:H:31:LEU:HD11	2.02	0.41
1:I:247:LEU:HD21	1:I:263:LEU:HD21	2.02	0.41
1:I:390:LEU:HA	1:I:393:GLU:CG	2.51	0.41
2:L:43:ASP:O	2:L:45:VAL:N	2.54	0.41
1:A:149:TRP:CZ3	1:A:167:LEU:HD22	2.56	0.40
1:A:194:ARG:O	1:A:195:VAL:C	2.64	0.40
1:C:367:GLU:C	1:C:369:GLY:N	2.79	0.40
1:E:131:LYS:HZ1	2:F:128:LYS:HA	1.86	0.40
1:E:141:ARG:HH11	1:E:141:ARG:CB	2.34	0.40
1:G:133:SER:O	1:G:140:TYR:HB2	2.20	0.40
1:I:137:LYS:H	2:J:144:ASP:HB2	1.85	0.40
1:I:141:ARG:HH11	1:I:141:ARG:HB3	1.85	0.40
2:L:133:GLU:HA	2:L:136:ARG:NH1	2.36	0.40
1:M:220:LEU:CB	1:M:241:ASN:ND2	2.82	0.40
1:M:337:ASN:C	1:M:339:CYS:H	2.28	0.40
2:N:136:ARG:O	2:N:140:ASN:HA	2.20	0.40
1:O:382:ILE:HG13	1:O:383:VAL:HG23	2.03	0.40
1:O:410:ARG:HG3	1:O:410:ARG:HH11	1.86	0.40
1:A:363:GLU:C	1:A:365:LEU:N	2.79	0.40
1:A:399:GLN:CG	1:A:402:CYS:SG	3.10	0.40
1:C:157:LYS:HG2	1:C:159:LEU:HD13	2.04	0.40
1:C:253:SER:HB3	1:C:280:ASP:HB2	1.95	0.40
1:E:181:PHE:CE1	1:G:384:PRO:HG3	2.57	0.40
1:G:176:ARG:NE	1:G:419:GLN:HA	2.36	0.40
1:I:109:TRP:HZ3	2:J:104:ILE:HD12	1.86	0.40
1:I:384:PRO:HD2	1:I:387:THR:OG1	2.21	0.40
2:J:45:VAL:HA	2:J:46:PRO:HD3	1.82	0.40
1:K:126:LEU:HD23	1:K:430:LEU:CG	2.52	0.40
2:L:19:GLU:OE2	2:L:19:GLU:HA	2.21	0.40
2:N:30:MET:C	2:N:32:GLU:N	2.79	0.40
1:O:421:ILE:O	1:O:421:ILE:HG23	2.21	0.40
1:A:109:TRP:HE1	2:B:97:GLN:HB3	1.86	0.40
1:A:151:THR:CG2	1:A:424:ILE:HG21	2.52	0.40
1:C:109:TRP:CE2	2:D:101:PHE:HD2	2.38	0.40
1:C:197:HIS:CE1	1:C:221:GLN:HE21	2.39	0.40
2:D:43:ASP:CB	2:D:44:PRO:CD	2.93	0.40
1:E:137:LYS:HA	2:F:145:PHE:CZ	2.56	0.40
1:G:109:TRP:C	1:G:111:SER:N	2.79	0.40
1:G:163:VAL:HG11	2:H:159:TRP:CG	2.57	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:281:PHE:CD2	1:I:285:HIS:CD2	3.09	0.40
1:K:320:ARG:CB	1:K:320:ARG:CZ	2.99	0.40
1:K:329:LEU:O	1:K:332:SER:OG	2.32	0.40
2:L:136:ARG:O	2:L:140:ASN:HA	2.21	0.40
1:M:253:SER:HB2	1:M:280:ASP:CG	2.46	0.40
1:M:299:GLN:HG2	1:M:326:HIS:ND1	2.36	0.40
1:M:415:ASN:O	1:M:416:LYS:CB	2.68	0.40
2:N:134:GLU:HA	2:N:137:LYS:HD2	2.03	0.40
1:O:109:TRP:CD2	2:P:101:PHE:HD2	2.40	0.40
1:A:116:LEU:CD1	2:B:105:LEU:HD23	2.52	0.40
1:A:141:ARG:CB	1:A:141:ARG:NH1	2.84	0.40
1:A:325:VAL:CG1	1:A:350:GLN:HG2	2.50	0.40
1:A:416:LYS:C	1:A:417:LYS:CD	2.92	0.40
2:B:130:LYS:HB3	2:B:134:GLU:HB2	2.03	0.40
1:C:141:ARG:HB3	1:C:141:ARG:NH1	2.37	0.40
1:E:137:LYS:HE2	2:F:149:GLU:OE1	2.21	0.40
1:E:247:LEU:HD21	1:E:263:LEU:HD21	2.02	0.40
1:E:379:VAL:O	1:E:379:VAL:HG23	2.22	0.40
1:G:141:ARG:HB3	1:G:141:ARG:NH1	2.36	0.40
2:H:45:VAL:HA	2:H:46:PRO:HD3	1.85	0.40
2:J:50:VAL:HG23	2:J:109:TYR:CD2	2.56	0.40
2:J:133:GLU:HA	2:J:136:ARG:HH11	1.87	0.40
1:K:129:LEU:O	1:K:132:VAL:HG22	2.22	0.40
1:K:179:ARG:HA	1:K:202:ASN:O	2.22	0.40
1:K:181:PHE:HA	1:K:204:VAL:O	2.21	0.40
1:K:399:GLN:CG	1:K:402:CYS:SG	3.10	0.40
2:L:88:TRP:O	2:L:91:GLU:HG3	2.21	0.40
2:N:54:ILE:O	2:N:58:VAL:HG23	2.22	0.40
2:N:131:THR:OG1	2:N:134:GLU:HG3	2.21	0.40
2:N:131:THR:O	2:N:135:ILE:HG13	2.21	0.40
1:O:260:LEU:O	1:O:264:LEU:HG	2.22	0.40
2:B:122:THR:O	2:B:126:MET:HG3	2.20	0.40
1:E:125:CYS:SG	1:E:127:PRO:HG2	2.62	0.40
1:E:129:LEU:O	1:E:132:VAL:HG22	2.21	0.40
1:E:329:LEU:O	1:E:330:SER:C	2.63	0.40
1:E:367:GLU:C	1:E:369:GLY:N	2.78	0.40
1:G:237:THR:HG23	1:G:240:LYS:HE3	2.04	0.40
1:G:281:PHE:HB2	1:G:285:HIS:CE1	2.56	0.40
2:H:101:PHE:O	2:H:105:LEU:HG	2.21	0.40
1:I:246:ARG:HG2	1:I:272:GLU:HB2	2.04	0.40
1:I:420:GLU:HG3	1:I:423:GLY:HA2	2.02	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:376:THR:HG22	1:K:399:GLN:HB2	2.01	0.40
1:M:119:GLY:O	1:M:122:SER:OG	2.34	0.40
2:N:43:ASP:O	2:N:45:VAL:N	2.54	0.40
1:O:184:GLN:O	1:O:209:THR:OG1	2.39	0.40
2:P:8:SER:C	2:P:10:ASP:N	2.79	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:49:ASN:ND2	2:L:43:ASP:OD1[2_656]	1.93	0.27

