



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 01:50 AM UTC

PDB ID : 1XFU / pdb_00001xfu
Title : Crystal structure of anthrax edema factor (EF) truncation mutant, EF-delta 64 in complex with calmodulin
Authors : Shen, Y.; Zhukovskaya, N.L.; Guo, Q.; Florian, J.; Tang, W.J.
Deposited on : 2004-09-15
Resolution : 3.35 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

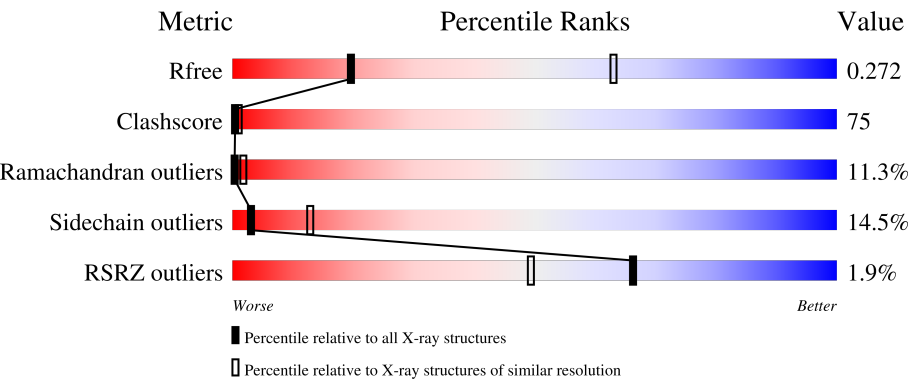
MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)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

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 3.35 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1099 (3.40-3.32)
Clashscore	190562	1116 (3.40-3.32)
Ramachandran outliers	187476	1101 (3.40-3.32)
Sidechain outliers	187428	1101 (3.40-3.32)
RSRZ outliers	180081	1099 (3.40-3.32)

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

Mol	Chain	Length	Quality of chain
1	A	747	3% 21% 54% 20%
1	B	747	2% 20% 55% 20%
1	C	747	2% 21% 55% 19%
1	D	747	2% 20% 54% 20%
1	E	747	2% 20% 55% 20%

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Mol	Chain	Length	Quality of chain				
1	F	747	2% 21% 55% 19% • •				
2	O	149	17% 59% 19% • •				
2	P	149	18% 60% 17% • •				
2	Q	149	% 17% 60% 18% • •				
2	R	149	17% 60% 18% • •				
2	S	149	18% 58% 18% • •				
2	T	149	% 18% 59% 17% • •				

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 42852 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 Calmodulin-sensitive adenylate cyclase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	735	Total	C	N	O	S	0	0	0
			5992	3828	995	1163	6			
1	B	735	Total	C	N	O	S	0	0	0
			5992	3828	995	1163	6			
1	C	735	Total	C	N	O	S	0	0	0
			5992	3828	995	1163	6			
1	D	735	Total	C	N	O	S	0	0	0
			5992	3828	995	1163	6			
1	E	735	Total	C	N	O	S	0	0	0
			5992	3828	995	1163	6			
1	F	735	Total	C	N	O	S	0	0	0
			5992	3828	995	1163	6			

There are 60 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	54	MET	-	initiating methionine	UNP P40136
A	55	HIS	-	expression tag	UNP P40136
A	56	HIS	-	expression tag	UNP P40136
A	57	HIS	-	expression tag	UNP P40136
A	58	HIS	-	expression tag	UNP P40136
A	59	HIS	-	expression tag	UNP P40136
A	60	HIS	-	expression tag	UNP P40136
A	61	ALA	-	cloning artifact	UNP P40136
A	62	ALA	-	cloning artifact	UNP P40136
A	63	ALA	-	cloning artifact	UNP P40136
B	54	MET	-	initiating methionine	UNP P40136
B	55	HIS	-	expression tag	UNP P40136
B	56	HIS	-	expression tag	UNP P40136
B	57	HIS	-	expression tag	UNP P40136
B	58	HIS	-	expression tag	UNP P40136
B	59	HIS	-	expression tag	UNP P40136
B	60	HIS	-	expression tag	UNP P40136

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Chain	Residue	Modelled	Actual	Comment	Reference
B	61	ALA	-	cloning artifact	UNP P40136
B	62	ALA	-	cloning artifact	UNP P40136
B	63	ALA	-	cloning artifact	UNP P40136
C	54	MET	-	initiating methionine	UNP P40136
C	55	HIS	-	expression tag	UNP P40136
C	56	HIS	-	expression tag	UNP P40136
C	57	HIS	-	expression tag	UNP P40136
C	58	HIS	-	expression tag	UNP P40136
C	59	HIS	-	expression tag	UNP P40136
C	60	HIS	-	expression tag	UNP P40136
C	61	ALA	-	cloning artifact	UNP P40136
C	62	ALA	-	cloning artifact	UNP P40136
C	63	ALA	-	cloning artifact	UNP P40136
D	54	MET	-	initiating methionine	UNP P40136
D	55	HIS	-	expression tag	UNP P40136
D	56	HIS	-	expression tag	UNP P40136
D	57	HIS	-	expression tag	UNP P40136
D	58	HIS	-	expression tag	UNP P40136
D	59	HIS	-	expression tag	UNP P40136
D	60	HIS	-	expression tag	UNP P40136
D	61	ALA	-	cloning artifact	UNP P40136
D	62	ALA	-	cloning artifact	UNP P40136
D	63	ALA	-	cloning artifact	UNP P40136
E	54	MET	-	initiating methionine	UNP P40136
E	55	HIS	-	expression tag	UNP P40136
E	56	HIS	-	expression tag	UNP P40136
E	57	HIS	-	expression tag	UNP P40136
E	58	HIS	-	expression tag	UNP P40136
E	59	HIS	-	expression tag	UNP P40136
E	60	HIS	-	expression tag	UNP P40136
E	61	ALA	-	cloning artifact	UNP P40136
E	62	ALA	-	cloning artifact	UNP P40136
E	63	ALA	-	cloning artifact	UNP P40136
F	54	MET	-	initiating methionine	UNP P40136
F	55	HIS	-	expression tag	UNP P40136
F	56	HIS	-	expression tag	UNP P40136
F	57	HIS	-	expression tag	UNP P40136
F	58	HIS	-	expression tag	UNP P40136
F	59	HIS	-	expression tag	UNP P40136
F	60	HIS	-	expression tag	UNP P40136
F	61	ALA	-	cloning artifact	UNP P40136
F	62	ALA	-	cloning artifact	UNP P40136

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Chain	Residue	Modelled	Actual	Comment	Reference
F	63	ALA	-	cloning artifact	UNP P40136

- Molecule 2 is a protein called Calmodulin 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	O	146	Total	C	N	O	S	0	0	0
			1146	702	186	249	9			
2	P	146	Total	C	N	O	S	0	0	0
			1146	702	186	249	9			
2	Q	146	Total	C	N	O	S	0	0	0
			1146	702	186	249	9			
2	R	146	Total	C	N	O	S	0	0	0
			1146	702	186	249	9			
2	S	146	Total	C	N	O	S	0	0	0
			1146	702	186	249	9			
2	T	146	Total	C	N	O	S	0	0	0
			1146	702	186	249	9			

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mg	0	0
			1	1		
3	B	1	Total	Mg	0	0
			1	1		
3	C	1	Total	Mg	0	0
			1	1		
3	D	1	Total	Mg	0	0
			1	1		
3	E	1	Total	Mg	0	0
			1	1		
3	F	1	Total	Mg	0	0
			1	1		

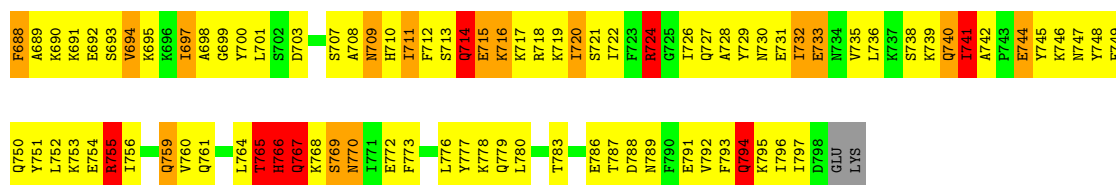
- Molecule 4 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	O	3	Total	Ca	0	0
			3	3		
4	P	3	Total	Ca	0	0
			3	3		

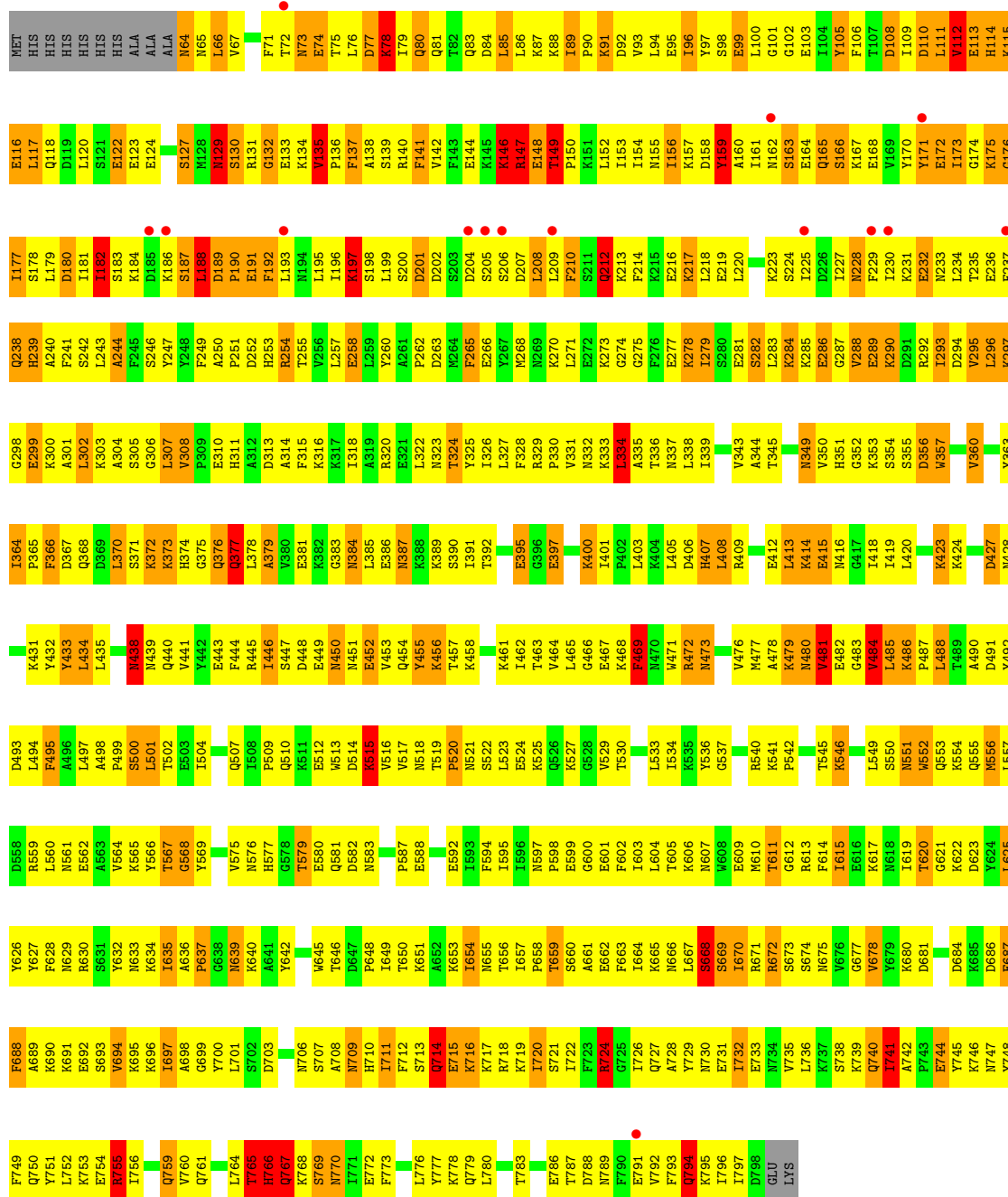
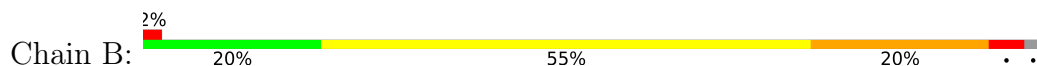
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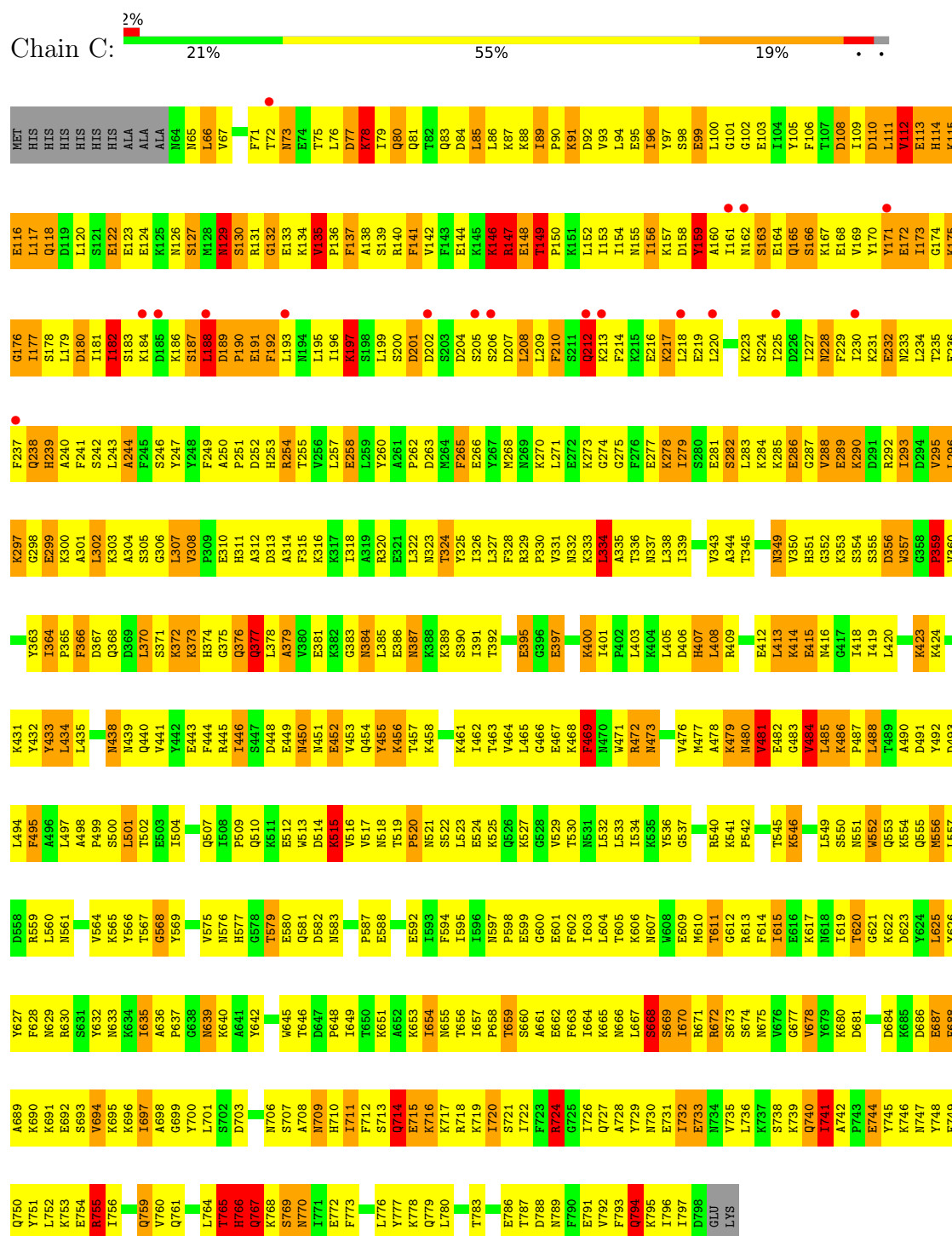
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	Q	3	Total 3	Ca 3	0	0
4	R	3	Total 3	Ca 3	0	0
4	S	3	Total 3	Ca 3	0	0
4	T	3	Total 3	Ca 3	0	0



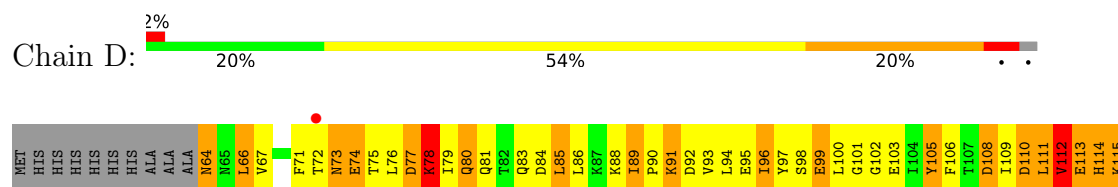
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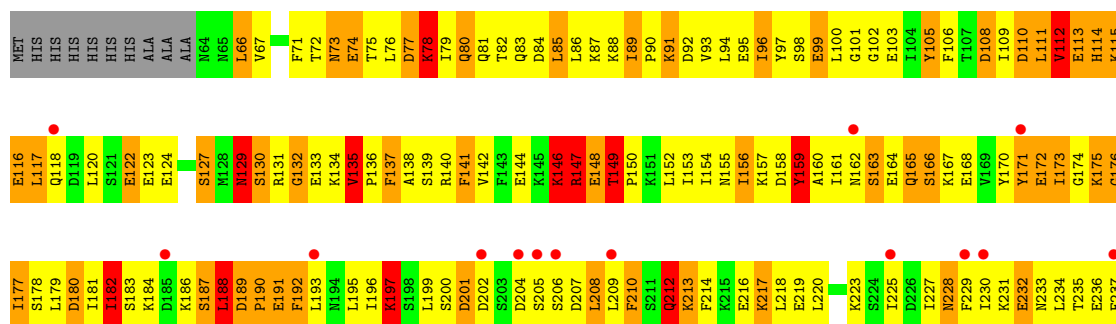
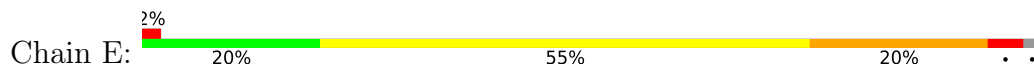
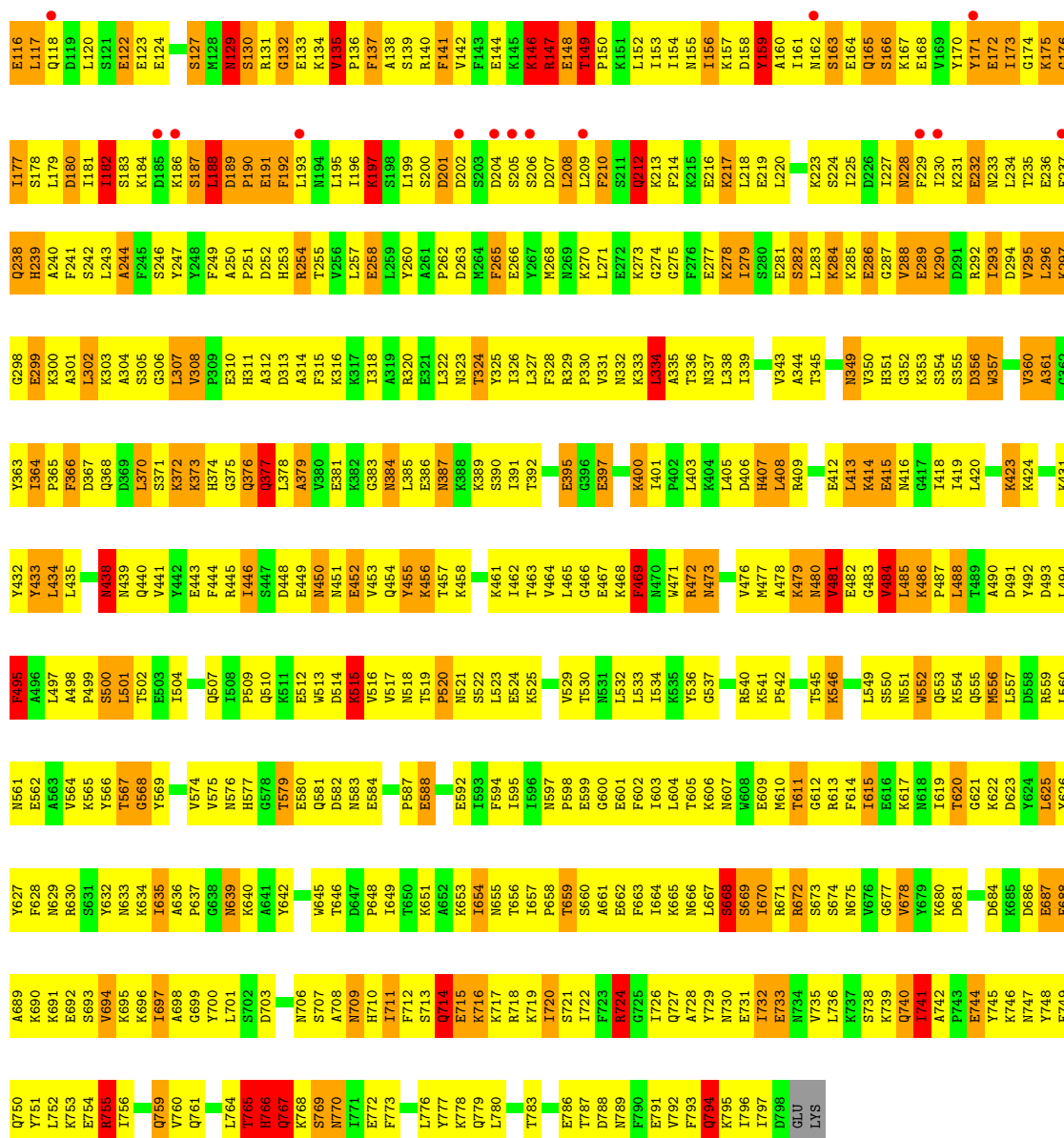


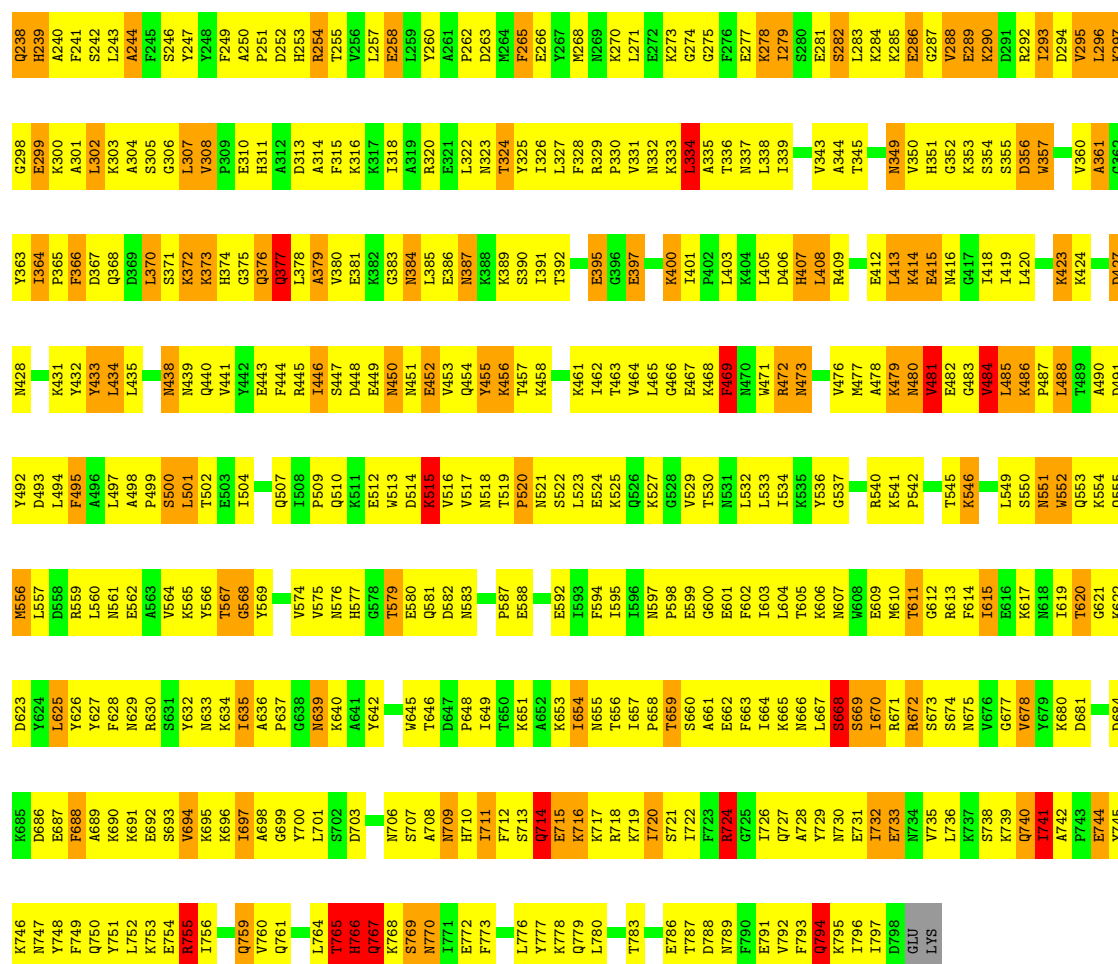
• Molecule 1: Calmodulin-sensitive adenylate cyclase



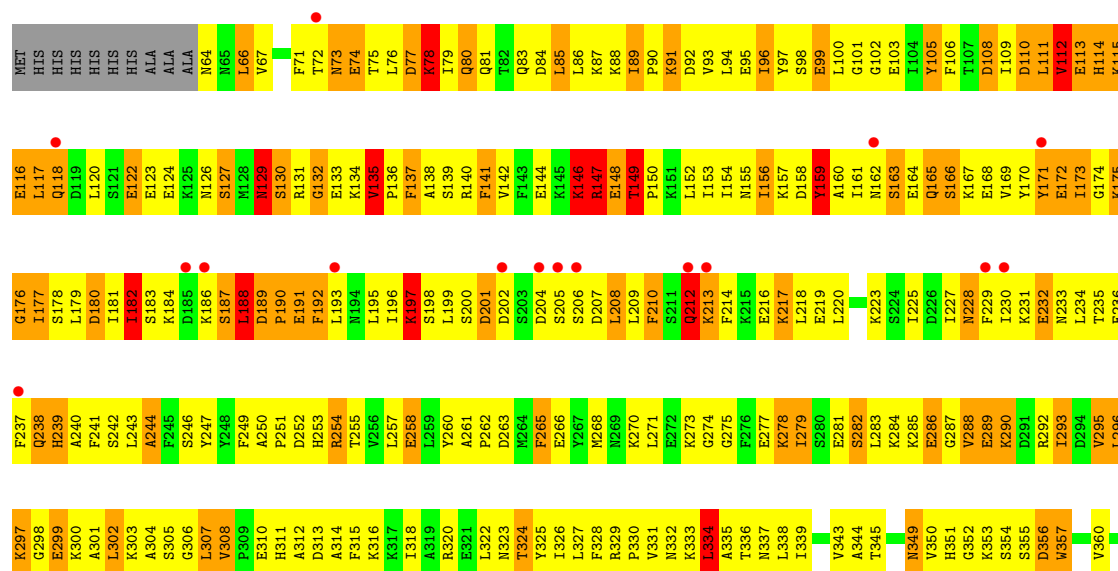
- Molecule 1: Calmodulin-sensitive adenylate cyclase