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	323/336 (96%)	294 (91%)	21 (6%)	8 (2%)	4	16
1	C	323/336 (96%)	292 (90%)	24 (7%)	7 (2%)	5	19
1	E	323/336 (96%)	289 (90%)	27 (8%)	7 (2%)	5	19
1	G	323/336 (96%)	292 (90%)	24 (7%)	7 (2%)	5	19
1	I	323/336 (96%)	295 (91%)	21 (6%)	7 (2%)	5	19
1	K	323/336 (96%)	291 (90%)	24 (7%)	8 (2%)	4	16
1	M	323/336 (96%)	293 (91%)	25 (8%)	5 (2%)	8	28
1	O	323/336 (96%)	294 (91%)	21 (6%)	8 (2%)	4	16
2	B	133/149 (89%)	117 (88%)	12 (9%)	4 (3%)	3	12
2	D	133/149 (89%)	118 (89%)	10 (8%)	5 (4%)	2	9
2	F	133/149 (89%)	118 (89%)	11 (8%)	4 (3%)	3	12
2	H	133/149 (89%)	118 (89%)	12 (9%)	3 (2%)	5	18

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	J	133/149 (89%)	120 (90%)	8 (6%)	5 (4%)	2	9
2	L	133/149 (89%)	119 (90%)	9 (7%)	5 (4%)	2	9
2	N	133/149 (89%)	116 (87%)	13 (10%)	4 (3%)	3	12
2	P	133/149 (89%)	117 (88%)	11 (8%)	5 (4%)	2	9
All	All	3648/3880 (94%)	3283 (90%)	273 (8%)	92 (2%)	4	16