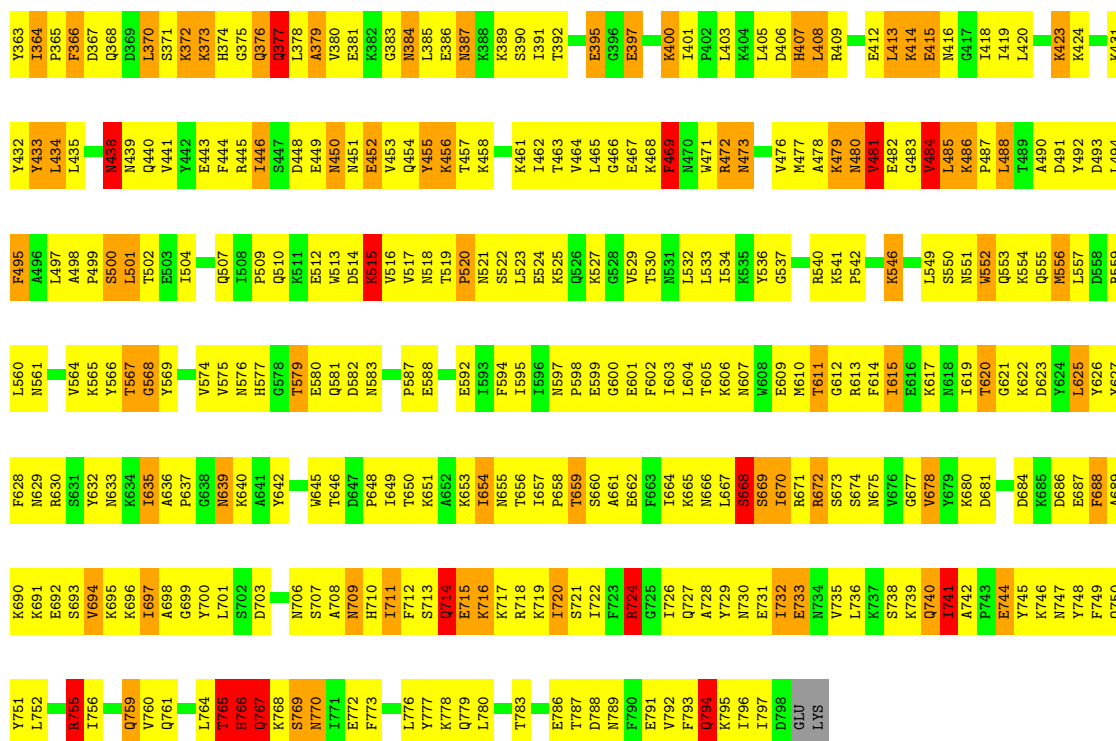




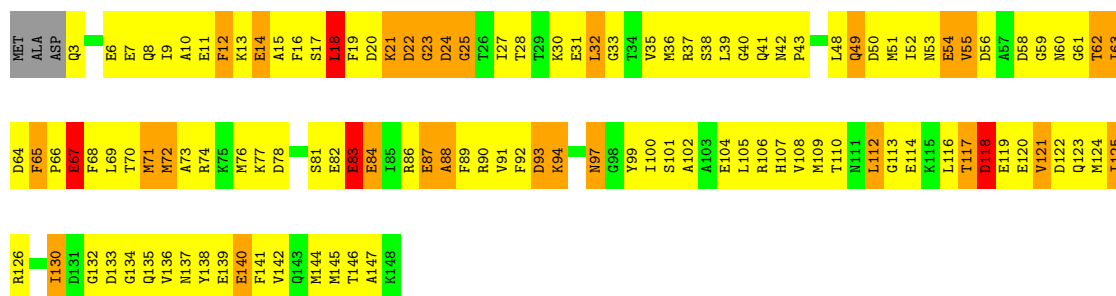


● Molecule 1: Calmodulin-sensitive adenylylate cyclase

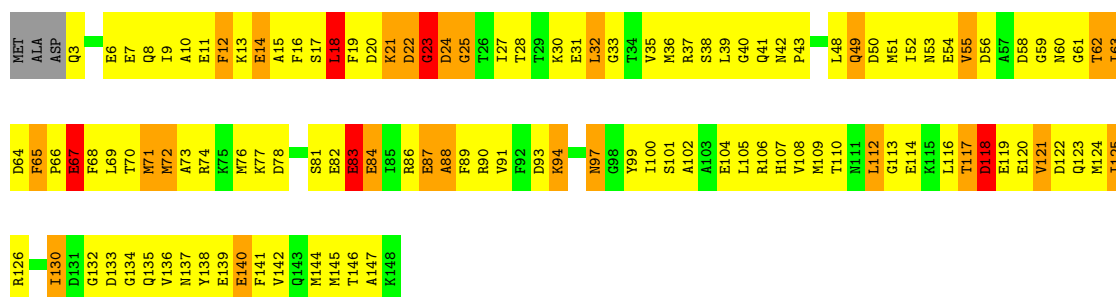
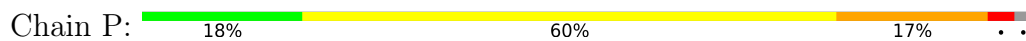




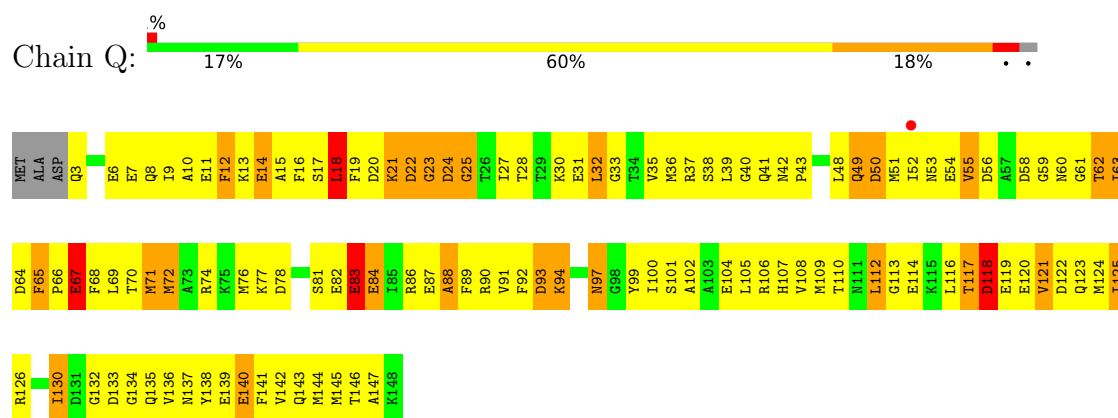
• Molecule 2: Calmodulin 2



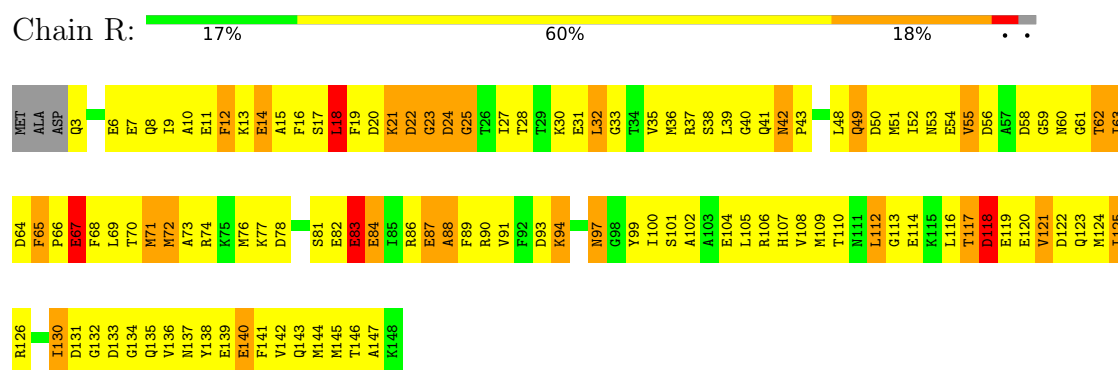
• Molecule 2: Calmodulin 2



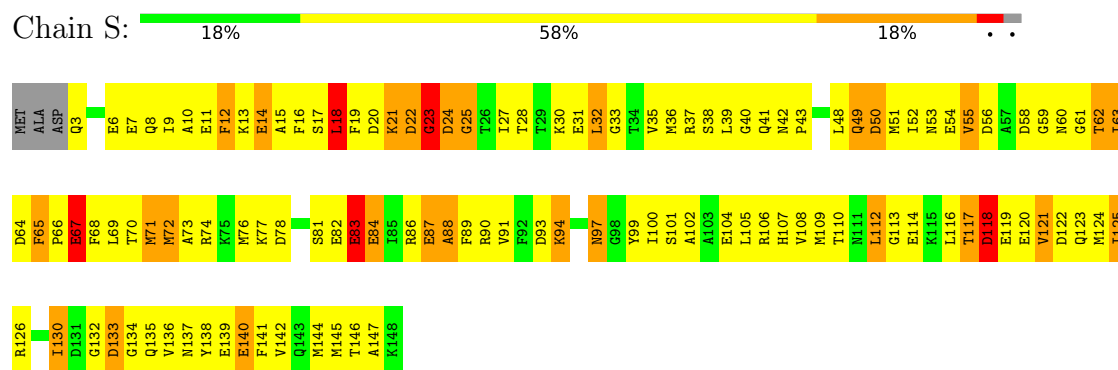
• Molecule 2: Calmodulin 2



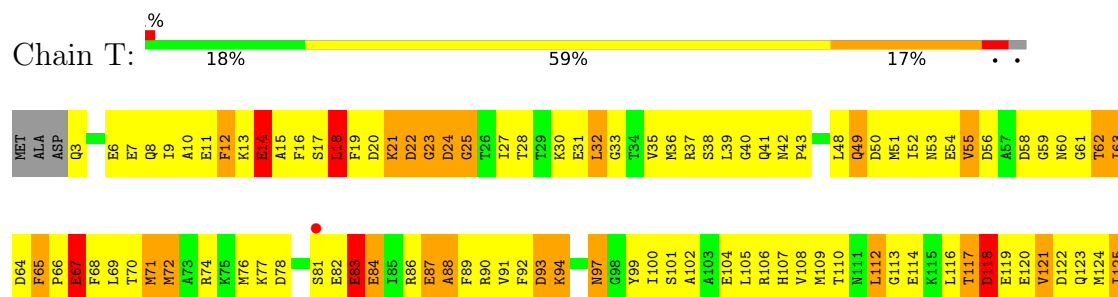
• Molecule 2: Calmodulin 2



• Molecule 2: Calmodulin 2



• Molecule 2: Calmodulin 2



R126	
I130	
D131	
G132	
D133	
G134	
Q135	
V136	
N137	
Y138	
E139	
E140	
F141	
V142	
Q143	
M144	
M145	
T146	
A147	
R148	

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	317.56Å 183.11Å 141.12Å 90.00° 90.05° 90.00°	Depositor
Resolution (Å)	29.99 – 3.35 29.99 – 3.35	Depositor EDS
% Data completeness (in resolution range)	95.8 (29.99-3.35) 95.7 (29.99-3.35)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	0.03	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.58 (at 3.32Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.283 , 0.300 0.255 , 0.272	Depositor DCC
R_{free} test set	6103 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	100.0	Xtriage
Anisotropy	0.184	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 67.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.468 for -1/2*h-3/2*k,-1/2*h+1/2*k,-l 0.468 for -1/2*h+3/2*k,1/2*h+1/2*k,-l 0.468 for 1/2*h-3/2*k,-1/2*h-1/2*k,-l 0.460 for 1/2*h+3/2*k,1/2*h-1/2*k,-l 0.468 for -h,-k,l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	42852	wwPDB-VP
Average B, all atoms (Å ²)	80.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.43% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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

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	0.65	1/6104 (0.0%)	1.22	79/8208 (1.0%)
1	B	0.66	2/6104 (0.0%)	1.22	80/8208 (1.0%)
1	C	0.68	4/6104 (0.1%)	1.24	83/8208 (1.0%)
1	D	0.67	2/6104 (0.0%)	1.22	81/8208 (1.0%)
1	E	0.67	2/6104 (0.0%)	1.22	80/8208 (1.0%)
1	F	0.68	4/6104 (0.1%)	1.22	79/8208 (1.0%)
2	O	0.68	0/1158	1.17	10/1553 (0.6%)
2	P	0.70	0/1158	1.18	10/1553 (0.6%)
2	Q	0.70	0/1158	1.16	9/1553 (0.6%)
2	R	0.71	0/1158	1.17	10/1553 (0.6%)
2	S	0.72	0/1158	1.17	10/1553 (0.6%)
2	T	0.70	0/1158	1.16	10/1553 (0.6%)
All	All	0.68	15/43572 (0.0%)	1.22	541/58566 (0.9%)

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	132	GLY	C-N	10.92	1.49	1.33
1	C	132	GLY	C-N	9.96	1.47	1.33
1	D	132	GLY	C-N	9.79	1.47	1.33
1	E	132	GLY	C-N	7.90	1.44	1.33
1	B	132	GLY	C-N	7.42	1.44	1.33
1	F	126	ASN	C-N	6.22	1.42	1.33
1	C	126	ASN	C-N	5.88	1.42	1.33
1	C	118	GLN	C-N	-5.87	1.25	1.33
1	F	118	GLN	C-N	-5.39	1.25	1.33
1	C	188	LEU	C-N	-5.16	1.25	1.33
1	B	188	LEU	C-N	-5.14	1.25	1.33
1	D	188	LEU	C-N	-5.14	1.25	1.33
1	E	188	LEU	C-N	-5.13	1.25	1.33
1	A	188	LEU	C-N	-5.13	1.25	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	188	LEU	C-N	-5.11	1.25	1.33

All (541) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	159	TYR	N-CA-C	14.45	126.72	110.97
1	C	159	TYR	N-CA-C	14.42	126.69	110.97
1	A	159	TYR	N-CA-C	14.40	126.67	110.97
1	F	159	TYR	N-CA-C	14.40	126.67	110.97
1	D	159	TYR	N-CA-C	14.36	126.62	110.97
1	B	159	TYR	N-CA-C	14.36	126.62	110.97
1	D	147	ARG	N-CA-C	12.36	129.09	112.90
1	F	147	ARG	N-CA-C	12.35	129.08	112.90
1	A	147	ARG	N-CA-C	12.33	129.05	112.90
1	C	147	ARG	N-CA-C	12.33	129.05	112.90
1	E	147	ARG	N-CA-C	12.32	129.04	112.90
1	B	147	ARG	N-CA-C	12.32	129.04	112.90
1	D	219	GLU	N-CA-C	-11.74	98.48	112.92
1	A	219	GLU	N-CA-C	-11.74	98.48	112.92
1	F	219	GLU	N-CA-C	-11.73	98.50	112.92
1	E	219	GLU	N-CA-C	-11.69	98.54	112.92
1	C	219	GLU	N-CA-C	-11.65	98.60	112.92
1	B	219	GLU	N-CA-C	-11.56	98.70	112.92
1	B	212	GLN	OE1-CD-NE2	-10.07	112.53	122.60
1	C	212	GLN	OE1-CD-NE2	-10.06	112.53	122.60
1	D	212	GLN	OE1-CD-NE2	-10.06	112.53	122.60
1	E	212	GLN	OE1-CD-NE2	-10.04	112.56	122.60
1	A	212	GLN	OE1-CD-NE2	-10.03	112.57	122.60
1	F	212	GLN	OE1-CD-NE2	-10.02	112.58	122.60
1	F	740	GLN	OE1-CD-NE2	-9.79	112.81	122.60
1	E	767	GLN	OE1-CD-NE2	-9.75	112.85	122.60
1	D	740	GLN	OE1-CD-NE2	-9.74	112.86	122.60
1	A	740	GLN	OE1-CD-NE2	-9.73	112.87	122.60
1	D	767	GLN	OE1-CD-NE2	-9.73	112.87	122.60
1	C	740	GLN	OE1-CD-NE2	-9.72	112.88	122.60
1	B	740	GLN	OE1-CD-NE2	-9.72	112.88	122.60
1	C	767	GLN	OE1-CD-NE2	-9.72	112.88	122.60
1	A	767	GLN	OE1-CD-NE2	-9.71	112.89	122.60
1	F	767	GLN	OE1-CD-NE2	-9.69	112.91	122.60
1	B	767	GLN	OE1-CD-NE2	-9.69	112.91	122.60
1	E	740	GLN	OE1-CD-NE2	-9.69	112.91	122.60
1	D	674	SER	N-CA-C	-9.62	95.44	109.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	674	SER	N-CA-C	-9.62	95.46	109.59
1	F	674	SER	N-CA-C	-9.62	95.45	109.59
1	A	674	SER	N-CA-C	-9.61	95.46	109.59
1	B	674	SER	N-CA-C	-9.61	95.47	109.59
1	E	674	SER	N-CA-C	-9.61	95.47	109.59
1	C	770	ASN	N-CA-C	-8.97	104.39	114.62
1	E	770	ASN	N-CA-C	-8.94	104.42	114.62
1	A	770	ASN	N-CA-C	-8.94	104.43	114.62
1	F	770	ASN	N-CA-C	-8.94	104.43	114.62
1	B	770	ASN	N-CA-C	-8.93	104.44	114.62
1	D	770	ASN	N-CA-C	-8.92	104.46	114.62
1	C	149	THR	N-CA-C	-8.77	98.18	110.29
1	D	149	THR	N-CA-C	-8.77	98.19	110.29
1	F	149	THR	N-CA-C	-8.76	98.21	110.29
1	A	149	THR	N-CA-C	-8.75	98.21	110.29
1	E	149	THR	N-CA-C	-8.75	98.21	110.29
1	B	149	THR	N-CA-C	-8.74	98.23	110.29
1	B	160	ALA	N-CA-C	8.66	122.51	109.25
1	E	160	ALA	N-CA-C	8.64	122.48	109.25
1	A	160	ALA	N-CA-C	8.64	122.47	109.25
1	D	160	ALA	N-CA-C	8.64	122.47	109.25
1	D	127	SER	N-CA-C	8.63	124.13	108.17
1	C	160	ALA	N-CA-C	8.62	122.45	109.25
1	A	127	SER	N-CA-C	8.61	124.10	108.17
1	F	160	ALA	N-CA-C	8.61	122.42	109.25
1	F	127	SER	N-CA-C	8.60	124.14	107.98
1	E	127	SER	N-CA-C	8.59	124.14	107.98
1	C	127	SER	N-CA-C	8.58	124.12	107.98
1	B	127	SER	N-CA-C	8.57	124.09	107.98
2	S	88	ALA	N-CA-C	-8.38	102.22	111.36
2	S	121	VAL	N-CA-C	-8.27	103.81	111.67
1	B	433	TYR	N-CA-C	-8.27	96.61	109.76
1	A	433	TYR	N-CA-C	-8.25	96.64	109.76
1	C	433	TYR	N-CA-C	-8.25	96.64	109.76
2	P	88	ALA	N-CA-C	-8.25	102.36	111.36
1	E	433	TYR	N-CA-C	-8.25	96.64	109.76
1	D	433	TYR	N-CA-C	-8.23	96.67	109.76
1	F	433	TYR	N-CA-C	-8.23	96.67	109.76
2	O	88	ALA	N-CA-C	-8.17	102.46	111.36
2	T	121	VAL	N-CA-C	-8.15	103.93	111.67
2	Q	121	VAL	N-CA-C	-8.10	103.97	111.67
2	R	88	ALA	N-CA-C	-8.09	102.54	111.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	Q	88	ALA	N-CA-C	-8.07	102.56	111.36
1	D	188	LEU	CA-C-N	-8.06	110.37	122.29
1	D	188	LEU	C-N-CA	-8.06	110.37	122.29
1	E	188	LEU	CA-C-N	-8.05	110.38	122.29
1	E	188	LEU	C-N-CA	-8.05	110.38	122.29
2	O	121	VAL	N-CA-C	-8.04	104.03	111.67
1	A	188	LEU	CA-C-N	-8.03	110.41	122.29
1	A	188	LEU	C-N-CA	-8.03	110.41	122.29
1	F	188	LEU	CA-C-N	-8.01	110.43	122.29
1	F	188	LEU	C-N-CA	-8.01	110.43	122.29
1	B	188	LEU	CA-C-N	-8.00	110.45	122.29
1	B	188	LEU	C-N-CA	-8.00	110.45	122.29
1	C	188	LEU	CA-C-N	-8.00	110.45	122.29
1	C	188	LEU	C-N-CA	-8.00	110.45	122.29
1	A	188	LEU	N-CA-C	-8.00	93.77	110.80
1	D	188	LEU	N-CA-C	-8.00	93.77	110.80
1	F	188	LEU	N-CA-C	-8.00	93.77	110.80
1	B	188	LEU	N-CA-C	-7.99	93.78	110.80
1	E	188	LEU	N-CA-C	-7.99	93.78	110.80
1	C	188	LEU	N-CA-C	-7.98	93.80	110.80
2	R	121	VAL	N-CA-C	-7.97	104.10	111.67
1	E	182	ILE	N-CA-C	-7.94	105.13	112.43
2	P	121	VAL	N-CA-C	-7.91	104.16	111.67
1	A	182	ILE	N-CA-C	-7.88	105.18	112.43
1	E	238	GLN	N-CA-C	-7.88	102.20	112.68
1	D	182	ILE	N-CA-C	-7.82	105.24	112.43
2	T	88	ALA	N-CA-C	-7.79	102.78	111.28
1	F	182	ILE	N-CA-C	-7.79	105.26	112.43
1	B	238	GLN	N-CA-C	-7.78	102.34	112.68
1	F	238	GLN	N-CA-C	-7.77	102.35	112.68
1	C	182	ILE	N-CA-C	-7.73	105.32	112.43
1	A	238	GLN	N-CA-C	-7.70	102.43	112.68
1	C	238	GLN	N-CA-C	-7.70	102.44	112.68
1	D	238	GLN	N-CA-C	-7.68	102.46	112.68
1	B	182	ILE	N-CA-C	-7.65	105.40	112.43
2	P	122	ASP	N-CA-C	-7.64	103.70	112.87
2	Q	122	ASP	N-CA-C	-7.47	103.90	112.87
2	S	122	ASP	N-CA-C	-7.47	103.91	112.87
2	O	122	ASP	N-CA-C	-7.46	103.92	112.87
2	R	122	ASP	N-CA-C	-7.43	103.95	112.87
2	T	122	ASP	N-CA-C	-7.40	103.99	112.87
1	A	173	ILE	N-CA-C	-7.32	103.65	110.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	173	ILE	N-CA-C	-7.29	103.67	110.53
1	C	173	ILE	N-CA-C	-7.29	103.68	110.53
1	D	765	THR	CA-C-N	-7.26	109.69	122.26
1	D	765	THR	C-N-CA	-7.26	109.69	122.26
1	B	765	THR	CA-C-N	-7.25	109.71	122.26
1	B	765	THR	C-N-CA	-7.25	109.71	122.26
1	E	765	THR	CA-C-N	-7.25	109.71	122.26
1	E	765	THR	C-N-CA	-7.25	109.71	122.26
1	F	765	THR	CA-C-N	-7.25	109.72	122.26
1	F	765	THR	C-N-CA	-7.25	109.72	122.26
1	A	765	THR	CA-C-N	-7.24	109.73	122.26
1	A	765	THR	C-N-CA	-7.24	109.73	122.26
1	B	173	ILE	N-CA-C	-7.24	103.73	110.53
1	C	765	THR	CA-C-N	-7.20	109.80	122.26
1	C	765	THR	C-N-CA	-7.20	109.80	122.26
1	E	173	ILE	N-CA-C	-7.20	103.77	110.53
1	A	566	TYR	N-CA-C	-7.16	104.91	112.93
1	F	173	ILE	N-CA-C	-7.12	103.84	110.53
1	D	566	TYR	N-CA-C	-7.11	104.96	112.93
1	C	566	TYR	N-CA-C	-7.10	104.97	112.93
1	E	566	TYR	N-CA-C	-6.97	105.12	112.93
1	E	212	GLN	CG-CD-NE2	6.90	126.75	116.40
1	C	212	GLN	CG-CD-NE2	6.90	126.74	116.40
1	D	212	GLN	CG-CD-NE2	6.89	126.74	116.40
1	F	212	GLN	CG-CD-NE2	6.88	126.73	116.40
1	A	99	GLU	N-CA-C	-6.88	104.76	113.43
1	B	212	GLN	CG-CD-NE2	6.88	126.72	116.40
1	A	212	GLN	CG-CD-NE2	6.88	126.72	116.40
1	F	566	TYR	N-CA-C	-6.87	105.24	112.93
1	B	99	GLU	N-CA-C	-6.86	104.78	113.43
1	B	566	TYR	N-CA-C	-6.85	105.25	112.93
1	D	693	SER	N-CA-C	6.83	119.57	111.71
1	C	189	ASP	CA-CB-CG	-6.82	105.78	112.60
1	F	189	ASP	CA-CB-CG	-6.82	105.78	112.60
1	A	189	ASP	CA-CB-CG	-6.81	105.79	112.60
1	D	277	GLU	N-CA-C	-6.81	104.44	112.89
1	B	189	ASP	CA-CB-CG	-6.81	105.79	112.60
1	F	99	GLU	N-CA-C	-6.80	104.87	113.43
1	F	769	SER	N-CA-C	6.79	121.34	107.69
1	C	769	SER	N-CA-C	6.79	121.34	107.69
1	E	189	ASP	CA-CB-CG	-6.79	105.81	112.60
1	D	769	SER	N-CA-C	6.78	121.33	107.69