All (92) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	190	PHE
1	A	306	ARG
1	C	187	ALA
1	C	306	ARG
1	C	421	ILE
1	E	187	ALA
1	E	190	PHE
1	E	306	ARG
1	G	188	GLU
1	G	190	PHE
1	G	306	ARG
2	H	33	ASP
1	I	190	PHE
1	I	306	ARG
1	I	421	ILE
1	K	188	GLU
1	K	190	PHE
1	K	306	ARG
2	L	145	PHE
1	M	188	GLU
1	M	190	PHE
1	M	306	ARG
1	O	190	PHE
1	O	306	ARG
1	O	421	ILE
1	A	188	GLU
1	A	421	ILE
2	B	33	ASP
2	D	33	ASP
1	E	188	GLU
1	E	421	ILE
1	G	189	HIS

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Mol	Chain	Res	Type
1	G	421	ILE
1	I	188	GLU
1	K	187	ALA
1	K	421	ILE
2	L	33	ASP
1	M	421	ILE
2	N	33	ASP
2	N	159	TRP
1	O	188	GLU
2	P	33	ASP
2	P	140	ASN
1	A	186	LEU
1	C	186	LEU
1	C	188	GLU
1	C	418	ASN
1	E	186	LEU
2	F	33	ASP
2	F	140	ASN
2	H	140	ASN
2	J	33	ASP
1	O	189	HIS
2	B	140	ASN
2	D	43	ASP
2	D	113	LYS
2	D	140	ASN
1	G	187	ALA
2	H	43	ASP
1	I	189	HIS
1	I	333	VAL
1	K	189	HIS
2	L	43	ASP
2	L	140	ASN
1	M	187	ALA
2	N	43	ASP
1	O	279	PHE
2	P	43	ASP
2	P	159	TRP
1	A	418	ASN
2	B	43	ASP
2	B	159	TRP
2	D	137	LYS
2	F	43	ASP

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Mol	Chain	Res	Type
1	G	186	LEU
2	J	43	ASP
2	J	138	THR
2	J	140	ASN
1	K	418	ASN
2	L	138	THR
2	N	138	THR
1	O	195	VAL
2	F	138	THR
2	J	137	LYS
2	P	138	THR
1	C	195	VAL
1	E	195	VAL
1	I	195	VAL
1	A	195	VAL
1	A	333	VAL
1	O	408	ILE
1	K	195	VAL

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	299/311 (96%)	283 (95%)	16 (5%)	20	51
1	C	299/311 (96%)	283 (95%)	16 (5%)	20	51
1	E	299/311 (96%)	286 (96%)	13 (4%)	26	60
1	G	299/311 (96%)	284 (95%)	15 (5%)	22	54
1	I	299/311 (96%)	283 (95%)	16 (5%)	20	51
1	K	299/311 (96%)	285 (95%)	14 (5%)	23	57
1	M	299/311 (96%)	284 (95%)	15 (5%)	22	54
1	O	299/311 (96%)	284 (95%)	15 (5%)	22	54
2	B	124/134 (92%)	121 (98%)	3 (2%)	43	77
2	D	124/134 (92%)	120 (97%)	4 (3%)	34	70

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	F	124/134 (92%)	119 (96%)	5 (4%)	28	63
2	H	124/134 (92%)	120 (97%)	4 (3%)	34	70
2	J	124/134 (92%)	120 (97%)	4 (3%)	34	70
2	L	124/134 (92%)	121 (98%)	3 (2%)	43	77
2	N	124/134 (92%)	121 (98%)	3 (2%)	43	77
2	P	124/134 (92%)	119 (96%)	5 (4%)	28	63
All	All	3384/3560 (95%)	3233 (96%)	151 (4%)	24	58

All (151) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	114	ASP
1	A	115	GLU
1	A	128	GLU
1	A	157	LYS
1	A	192	PRO
1	A	207	VAL
1	A	262	THR
1	A	282	THR
1	A	307	LYS
1	A	314	LEU
1	A	317	LEU
1	A	342	GLU
1	A	389	GLN
1	A	399	GLN
1	A	418	ASN
1	A	420	GLU
2	B	55	LEU
2	B	94	LYS
2	B	116	LEU
1	C	114	ASP
1	C	115	GLU
1	C	128	GLU
1	C	157	LYS
1	C	191	SER
1	C	207	VAL
1	C	262	THR
1	C	274	ASN
1	C	307	LYS
1	C	314	LEU

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Mol	Chain	Res	Type
1	C	317	LEU
1	C	342	GLU
1	C	389	GLN
1	C	399	GLN
1	C	418	ASN
1	C	420	GLU
2	D	51	ASN
2	D	55	LEU
2	D	94	LYS
2	D	116	LEU
1	E	114	ASP
1	E	115	GLU
1	E	128	GLU
1	E	157	LYS
1	E	207	VAL
1	E	262	THR
1	E	307	LYS
1	E	314	LEU
1	E	317	LEU
1	E	342	GLU
1	E	399	GLN
1	E	418	ASN
1	E	420	GLU
2	F	51	ASN
2	F	55	LEU
2	F	94	LYS
2	F	116	LEU
2	F	118	VAL
1	G	114	ASP
1	G	115	GLU
1	G	128	GLU
1	G	157	LYS
1	G	207	VAL
1	G	262	THR
1	G	282	THR
1	G	307	LYS
1	G	314	LEU
1	G	317	LEU
1	G	342	GLU
1	G	389	GLN
1	G	399	GLN
1	G	418	ASN

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Mol	Chain	Res	Type
1	G	420	GLU
2	H	51	ASN
2	H	55	LEU
2	H	94	LYS
2	H	118	VAL
1	I	114	ASP
1	I	115	GLU
1	I	128	GLU
1	I	157	LYS
1	I	207	VAL
1	I	237	THR
1	I	262	THR
1	I	274	ASN
1	I	282	THR
1	I	307	LYS
1	I	314	LEU
1	I	317	LEU
1	I	342	GLU
1	I	399	GLN
1	I	418	ASN
1	I	420	GLU
2	J	55	LEU
2	J	94	LYS
2	J	116	LEU
2	J	118	VAL
1	K	114	ASP
1	K	115	GLU
1	K	128	GLU
1	K	157	LYS
1	K	207	VAL
1	K	262	THR
1	K	307	LYS
1	K	314	LEU
1	K	317	LEU
1	K	342	GLU
1	K	389	GLN
1	K	399	GLN
1	K	418	ASN
1	K	424	ILE
2	L	55	LEU
2	L	94	LYS
2	L	116	LEU

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Mol	Chain	Res	Type
1	M	114	ASP
1	M	115	GLU
1	M	128	GLU
1	M	157	LYS
1	M	207	VAL
1	M	262	THR
1	M	282	THR
1	M	307	LYS
1	M	314	LEU
1	M	317	LEU
1	M	342	GLU
1	M	389	GLN
1	M	399	GLN
1	M	418	ASN
1	M	420	GLU
2	N	55	LEU
2	N	94	LYS
2	N	116	LEU
1	O	114	ASP
1	O	115	GLU
1	O	128	GLU
1	O	157	LYS
1	O	207	VAL
1	O	262	THR
1	O	274	ASN
1	O	307	LYS
1	O	314	LEU
1	O	317	LEU
1	O	342	GLU
1	O	389	GLN
1	O	399	GLN
1	O	418	ASN
1	O	420	GLU
2	P	51	ASN
2	P	55	LEU
2	P	94	LYS
2	P	116	LEU
2	P	118	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (208) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	150	GLN
1	A	170	GLN
1	A	196	GLN
1	A	197	HIS
1	A	211	HIS
1	A	216	GLN
1	A	221	GLN
1	A	241	ASN
1	A	274	ASN
1	A	287	GLN
1	A	323	ASN
1	A	341	GLN
1	A	350	GLN
1	A	351	HIS
1	A	378	GLN
1	A	389	GLN
1	A	399	GLN
1	A	431	GLN
2	B	7	GLN
2	B	23	GLN
2	B	51	ASN
2	B	60	GLN
2	B	64	HIS
2	B	90	GLN
2	B	125	ASN
2	B	140	ASN
2	B	158	GLN
1	C	196	GLN
1	C	197	HIS
1	C	211	HIS
1	C	221	GLN
1	C	241	ASN
1	C	261	GLN
1	C	274	ASN
1	C	287	GLN
1	C	310	GLN
1	C	350	GLN
1	C	351	HIS
1	C	378	GLN
1	C	389	GLN
1	C	397	HIS
1	C	399	GLN
1	C	404	HIS