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	99	GLU	N-CA-C	-6.78	104.88	113.43
1	E	769	SER	N-CA-C	6.78	121.32	107.69
1	A	769	SER	N-CA-C	6.78	121.31	107.69
1	C	96	ILE	N-CA-C	-6.78	103.88	110.72
1	D	189	ASP	CA-CB-CG	-6.78	105.82	112.60
1	D	99	GLU	N-CA-C	-6.77	104.90	113.43
1	B	769	SER	N-CA-C	6.77	121.29	107.69
1	D	469	PHE	N-CA-C	6.75	119.69	109.23
1	E	469	PHE	N-CA-C	6.74	119.68	109.23
1	C	469	PHE	N-CA-C	6.74	119.67	109.23
1	F	469	PHE	N-CA-C	6.74	119.67	109.23
1	A	579	THR	N-CA-C	6.73	120.46	109.76
1	B	277	GLU	N-CA-C	-6.70	104.58	112.89
1	A	277	GLU	N-CA-C	-6.70	104.58	112.89
1	A	469	PHE	N-CA-C	6.70	119.61	109.23
1	F	579	THR	N-CA-C	6.67	120.37	109.76
1	C	99	GLU	N-CA-C	-6.67	105.03	113.43
1	B	469	PHE	N-CA-C	6.66	119.56	109.23
1	A	133	GLU	N-CA-C	6.66	118.64	108.52
1	A	693	SER	N-CA-C	6.66	119.37	111.71
1	E	129	ASN	N-CA-C	6.65	124.96	110.80
1	C	129	ASN	N-CA-C	6.64	124.95	110.80
1	F	129	ASN	N-CA-C	6.64	124.94	110.80
1	C	277	GLU	N-CA-C	-6.64	104.66	112.89
1	A	129	ASN	N-CA-C	6.63	124.93	110.80
1	D	129	ASN	N-CA-C	6.63	124.92	110.80
1	B	129	ASN	N-CA-C	6.62	124.90	110.80
1	E	277	GLU	N-CA-C	-6.61	104.69	112.89
1	C	693	SER	N-CA-C	6.61	119.31	111.71
1	C	579	THR	N-CA-C	6.60	120.25	109.76
1	D	579	THR	N-CA-C	6.60	120.25	109.76
1	F	767	GLN	CG-CD-NE2	6.58	126.27	116.40
1	A	767	GLN	CG-CD-NE2	6.58	126.26	116.40
1	E	767	GLN	CG-CD-NE2	6.57	126.26	116.40
1	F	732	ILE	N-CA-C	-6.57	104.45	110.82
1	F	277	GLU	N-CA-C	-6.57	104.75	112.89
1	C	767	GLN	CG-CD-NE2	6.56	126.25	116.40
1	E	693	SER	N-CA-C	6.56	119.26	111.71
1	B	767	GLN	CG-CD-NE2	6.56	126.25	116.40
1	D	767	GLN	CG-CD-NE2	6.56	126.24	116.40
1	B	579	THR	N-CA-C	6.54	120.15	109.76
1	E	579	THR	N-CA-C	6.54	120.16	109.76

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	693	SER	N-CA-C	6.53	119.21	111.71
1	B	693	SER	N-CA-C	6.52	119.21	111.71
1	F	740	GLN	CG-CD-NE2	6.50	126.14	116.40
1	F	158	ASP	CA-C-N	6.48	129.21	120.65
1	F	158	ASP	C-N-CA	6.48	129.21	120.65
1	C	740	GLN	CG-CD-NE2	6.48	126.12	116.40
1	C	158	ASP	CA-C-N	6.48	129.20	120.65
1	C	158	ASP	C-N-CA	6.48	129.20	120.65
1	A	740	GLN	CG-CD-NE2	6.47	126.11	116.40
1	B	158	ASP	CA-C-N	6.47	129.19	120.65
1	B	158	ASP	C-N-CA	6.47	129.19	120.65
1	D	740	GLN	CG-CD-NE2	6.47	126.10	116.40
1	E	740	GLN	CG-CD-NE2	6.46	126.09	116.40
1	B	740	GLN	CG-CD-NE2	6.46	126.09	116.40
1	A	158	ASP	CA-C-N	6.45	129.16	120.65
1	A	158	ASP	C-N-CA	6.45	129.16	120.65
1	D	158	ASP	CA-C-N	6.45	129.16	120.65
1	D	158	ASP	C-N-CA	6.45	129.16	120.65
1	E	158	ASP	CA-C-N	6.45	129.16	120.65
1	E	158	ASP	C-N-CA	6.45	129.16	120.65
1	D	96	ILE	N-CA-C	-6.44	104.21	110.72
1	A	96	ILE	N-CA-C	-6.43	104.23	110.72
1	F	96	ILE	N-CA-C	-6.40	104.25	110.72
1	E	96	ILE	N-CA-C	-6.32	104.33	110.72
1	C	73	ASN	N-CA-C	6.32	118.73	111.02
1	A	732	ILE	N-CA-C	-6.29	104.72	110.82
1	B	96	ILE	N-CA-C	-6.28	104.38	110.72
1	E	427	ASP	N-CA-C	6.28	118.26	107.28
1	E	732	ILE	N-CA-C	-6.27	104.74	110.82
1	D	732	ILE	N-CA-C	-6.26	104.74	110.82
1	B	732	ILE	N-CA-C	-6.26	104.75	110.82
2	P	87	GLU	N-CA-C	-6.23	105.81	113.41
1	B	611	THR	N-CA-C	-6.20	104.43	111.07
2	S	87	GLU	N-CA-C	-6.20	105.84	113.41
1	C	552	TRP	N-CA-C	-6.18	104.54	111.28
1	C	732	ILE	N-CA-C	-6.18	104.82	110.82
1	E	611	THR	N-CA-C	-6.16	104.48	111.07
1	B	282	SER	N-CA-C	-6.14	105.77	113.20
1	F	552	TRP	N-CA-C	-6.12	104.61	111.28
1	A	552	TRP	N-CA-C	-6.12	104.61	111.28
1	E	304	ALA	N-CA-C	-6.12	105.85	113.38
1	D	552	TRP	N-CA-C	-6.11	104.62	111.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	O	83	GLU	N-CA-C	6.11	117.74	111.14
1	D	440	GLN	N-CA-C	6.11	119.68	112.72
1	C	488	LEU	N-CA-C	6.10	119.41	109.59
1	F	282	SER	N-CA-C	-6.10	105.82	113.20
1	F	440	GLN	N-CA-C	6.08	119.65	112.72
1	A	611	THR	N-CA-C	-6.06	104.59	111.07
1	D	282	SER	N-CA-C	-6.03	105.91	113.20
2	T	83	GLU	N-CA-C	6.02	117.65	111.14
2	R	83	GLU	N-CA-C	6.02	117.64	111.14
1	D	304	ALA	N-CA-C	-6.01	105.99	113.38
1	A	488	LEU	N-CA-C	6.01	119.26	109.59
1	A	440	GLN	N-CA-C	6.01	119.57	112.72
1	F	304	ALA	N-CA-C	-6.01	105.99	113.38
2	Q	83	GLU	N-CA-C	6.00	117.62	111.14
1	F	791	GLU	N-CA-C	-6.00	104.74	111.28
1	C	440	GLN	N-CA-C	6.00	119.56	112.72
1	A	282	SER	N-CA-C	-5.99	105.95	113.20
1	E	282	SER	N-CA-C	-5.99	105.95	113.20
1	B	488	LEU	N-CA-C	5.99	119.23	109.59
1	C	282	SER	N-CA-C	-5.98	105.97	113.20
1	B	304	ALA	N-CA-C	-5.97	106.04	113.38
1	C	304	ALA	N-CA-C	-5.97	106.04	113.38
1	F	611	THR	N-CA-C	-5.97	104.69	111.07
1	E	440	GLN	N-CA-C	5.96	119.52	112.72
1	A	258	GLU	N-CA-C	-5.96	105.84	114.12
1	F	697	ILE	N-CA-C	-5.95	104.52	111.00
1	E	724	ARG	N-CA-C	-5.94	104.14	111.33
1	A	208	LEU	N-CA-C	-5.94	104.64	112.72
1	B	791	GLU	N-CA-C	-5.94	104.81	111.28
1	B	440	GLN	N-CA-C	5.93	119.48	112.72
2	P	83	GLU	N-CA-C	5.93	117.54	111.14
1	C	791	GLU	N-CA-C	-5.92	104.83	111.28
1	B	258	GLU	N-CA-C	-5.91	105.90	114.12
1	B	366	PHE	N-CA-C	-5.91	104.18	111.33
1	D	488	LEU	N-CA-C	5.91	119.11	109.59
1	E	488	LEU	N-CA-C	5.91	119.11	109.59
1	E	258	GLU	N-CA-C	-5.91	105.91	114.12
1	F	208	LEU	N-CA-C	-5.91	104.69	112.72
1	D	611	THR	N-CA-C	-5.91	104.75	111.07
2	S	83	GLU	N-CA-C	5.91	117.52	111.14
1	B	552	TRP	N-CA-C	-5.90	104.85	111.28
1	B	265	PHE	N-CA-C	-5.90	104.76	111.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	208	LEU	N-CA-C	-5.90	104.70	112.72
1	E	552	TRP	N-CA-C	-5.89	104.86	111.28
1	F	286	GLU	N-CA-C	-5.89	106.14	113.15
1	A	791	GLU	N-CA-C	-5.89	104.86	111.28
1	C	611	THR	N-CA-C	-5.88	104.78	111.07
1	C	66	LEU	N-CA-C	5.88	117.77	111.36
1	E	208	LEU	N-CA-C	-5.88	104.72	112.72
1	E	791	GLU	N-CA-C	-5.87	104.88	111.28
1	F	366	PHE	N-CA-C	-5.87	104.23	111.33
1	E	286	GLU	N-CA-C	-5.87	106.17	113.15
1	A	265	PHE	N-CA-C	-5.86	104.80	111.07
1	E	366	PHE	N-CA-C	-5.86	104.23	111.33
1	A	304	ALA	N-CA-C	-5.86	106.17	113.38
1	C	366	PHE	N-CA-C	-5.85	104.25	111.33
1	D	791	GLU	N-CA-C	-5.84	104.91	111.28
1	C	208	LEU	N-CA-C	-5.84	104.78	112.72
1	F	488	LEU	N-CA-C	5.83	118.98	109.59
1	E	766	HIS	N-CA-C	-5.83	106.21	113.38
1	C	286	GLU	N-CA-C	-5.83	106.21	113.15
1	D	171	TYR	N-CA-C	-5.83	106.02	113.01
2	O	32	LEU	N-CA-C	-5.83	104.83	111.07
1	C	265	PHE	N-CA-C	-5.83	104.84	111.07
1	E	265	PHE	N-CA-C	-5.82	104.84	111.07
1	A	286	GLU	N-CA-C	-5.82	106.23	113.15
1	C	258	GLU	N-CA-C	-5.82	106.03	114.12
1	F	66	LEU	N-CA-C	5.82	117.62	111.28
1	F	766	HIS	N-CA-C	-5.82	106.23	113.38
1	B	171	TYR	N-CA-C	-5.81	106.03	113.01
1	C	724	ARG	N-CA-C	-5.81	104.30	111.33
1	A	366	PHE	N-CA-C	-5.81	104.30	111.33
1	B	208	LEU	N-CA-C	-5.81	104.82	112.72
1	D	694	VAL	N-CA-C	-5.81	104.67	111.00
1	C	694	VAL	N-CA-C	-5.81	104.67	111.00
1	A	766	HIS	N-CA-C	-5.80	106.24	113.38
1	C	766	HIS	N-CA-C	-5.80	106.24	113.38
2	T	32	LEU	N-CA-C	-5.80	104.86	111.07
2	S	32	LEU	N-CA-C	-5.80	104.86	111.07
1	B	766	HIS	N-CA-C	-5.80	106.25	113.38
1	F	265	PHE	N-CA-C	-5.79	104.87	111.07
1	D	258	GLU	N-CA-C	-5.79	106.08	114.12
1	D	766	HIS	N-CA-C	-5.78	106.27	113.38
1	C	359	PRO	CA-N-CD	-5.78	103.91	112.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	286	GLU	N-CA-C	-5.77	106.28	113.15
1	B	724	ARG	N-CA-C	-5.77	104.35	111.33
1	F	201	ASP	N-CA-C	-5.77	103.80	111.24
1	D	265	PHE	N-CA-C	-5.77	104.90	111.07
1	B	360	VAL	CB-CA-C	-5.77	101.83	111.29
2	P	32	LEU	N-CA-C	-5.77	104.90	111.07
1	D	201	ASP	N-CA-C	-5.76	103.80	111.24
1	D	724	ARG	N-CA-C	-5.76	104.36	111.33
1	A	171	TYR	N-CA-C	-5.75	106.11	113.01
1	C	171	TYR	N-CA-C	-5.75	106.11	113.01
1	F	546	LYS	N-CA-C	5.75	120.53	112.72
1	C	201	ASP	N-CA-C	-5.74	103.83	111.24
1	C	546	LYS	N-CA-C	5.74	120.53	112.72
1	D	546	LYS	N-CA-C	5.74	120.52	112.72
1	A	201	ASP	N-CA-C	-5.74	103.84	111.24
1	D	197	LYS	N-CA-C	-5.73	105.39	112.38
1	D	366	PHE	N-CA-C	-5.73	104.40	111.33
1	E	546	LYS	N-CA-C	5.73	120.51	112.72
1	B	286	GLU	N-CA-C	-5.72	106.34	113.15
1	E	201	ASP	N-CA-C	-5.72	103.86	111.24
1	A	724	ARG	N-CA-C	-5.72	104.41	111.33
1	E	697	ILE	N-CA-C	-5.72	104.77	111.00
1	D	697	ILE	N-CA-C	-5.71	104.77	111.00
1	A	197	LYS	N-CA-C	-5.70	105.43	112.38
1	D	146	LYS	N-CA-C	5.70	122.93	110.80
2	S	74	ARG	N-CA-C	-5.70	105.21	111.82
1	B	197	LYS	N-CA-C	-5.69	105.44	112.38
2	Q	74	ARG	N-CA-C	-5.69	105.22	111.82
1	B	146	LYS	N-CA-C	5.69	122.91	110.80
1	B	201	ASP	N-CA-C	-5.69	103.90	111.24
2	Q	32	LEU	N-CA-C	-5.69	104.98	111.07
1	C	146	LYS	N-CA-C	5.68	122.91	110.80
1	C	359	PRO	O-C-N	5.68	128.59	122.17
1	E	146	LYS	N-CA-C	5.68	122.90	110.80
1	E	694	VAL	N-CA-C	-5.68	104.81	111.00
1	F	724	ARG	N-CA-C	-5.68	104.46	111.33
1	A	146	LYS	N-CA-C	5.68	122.90	110.80
1	F	146	LYS	N-CA-C	5.68	122.89	110.80
1	C	741	ILE	N-CA-C	-5.67	102.21	109.30
1	D	66	LEU	N-CA-C	5.67	117.46	111.28
2	T	74	ARG	N-CA-C	-5.67	105.25	111.82
1	C	197	LYS	N-CA-C	-5.67	105.47	112.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	567	THR	N-CA-C	-5.66	105.40	112.93
1	E	171	TYR	N-CA-C	-5.66	106.22	113.01
2	O	74	ARG	N-CA-C	-5.66	105.25	111.82
1	B	384	ASN	N-CA-C	-5.66	105.02	111.07
2	R	32	LEU	N-CA-C	-5.66	105.02	111.07
1	C	433	TYR	CB-CA-C	5.66	119.06	109.84
1	A	433	TYR	CB-CA-C	5.65	119.06	109.84
1	B	433	TYR	CB-CA-C	5.65	119.05	109.84
1	B	697	ILE	N-CA-C	-5.65	104.84	111.00
1	F	384	ASN	N-CA-C	-5.65	105.03	111.07
2	R	74	ARG	N-CA-C	-5.65	105.27	111.82
1	E	433	TYR	CB-CA-C	5.64	119.03	109.84
1	D	567	THR	N-CA-C	-5.64	105.43	112.93
2	P	74	ARG	N-CA-C	-5.63	105.29	111.82
1	B	78	LYS	N-CA-C	-5.63	104.83	110.97
1	F	171	TYR	N-CA-C	-5.63	106.25	113.01
1	F	694	VAL	N-CA-C	-5.63	104.86	111.00
1	F	433	TYR	CB-CA-C	5.62	119.00	109.84
1	D	433	TYR	CB-CA-C	5.62	118.99	109.84
1	D	741	ILE	N-CA-C	-5.60	102.30	109.30
1	B	546	LYS	N-CA-C	5.59	120.32	112.72
1	E	66	LEU	N-CA-C	5.58	117.44	111.36
1	A	546	LYS	N-CA-C	5.58	120.31	112.72
1	E	567	THR	N-CA-C	-5.57	105.52	112.93
1	B	72	THR	N-CA-C	-5.57	106.69	112.93
1	F	567	THR	N-CA-C	-5.57	105.53	112.93
1	F	741	ILE	N-CA-C	-5.56	102.35	109.30
1	E	741	ILE	N-CA-C	-5.55	102.36	109.30
1	E	384	ASN	N-CA-C	-5.55	105.14	111.07
1	B	308	VAL	CA-C-N	5.54	126.77	119.84
1	B	308	VAL	C-N-CA	5.54	126.77	119.84
1	D	78	LYS	N-CA-C	-5.54	104.93	110.97
1	B	741	ILE	N-CA-C	-5.53	102.39	109.30
1	F	72	THR	N-CA-C	-5.52	106.75	112.93
1	F	761	GLN	N-CA-C	-5.51	105.18	111.07
1	C	78	LYS	N-CA-C	-5.50	104.92	111.03
1	C	308	VAL	CA-C-N	5.50	126.72	119.84
1	C	308	VAL	C-N-CA	5.50	126.72	119.84
1	F	197	LYS	N-CA-C	-5.50	105.39	111.71
1	A	694	VAL	N-CA-C	-5.48	105.02	111.00
1	F	244	ALA	N-CA-C	-5.48	105.22	111.14
1	C	384	ASN	N-CA-C	-5.47	105.21	111.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	741	ILE	N-CA-C	-5.47	102.46	109.30
1	D	244	ALA	N-CA-C	-5.47	105.23	111.14
1	A	72	THR	N-CA-C	-5.47	106.81	112.93
1	B	427	ASP	N-CA-C	5.47	116.49	107.20
1	A	567	THR	N-CA-C	-5.46	105.66	112.93
1	D	72	THR	N-CA-C	-5.46	106.81	112.93
1	C	244	ALA	N-CA-C	-5.46	105.24	111.14
1	E	72	THR	N-CA-C	-5.46	106.81	112.93
1	C	72	THR	N-CA-C	-5.46	106.82	112.93
1	C	761	GLN	N-CA-C	-5.45	105.24	111.07
1	A	427	ASP	N-CA-C	5.43	116.44	107.20
1	E	78	LYS	N-CA-C	-5.43	105.05	110.97
1	C	697	ILE	N-CA-C	-5.43	105.08	111.00
1	F	308	VAL	CA-C-N	5.42	126.62	119.84
1	F	308	VAL	C-N-CA	5.42	126.62	119.84
1	A	308	VAL	CA-C-N	5.42	126.62	119.84
1	A	308	VAL	C-N-CA	5.42	126.62	119.84
1	A	78	LYS	N-CA-C	-5.42	105.01	111.03
1	E	197	LYS	N-CA-C	-5.41	105.48	111.71
1	C	379	ALA	N-CA-C	-5.41	105.46	111.36
1	D	761	GLN	N-CA-C	-5.41	105.29	111.07
1	C	72	THR	CA-C-N	5.40	128.30	120.79
1	C	72	THR	C-N-CA	5.40	128.30	120.79
1	D	308	VAL	CA-C-N	5.40	126.59	119.84
1	D	308	VAL	C-N-CA	5.40	126.59	119.84
1	F	78	LYS	N-CA-C	-5.40	105.03	111.03
1	A	384	ASN	N-CA-C	-5.39	105.30	111.07
1	B	105	TYR	N-CA-C	5.39	117.20	108.41
1	B	244	ALA	N-CA-C	-5.38	105.33	111.14
1	A	697	ILE	N-CA-C	-5.37	105.14	111.00
2	O	56	ASP	N-CA-C	5.37	117.00	108.67
1	B	694	VAL	N-CA-C	-5.36	105.16	111.00
1	E	244	ALA	N-CA-C	-5.35	105.36	111.14
1	E	308	VAL	CA-C-N	5.34	126.52	119.84
1	E	308	VAL	C-N-CA	5.34	126.52	119.84
2	P	23	GLY	N-CA-C	5.32	125.79	113.18
1	D	384	ASN	N-CA-C	-5.32	105.38	111.07
1	A	379	ALA	N-CA-C	-5.32	105.56	111.36
1	A	761	GLN	N-CA-C	-5.32	105.38	111.07
1	D	486	LYS	CA-C-N	5.32	125.61	119.92
1	D	486	LYS	C-N-CA	5.32	125.61	119.92
1	B	687	GLU	N-CA-C	-5.31	105.13	111.03

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	105	TYR	N-CA-C	5.31	117.07	108.41
1	E	379	ALA	N-CA-C	-5.31	105.57	111.36
1	E	228	ASN	CB-CA-C	5.31	119.55	110.74
2	O	23	GLY	N-CA-C	5.30	125.75	113.18
2	T	56	ASP	N-CA-C	5.30	116.89	108.67
2	T	112	LEU	N-CA-C	-5.30	106.66	113.02
2	P	56	ASP	N-CA-C	5.30	116.89	108.67
1	A	66	LEU	N-CA-C	5.29	117.13	111.36
1	A	228	ASN	CB-CA-C	5.29	119.52	110.74
1	B	228	ASN	CB-CA-C	5.29	119.52	110.74
1	C	228	ASN	CB-CA-C	5.29	119.52	110.74
1	F	228	ASN	CB-CA-C	5.29	119.52	110.74
2	O	87	GLU	N-CA-C	-5.28	105.98	112.90
2	T	87	GLU	N-CA-C	-5.28	105.98	112.90
1	D	228	ASN	CB-CA-C	5.28	119.51	110.74
1	B	379	ALA	N-CA-C	-5.27	105.62	111.36
2	T	23	GLY	N-CA-C	5.27	125.66	113.18
1	E	761	GLN	N-CA-C	-5.26	105.44	111.07
2	P	112	LEU	N-CA-C	-5.26	106.71	113.02
2	Q	23	GLY	N-CA-C	5.26	125.64	113.18
2	R	112	LEU	N-CA-C	-5.26	106.71	113.02
2	S	56	ASP	N-CA-C	5.26	116.82	108.67
1	B	761	GLN	N-CA-C	-5.25	105.45	111.07
2	R	56	ASP	N-CA-C	5.25	116.81	108.67
2	S	23	GLY	N-CA-C	5.25	125.61	113.18
1	C	486	LYS	CA-C-N	5.24	125.53	119.92
1	C	486	LYS	C-N-CA	5.24	125.53	119.92
2	R	23	GLY	N-CA-C	5.24	125.59	113.18
2	Q	56	ASP	N-CA-C	5.23	116.78	108.67
1	E	105	TYR	N-CA-C	5.22	116.92	108.41
1	D	149	THR	CB-CA-C	5.21	116.37	108.86
2	O	112	LEU	N-CA-C	-5.21	106.77	113.02
1	F	149	THR	CB-CA-C	5.21	116.36	108.86
1	A	149	THR	CB-CA-C	5.20	116.35	108.86
1	B	149	THR	CB-CA-C	5.20	116.35	108.86
1	C	149	THR	CB-CA-C	5.20	116.34	108.86
1	E	149	THR	CB-CA-C	5.20	116.34	108.86
1	B	66	LEU	N-CA-C	5.19	117.01	111.36
1	F	258	GLU	N-CA-C	-5.19	105.83	113.61
1	F	379	ALA	N-CA-C	-5.19	105.71	111.36
1	A	105	TYR	N-CA-C	5.18	116.85	108.41
1	D	379	ALA	N-CA-C	-5.17	105.55	111.14

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	450	ASN	N-CA-C	-5.17	101.26	109.59
1	F	486	LYS	CA-C-N	5.17	125.45	119.92
1	F	486	LYS	C-N-CA	5.17	125.45	119.92
1	B	545	THR	N-CA-C	5.16	117.90	111.24
1	C	450	ASN	N-CA-C	-5.15	101.29	109.59
1	D	105	TYR	N-CA-C	5.15	116.81	108.41
1	A	450	ASN	N-CA-C	-5.15	101.30	109.59
1	E	486	LYS	CA-C-N	5.15	125.43	119.92
1	E	486	LYS	C-N-CA	5.15	125.43	119.92
1	F	733	GLU	N-CA-C	-5.14	105.10	111.33
1	F	450	ASN	N-CA-C	-5.14	101.32	109.59
1	E	450	ASN	N-CA-C	-5.13	101.33	109.59
1	B	450	ASN	N-CA-C	-5.13	101.33	109.59
1	C	733	GLU	N-CA-C	-5.12	105.13	111.33
1	D	495	PHE	N-CA-C	-5.12	106.19	112.90
1	E	545	THR	N-CA-C	5.11	117.84	111.24
1	B	486	LYS	CA-C-N	5.11	125.39	119.92
1	B	486	LYS	C-N-CA	5.11	125.39	119.92
2	S	112	LEU	N-CA-C	-5.11	106.89	113.02
2	R	87	GLU	N-CA-C	-5.09	106.23	112.90
1	D	687	GLU	N-CA-C	-5.09	105.38	111.03
1	D	545	THR	N-CA-C	5.08	117.80	111.24
2	Q	112	LEU	N-CA-C	-5.08	106.92	113.02
1	A	733	GLU	N-CA-C	-5.08	105.19	111.33
1	C	687	GLU	N-CA-C	-5.07	105.40	111.03
1	F	135	VAL	N-CA-C	5.07	119.83	108.88
1	A	135	VAL	N-CA-C	5.06	119.81	108.88
1	D	135	VAL	N-CA-C	5.06	119.81	108.88
1	C	135	VAL	N-CA-C	5.05	119.80	108.88
1	E	135	VAL	N-CA-C	5.05	119.80	108.88
1	E	733	GLU	N-CA-C	-5.05	105.22	111.33
1	B	135	VAL	N-CA-C	5.04	119.78	108.88
1	E	82	THR	N-CA-C	-5.04	105.78	111.28
1	A	132	GLY	N-CA-C	-5.04	101.24	113.18
1	A	486	LYS	CA-C-N	5.04	125.31	119.92
1	A	486	LYS	C-N-CA	5.04	125.31	119.92
1	C	545	THR	N-CA-C	5.02	117.71	111.24
1	C	359	PRO	CA-C-O	5.01	127.30	120.03
1	D	733	GLU	N-CA-C	-5.01	105.27	111.33
1	D	360	VAL	N-CA-C	5.00	117.34	111.09
1	F	261	ALA	CA-C-N	5.00	124.49	119.19
1	F	261	ALA	C-N-CA	5.00	124.49	119.19

There are no chirality outliers.

There are no planarity outliers.