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Mol	Chain	Res	Type
1	C	431	GLN
2	D	7	GLN
2	D	23	GLN
2	D	51	ASN
2	D	64	HIS
2	D	65	HIS
2	D	90	GLN
2	D	125	ASN
2	D	158	GLN
1	E	170	GLN
1	E	196	GLN
1	E	197	HIS
1	E	211	HIS
1	E	216	GLN
1	E	221	GLN
1	E	236	ASN
1	E	241	ASN
1	E	261	GLN
1	E	274	ASN
1	E	287	GLN
1	E	350	GLN
1	E	351	HIS
1	E	389	GLN
1	E	399	GLN
1	E	404	HIS
1	E	431	GLN
2	F	7	GLN
2	F	23	GLN
2	F	64	HIS
2	F	65	HIS
2	F	90	GLN
2	F	125	ASN
2	F	143	ASN
2	F	158	GLN
1	G	150	GLN
1	G	170	GLN
1	G	196	GLN
1	G	197	HIS
1	G	211	HIS
1	G	216	GLN
1	G	221	GLN
1	G	236	ASN

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Mol	Chain	Res	Type
1	G	241	ASN
1	G	274	ASN
1	G	287	GLN
1	G	310	GLN
1	G	350	GLN
1	G	351	HIS
1	G	378	GLN
1	G	389	GLN
1	G	399	GLN
1	G	404	HIS
1	G	431	GLN
2	H	7	GLN
2	H	23	GLN
2	H	51	ASN
2	H	60	GLN
2	H	90	GLN
2	H	108	ASN
2	H	125	ASN
2	H	140	ASN
2	H	158	GLN
1	I	196	GLN
1	I	197	HIS
1	I	211	HIS
1	I	221	GLN
1	I	236	ASN
1	I	241	ASN
1	I	261	GLN
1	I	274	ASN
1	I	310	GLN
1	I	326	HIS
1	I	351	HIS
1	I	378	GLN
1	I	389	GLN
1	I	397	HIS
1	I	399	GLN
1	I	431	GLN
2	J	7	GLN
2	J	23	GLN
2	J	51	ASN
2	J	60	GLN
2	J	64	HIS
2	J	90	GLN

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Mol	Chain	Res	Type
2	J	108	ASN
2	J	125	ASN
2	J	140	ASN
2	J	158	GLN
1	K	170	GLN
1	K	197	HIS
1	K	211	HIS
1	K	216	GLN
1	K	221	GLN
1	K	241	ASN
1	K	261	GLN
1	K	274	ASN
1	K	287	GLN
1	K	310	GLN
1	K	326	HIS
1	K	350	GLN
1	K	351	HIS
1	K	378	GLN
1	K	389	GLN
1	K	397	HIS
1	K	399	GLN
1	K	404	HIS
1	K	431	GLN
2	L	7	GLN
2	L	23	GLN
2	L	51	ASN
2	L	64	HIS
2	L	90	GLN
2	L	108	ASN
2	L	125	ASN
2	L	143	ASN
2	L	158	GLN
1	M	170	GLN
1	M	196	GLN
1	M	197	HIS
1	M	211	HIS
1	M	216	GLN
1	M	221	GLN
1	M	236	ASN
1	M	241	ASN
1	M	261	GLN
1	M	274	ASN

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Mol	Chain	Res	Type
1	M	285	HIS
1	M	287	GLN
1	M	350	GLN
1	M	351	HIS
1	M	378	GLN
1	M	389	GLN
1	M	399	GLN
1	M	431	GLN
2	N	7	GLN
2	N	23	GLN
2	N	60	GLN
2	N	64	HIS
2	N	90	GLN
2	N	125	ASN
2	N	140	ASN
1	O	150	GLN
1	O	170	GLN
1	O	196	GLN
1	O	197	HIS
1	O	211	HIS
1	O	216	GLN
1	O	221	GLN
1	O	241	ASN
1	O	274	ASN
1	O	287	GLN
1	O	299	GLN
1	O	310	GLN
1	O	350	GLN
1	O	351	HIS
1	O	378	GLN
1	O	389	GLN
1	O	399	GLN
1	O	431	GLN
2	P	7	GLN
2	P	23	GLN
2	P	60	GLN
2	P	64	HIS
2	P	90	GLN
2	P	125	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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	G	2
1	I	1
1	C	1
1	O	1
1	K	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	G	186:LEU	C	187:ALA	N	1.70
1	I	186:LEU	C	187:ALA	N	1.63
1	C	189:HIS	C	190:PHE	N	1.19
1	O	189:HIS	C	190:PHE	N	1.17
1	K	189:HIS	C	190:PHE	N	1.13
1	G	189:HIS	C	190:PHE	N	1.12

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.