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	5992	0	6010	878	0
1	B	5992	0	6010	883	0
1	C	5992	0	6010	880	0
1	D	5992	0	6010	888	0
1	E	5992	0	6010	881	0
1	F	5992	0	6010	892	0
2	O	1146	0	1071	210	0
2	P	1146	0	1071	203	0
2	Q	1146	0	1071	212	0
2	R	1146	0	1071	210	0
2	S	1146	0	1071	214	0
2	T	1146	0	1071	212	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
3	E	1	0	0	0	0
3	F	1	0	0	0	0
4	O	3	0	0	0	0
4	P	3	0	0	0	0
4	Q	3	0	0	0	0
4	R	3	0	0	0	0
4	S	3	0	0	0	0
4	T	3	0	0	0	0
All	All	42852	0	42486	6413	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 75.

All (6413) 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:A:154:ILE:HG13	1:A:171:TYR:HE1	1.02	1.19
1:E:165:GLN:HE21	1:E:251:PRO:HG2	1.07	1.17
1:E:154:ILE:HG13	1:E:171:TYR:HE1	1.01	1.15
1:B:165:GLN:HE21	1:B:251:PRO:HG2	1.06	1.14
1:F:154:ILE:HG13	1:F:171:TYR:HE1	1.03	1.14
1:A:165:GLN:HE21	1:A:251:PRO:HG2	1.05	1.13
1:B:122:GLU:HG3	1:B:147:ARG:HB2	1.30	1.13
1:D:122:GLU:HG3	1:D:147:ARG:HB2	1.30	1.13
1:C:154:ILE:HG13	1:C:171:TYR:HE1	1.01	1.13
1:D:165:GLN:HE21	1:D:251:PRO:HG2	1.05	1.13
1:A:324:THR:HB	1:A:499:PRO:HA	1.31	1.12
1:B:154:ILE:HG13	1:B:171:TYR:HE1	0.98	1.11
1:E:188:LEU:HD23	1:E:188:LEU:H	1.14	1.10
1:E:324:THR:HB	1:E:499:PRO:HA	1.30	1.10
1:C:122:GLU:HG3	1:C:147:ARG:HB2	1.30	1.10
1:C:165:GLN:HE21	1:C:251:PRO:HG2	1.06	1.10
1:D:324:THR:HB	1:D:499:PRO:HA	1.33	1.10
1:B:188:LEU:HD23	1:B:188:LEU:H	1.14	1.10
1:B:408:LEU:HD12	1:B:408:LEU:H	1.16	1.09
2:O:55:VAL:HG21	2:O:67:GLU:OE1	1.53	1.09
1:D:123:GLU:HG2	1:D:124:GLU:H	1.14	1.09
1:D:154:ILE:HG13	1:D:171:TYR:HE1	1.01	1.09
1:F:165:GLN:HE21	1:F:251:PRO:HG2	1.06	1.09
1:A:122:GLU:HG3	1:A:147:ARG:HB2	1.30	1.09
1:B:324:THR:HB	1:B:499:PRO:HA	1.31	1.08
1:E:123:GLU:HG2	1:E:124:GLU:H	1.14	1.08
1:E:408:LEU:HD12	1:E:408:LEU:H	1.16	1.08
1:F:324:THR:HB	1:F:499:PRO:HA	1.32	1.08
1:F:188:LEU:HD23	1:F:188:LEU:H	1.14	1.08
1:A:123:GLU:HG2	1:A:124:GLU:H	1.14	1.08
1:A:188:LEU:HD23	1:A:188:LEU:H	1.14	1.08
1:B:154:ILE:HG13	1:B:171:TYR:CE1	1.89	1.08
1:E:122:GLU:HG3	1:E:147:ARG:HB2	1.30	1.07
1:B:550:SER:HB3	1:B:553:GLN:HG3	1.09	1.07
1:D:550:SER:HB3	1:D:553:GLN:HG3	1.11	1.07
1:E:154:ILE:HG13	1:E:171:TYR:CE1	1.90	1.07
1:E:550:SER:HB3	1:E:553:GLN:HG3	1.10	1.06
1:F:123:GLU:HG2	1:F:124:GLU:H	1.14	1.06
1:A:550:SER:HB3	1:A:553:GLN:HG3	1.10	1.06
1:F:550:SER:HB3	1:F:553:GLN:HG3	1.08	1.06
1:A:154:ILE:HG13	1:A:171:TYR:CE1	1.91	1.05
1:B:123:GLU:HG2	1:B:124:GLU:H	1.14	1.05

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:154:ILE:HG13	1:C:171:TYR:CE1	1.91	1.05
1:C:188:LEU:HD23	1:C:188:LEU:H	1.14	1.05
1:D:188:LEU:H	1:D:188:LEU:HD23	1.14	1.05
1:A:408:LEU:H	1:A:408:LEU:HD12	1.17	1.05
1:C:550:SER:HB3	1:C:553:GLN:HG3	1.09	1.05
1:D:154:ILE:HG13	1:D:171:TYR:CE1	1.91	1.05
1:F:154:ILE:HG13	1:F:171:TYR:CE1	1.92	1.05
1:C:123:GLU:HG2	1:C:124:GLU:H	1.13	1.04
1:F:122:GLU:HG3	1:F:147:ARG:HB2	1.30	1.04
1:F:408:LEU:H	1:F:408:LEU:HD12	1.16	1.04
1:B:360:VAL:O	1:B:360:VAL:HG22	1.53	1.04
1:C:324:THR:HB	1:C:499:PRO:HA	1.32	1.04
2:R:100:ILE:HB	2:R:136:VAL:HG23	1.40	1.03
2:S:100:ILE:HB	2:S:136:VAL:HG23	1.40	1.03
1:D:408:LEU:H	1:D:408:LEU:HD12	1.17	1.03
2:T:100:ILE:HB	2:T:136:VAL:HG23	1.41	1.02
1:A:187:SER:C	1:A:188:LEU:O	1.99	1.01
2:O:100:ILE:HB	2:O:136:VAL:HG23	1.41	1.01
2:Q:100:ILE:HB	2:Q:136:VAL:HG23	1.40	1.01
1:C:408:LEU:H	1:C:408:LEU:HD12	1.18	1.00
1:D:184:LYS:HE2	1:D:193:LEU:HD12	1.42	1.00
1:F:187:SER:C	1:F:188:LEU:O	1.99	1.00
2:P:100:ILE:HB	2:P:136:VAL:HG23	1.41	1.00
2:T:72:MET:O	2:T:76:MET:HB3	1.62	1.00
2:R:72:MET:O	2:R:76:MET:HB3	1.62	1.00
2:O:72:MET:O	2:O:76:MET:HB3	1.62	1.00
1:A:188:LEU:H	1:A:188:LEU:CD2	1.74	0.99
1:B:188:LEU:H	1:B:188:LEU:CD2	1.74	0.99
2:R:30:LYS:H	2:R:30:LYS:HD3	1.27	0.99
2:S:72:MET:O	2:S:76:MET:HB3	1.62	0.99
1:D:188:LEU:H	1:D:188:LEU:CD2	1.74	0.99
1:C:188:LEU:H	1:C:188:LEU:CD2	1.74	0.99
1:D:187:SER:C	1:D:188:LEU:O	1.99	0.99
1:F:188:LEU:H	1:F:188:LEU:CD2	1.74	0.98
1:E:188:LEU:H	1:E:188:LEU:CD2	1.74	0.98
1:C:187:SER:C	1:C:188:LEU:O	1.99	0.98
1:E:184:LYS:HE2	1:E:193:LEU:HD12	1.43	0.98
1:A:184:LYS:HE2	1:A:193:LEU:HD12	1.43	0.98
2:Q:72:MET:O	2:Q:76:MET:HB3	1.62	0.98
2:S:30:LYS:HD3	2:S:30:LYS:H	1.28	0.98
2:P:72:MET:O	2:P:76:MET:HB3	1.62	0.98

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:550:SER:CB	1:F:553:GLN:HG3	1.93	0.98
2:Q:30:LYS:HD3	2:Q:30:LYS:H	1.27	0.98
1:B:184:LYS:HE2	1:B:193:LEU:HD12	1.43	0.97
1:F:184:LYS:HE2	1:F:193:LEU:HD12	1.43	0.97
1:F:89:ILE:HG22	1:F:93:VAL:HG11	1.46	0.97
2:T:30:LYS:H	2:T:30:LYS:HD3	1.29	0.97
1:B:550:SER:CB	1:B:553:GLN:HG3	1.94	0.97
1:C:550:SER:CB	1:C:553:GLN:HG3	1.94	0.97
1:E:550:SER:CB	1:E:553:GLN:HG3	1.95	0.97
1:C:184:LYS:HE2	1:C:193:LEU:HD12	1.43	0.97
2:P:30:LYS:HD3	2:P:30:LYS:H	1.29	0.96
2:O:30:LYS:H	2:O:30:LYS:HD3	1.28	0.96
2:T:36:MET:HE3	2:T:43:PRO:HG3	1.47	0.96
1:B:581:GLN:NE2	1:B:629:ASN:H	1.63	0.96
1:D:550:SER:CB	1:D:553:GLN:HG3	1.96	0.96
2:S:36:MET:HE3	2:S:43:PRO:HG3	1.48	0.96
1:A:550:SER:CB	1:A:553:GLN:HG3	1.95	0.96
1:E:187:SER:C	1:E:188:LEU:O	1.99	0.96
2:Q:36:MET:HE3	2:Q:43:PRO:HG3	1.47	0.96
1:A:89:ILE:HG22	1:A:93:VAL:HG11	1.46	0.96
2:O:36:MET:HE3	2:O:43:PRO:HG3	1.48	0.96
2:P:36:MET:HE3	2:P:43:PRO:HG3	1.47	0.96
1:A:581:GLN:NE2	1:A:629:ASN:H	1.63	0.96
2:R:36:MET:HE3	2:R:43:PRO:HG3	1.48	0.96
1:A:217:LYS:HZ2	1:A:217:LYS:HB3	1.30	0.96
1:C:89:ILE:HG22	1:C:93:VAL:HG11	1.46	0.96
1:F:678:VAL:HG22	1:F:745:TYR:HE2	1.31	0.96
1:E:89:ILE:HG22	1:E:93:VAL:HG11	1.48	0.95
1:B:173:ILE:HG13	1:B:242:SER:HB3	1.48	0.95
1:B:217:LYS:HZ2	1:B:217:LYS:HB3	1.31	0.95
1:E:581:GLN:NE2	1:E:629:ASN:H	1.64	0.95
1:F:581:GLN:NE2	1:F:629:ASN:H	1.64	0.95
2:P:12:PHE:HE1	2:P:72:MET:HE3	1.32	0.95
1:A:715:GLU:HA	1:A:718:ARG:NH1	1.81	0.95
1:D:89:ILE:HG22	1:D:93:VAL:HG11	1.44	0.95
1:D:173:ILE:HG13	1:D:242:SER:HB3	1.49	0.95
1:B:678:VAL:HG22	1:B:745:TYR:HE2	1.32	0.95
1:E:173:ILE:HG13	1:E:242:SER:HB3	1.49	0.95
1:C:173:ILE:HG13	1:C:242:SER:HB3	1.49	0.95
1:C:161:ILE:HA	1:C:167:LYS:HD2	1.49	0.94
1:D:217:LYS:HZ2	1:D:217:LYS:HB3	1.32	0.94

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:715:GLU:HA	1:F:718:ARG:NH1	1.83	0.94
1:A:173:ILE:HG13	1:A:242:SER:HB3	1.49	0.94
1:D:581:GLN:NE2	1:D:629:ASN:H	1.64	0.94
1:F:173:ILE:HG13	1:F:242:SER:HB3	1.49	0.94
1:E:217:LYS:HZ2	1:E:217:LYS:HB3	1.32	0.94
1:C:678:VAL:HG22	1:C:745:TYR:HE2	1.32	0.94
1:D:678:VAL:HG22	1:D:745:TYR:HE2	1.32	0.94
1:E:715:GLU:HA	1:E:718:ARG:NH1	1.82	0.94
1:F:217:LYS:HZ2	1:F:217:LYS:HB3	1.33	0.94
1:D:715:GLU:HA	1:D:718:ARG:NH1	1.83	0.93
1:B:187:SER:C	1:B:188:LEU:O	1.99	0.93
1:B:325:TYR:HB2	1:B:498:ALA:HB3	1.50	0.93
1:E:767:GLN:HG2	1:E:768:LYS:HG2	1.51	0.93
1:C:715:GLU:HA	1:C:718:ARG:NH1	1.84	0.93
2:S:12:PHE:HE1	2:S:72:MET:HE3	1.33	0.93
1:A:122:GLU:CG	1:A:147:ARG:HB2	1.99	0.93
2:O:55:VAL:CG2	2:O:67:GLU:OE1	2.17	0.93
2:R:12:PHE:HE1	2:R:72:MET:HE3	1.34	0.93
1:B:122:GLU:CG	1:B:147:ARG:HB2	1.99	0.92
1:A:472:ARG:HB3	1:A:472:ARG:HH11	1.34	0.92
1:C:122:GLU:CG	1:C:147:ARG:HB2	1.99	0.92
1:D:122:GLU:CG	1:D:147:ARG:HB2	1.99	0.92
1:E:678:VAL:HG22	1:E:745:TYR:HE2	1.32	0.92
2:O:12:PHE:HE1	2:O:72:MET:HE3	1.34	0.92
1:A:678:VAL:HG22	1:A:745:TYR:HE2	1.33	0.92
1:B:715:GLU:HA	1:B:718:ARG:NH1	1.83	0.92
1:B:89:ILE:HG22	1:B:93:VAL:HG11	1.47	0.92
1:F:161:ILE:HA	1:F:167:LYS:HD2	1.49	0.92
1:A:325:TYR:HB2	1:A:498:ALA:HB3	1.51	0.92
1:C:581:GLN:NE2	1:C:629:ASN:H	1.66	0.92
1:D:165:GLN:NE2	1:D:251:PRO:HG2	1.84	0.92
1:F:767:GLN:HG2	1:F:768:LYS:HG2	1.51	0.92
1:F:122:GLU:CG	1:F:147:ARG:HB2	1.99	0.92
1:F:472:ARG:HB3	1:F:472:ARG:HH11	1.35	0.92
1:E:161:ILE:HA	1:E:167:LYS:HD2	1.49	0.92
1:B:161:ILE:HA	1:B:167:LYS:HD2	1.50	0.91
1:C:165:GLN:NE2	1:C:251:PRO:HG2	1.85	0.91
1:D:767:GLN:HG2	1:D:768:LYS:HG2	1.51	0.91
1:B:472:ARG:HB3	1:B:472:ARG:HH11	1.34	0.91
1:F:165:GLN:NE2	1:F:251:PRO:HG2	1.85	0.91
1:A:165:GLN:NE2	1:A:251:PRO:HG2	1.85	0.91

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:472:ARG:HB3	1:C:472:ARG:HH11	1.35	0.91
1:E:122:GLU:CG	1:E:147:ARG:HB2	1.99	0.91
2:Q:12:PHE:HE1	2:Q:72:MET:HE3	1.34	0.91
1:B:165:GLN:NE2	1:B:251:PRO:HG2	1.85	0.91
1:B:767:GLN:HG2	1:B:768:LYS:HG2	1.51	0.91
1:C:767:GLN:HG2	1:C:768:LYS:HG2	1.51	0.91
1:F:188:LEU:HD23	1:F:188:LEU:N	1.85	0.91
1:A:546:LYS:HD2	1:A:554:LYS:HE2	1.52	0.91
1:C:188:LEU:HD23	1:C:188:LEU:N	1.85	0.91
1:C:550:SER:HB3	1:C:553:GLN:CG	2.01	0.90
1:E:188:LEU:HD23	1:E:188:LEU:N	1.85	0.90
2:T:12:PHE:HE1	2:T:72:MET:HE3	1.33	0.90
1:E:325:TYR:HB2	1:E:498:ALA:HB3	1.51	0.90
1:A:479:LYS:HG2	1:A:488:LEU:HD21	1.53	0.90
1:B:154:ILE:CG1	1:B:171:TYR:HE1	1.83	0.90
1:C:217:LYS:HB3	1:C:217:LYS:HZ2	1.35	0.90
1:D:472:ARG:HB3	1:D:472:ARG:HH11	1.35	0.90
1:F:441:VAL:HG22	1:F:461:LYS:HG2	1.54	0.90
1:E:722:ILE:HG23	1:E:760:VAL:HG13	1.54	0.90
1:A:161:ILE:HA	1:A:167:LYS:HD2	1.51	0.90
1:D:479:LYS:HG2	1:D:488:LEU:HD21	1.54	0.90
1:C:154:ILE:CG1	1:C:171:TYR:HE1	1.85	0.90
1:D:441:VAL:HG22	1:D:461:LYS:HG2	1.54	0.90
1:A:767:GLN:HG2	1:A:768:LYS:HG2	1.51	0.89
1:E:154:ILE:CG1	1:E:171:TYR:HE1	1.84	0.89
1:E:472:ARG:HB3	1:E:472:ARG:HH11	1.34	0.89
1:F:360:VAL:HG11	1:F:370:LEU:HD22	1.51	0.89
2:S:12:PHE:CD1	2:S:72:MET:HG3	2.07	0.89
1:B:546:LYS:HD2	1:B:554:LYS:HE2	1.53	0.89
1:C:441:VAL:HG22	1:C:461:LYS:HG2	1.54	0.89
2:R:12:PHE:CD1	2:R:72:MET:HG3	2.07	0.89
2:T:12:PHE:CD1	2:T:72:MET:HG3	2.07	0.89
1:C:479:LYS:HG2	1:C:488:LEU:HD21	1.54	0.89
1:F:325:TYR:HB2	1:F:498:ALA:HB3	1.52	0.89
2:O:12:PHE:CD1	2:O:72:MET:HG3	2.07	0.89
1:A:441:VAL:HG22	1:A:461:LYS:HG2	1.54	0.89
1:C:409:ARG:NE	1:C:413:LEU:HD21	1.87	0.89
1:D:154:ILE:CG1	1:D:171:TYR:HE1	1.85	0.89
1:D:161:ILE:HA	1:D:167:LYS:HD2	1.53	0.89
1:E:441:VAL:HG22	1:E:461:LYS:HG2	1.54	0.89
1:B:409:ARG:NE	1:B:413:LEU:HD21	1.87	0.89

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:479:LYS:HG2	1:E:488:LEU:HD21	1.54	0.89
1:F:722:ILE:HG23	1:F:760:VAL:HG13	1.55	0.89
1:A:722:ILE:HG23	1:A:760:VAL:HG13	1.54	0.89
1:B:188:LEU:HD23	1:B:188:LEU:N	1.85	0.89
2:O:58:ASP:HB3	2:O:62:THR:HG23	1.54	0.89
1:F:546:LYS:HD2	1:F:554:LYS:HE2	1.53	0.89
1:B:360:VAL:O	1:B:360:VAL:CG2	2.18	0.88
1:B:479:LYS:HG2	1:B:488:LEU:HD21	1.54	0.88
1:B:550:SER:HB3	1:B:553:GLN:CG	2.01	0.88
1:E:275:GLY:HA2	1:E:278:LYS:HG3	1.55	0.88
2:Q:12:PHE:CD1	2:Q:72:MET:HG3	2.08	0.88
1:D:325:TYR:HB2	1:D:498:ALA:HB3	1.54	0.88
1:D:409:ARG:NE	1:D:413:LEU:HD21	1.88	0.88
1:D:722:ILE:HG23	1:D:760:VAL:HG13	1.55	0.88
2:Q:63:ILE:HG13	2:Q:67:GLU:HB3	1.55	0.88
1:A:154:ILE:CG1	1:A:171:TYR:HE1	1.85	0.88
1:E:409:ARG:NE	1:E:413:LEU:HD21	1.87	0.88
1:C:360:VAL:HG11	1:C:370:LEU:HD22	1.54	0.88
1:F:275:GLY:HA2	1:F:278:LYS:HG3	1.56	0.88
2:P:12:PHE:CD1	2:P:72:MET:HG3	2.08	0.88
1:E:165:GLN:NE2	1:E:251:PRO:HG2	1.86	0.88
2:P:58:ASP:HB3	2:P:62:THR:HG23	1.55	0.88
1:A:409:ARG:NE	1:A:413:LEU:HD21	1.87	0.88
1:C:546:LYS:HD2	1:C:554:LYS:HE2	1.54	0.88
1:F:479:LYS:HG2	1:F:488:LEU:HD21	1.55	0.88
1:C:275:GLY:HA2	1:C:278:LYS:HG3	1.55	0.88
1:E:546:LYS:HD2	1:E:554:LYS:HE2	1.54	0.88
2:P:63:ILE:HG13	2:P:67:GLU:HB3	1.56	0.88
1:C:722:ILE:HG23	1:C:760:VAL:HG13	1.56	0.88
1:E:550:SER:HB3	1:E:553:GLN:CG	2.02	0.88
1:B:441:VAL:HG22	1:B:461:LYS:HG2	1.54	0.87
1:C:325:TYR:HB2	1:C:498:ALA:HB3	1.52	0.87
1:F:154:ILE:CG1	1:F:171:TYR:HE1	1.86	0.87
1:F:409:ARG:NE	1:F:413:LEU:HD21	1.88	0.87
2:S:63:ILE:HG13	2:S:67:GLU:HB3	1.56	0.87
2:T:58:ASP:HB3	2:T:62:THR:HG23	1.55	0.87
1:B:722:ILE:HG23	1:B:760:VAL:HG13	1.55	0.87
1:F:550:SER:HB3	1:F:553:GLN:CG	2.00	0.87
1:D:546:LYS:HD2	1:D:554:LYS:HE2	1.54	0.87
1:B:275:GLY:HA2	1:B:278:LYS:HG3	1.56	0.87
2:R:63:ILE:HG13	2:R:67:GLU:HB3	1.57	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:678:VAL:HG22	1:F:745:TYR:CE2	2.08	0.87
1:A:188:LEU:HD23	1:A:188:LEU:N	1.85	0.87
1:B:678:VAL:HG22	1:B:745:TYR:CE2	2.09	0.87
1:E:324:THR:HB	1:E:499:PRO:CA	2.06	0.86
1:C:678:VAL:HG22	1:C:745:TYR:CE2	2.09	0.86
1:D:275:GLY:HA2	1:D:278:LYS:HG3	1.56	0.86
1:D:678:VAL:HG22	1:D:745:TYR:CE2	2.10	0.86
2:S:58:ASP:HB3	2:S:62:THR:HG23	1.55	0.86
2:R:58:ASP:HB3	2:R:62:THR:HG23	1.55	0.86
1:A:678:VAL:HG22	1:A:745:TYR:CE2	2.10	0.86
1:E:360:VAL:HG11	1:E:370:LEU:HD22	1.58	0.86
1:A:275:GLY:HA2	1:A:278:LYS:HG3	1.56	0.86
2:P:49:GLN:HE21	2:P:49:GLN:N	1.74	0.86
1:A:615:ILE:HG23	1:A:619:ILE:HD12	1.57	0.85
1:B:324:THR:HB	1:B:499:PRO:CA	2.06	0.85
1:B:355:SER:HB2	1:B:371:SER:HA	1.59	0.85
1:E:678:VAL:HG22	1:E:745:TYR:CE2	2.09	0.85
1:F:324:THR:HB	1:F:499:PRO:CA	2.07	0.85
2:O:63:ILE:HG13	2:O:67:GLU:HB3	1.56	0.85
2:Q:58:ASP:HB3	2:Q:62:THR:HG23	1.55	0.85
1:D:550:SER:HB3	1:D:553:GLN:CG	2.02	0.85
1:B:360:VAL:HG11	1:B:370:LEU:HD22	1.56	0.85
1:E:355:SER:HB2	1:E:371:SER:HA	1.59	0.85
2:P:12:PHE:CE1	2:P:72:MET:HE3	2.11	0.85
2:T:63:ILE:HG13	2:T:67:GLU:HB3	1.56	0.85
1:E:302:LEU:HD22	1:E:602:PHE:CE1	2.12	0.85
1:E:615:ILE:HG23	1:E:619:ILE:HD12	1.58	0.85
1:D:165:GLN:HE21	1:D:251:PRO:CG	1.88	0.84
2:S:49:GLN:N	2:S:49:GLN:HE21	1.74	0.84
1:B:302:LEU:HD22	1:B:602:PHE:CE1	2.13	0.84
2:Q:12:PHE:CE1	2:Q:72:MET:HE3	2.12	0.84
2:R:12:PHE:CE1	2:R:72:MET:HE3	2.12	0.84
1:D:360:VAL:HG11	1:D:370:LEU:HD22	1.59	0.84
1:D:615:ILE:HG23	1:D:619:ILE:HD12	1.56	0.84
1:F:153:ILE:O	1:F:154:ILE:HD13	1.77	0.84
2:P:49:GLN:HE21	2:P:49:GLN:H	1.23	0.84
1:D:302:LEU:HD22	1:D:602:PHE:CE1	2.13	0.84
1:B:517:VAL:HB	1:B:525:LYS:HZ1	1.43	0.84
1:C:270:LYS:O	1:C:273:LYS:HB2	1.77	0.84
1:C:324:THR:HB	1:C:499:PRO:CA	2.08	0.84
2:T:49:GLN:N	2:T:49:GLN:HE21	1.74	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:153:ILE:O	1:A:154:ILE:HD13	1.78	0.84
1:A:302:LEU:HD22	1:A:602:PHE:HE1	1.43	0.84
1:C:302:LEU:HD22	1:C:602:PHE:CE1	2.12	0.84
1:D:188:LEU:HD23	1:D:188:LEU:N	1.85	0.84
1:A:302:LEU:HD22	1:A:602:PHE:CE1	2.12	0.84
1:C:517:VAL:HB	1:C:525:LYS:HZ1	1.42	0.84
1:D:324:THR:HB	1:D:499:PRO:CA	2.07	0.84
1:A:360:VAL:HG11	1:A:370:LEU:HD22	1.59	0.84
2:T:12:PHE:CE1	2:T:72:MET:HE3	2.12	0.84
1:F:355:SER:HB2	1:F:371:SER:HA	1.60	0.84
1:A:324:THR:HB	1:A:499:PRO:CA	2.07	0.84
1:B:615:ILE:HG23	1:B:619:ILE:HD12	1.57	0.84
1:C:355:SER:HB2	1:C:371:SER:HA	1.60	0.84
1:A:355:SER:HB2	1:A:371:SER:HA	1.60	0.83
1:C:165:GLN:HE21	1:C:251:PRO:CG	1.89	0.83
1:F:165:GLN:HE21	1:F:251:PRO:CG	1.89	0.83
2:O:49:GLN:N	2:O:49:GLN:HE21	1.75	0.83
1:D:517:VAL:HB	1:D:525:LYS:HZ1	1.43	0.83
1:F:302:LEU:HD22	1:F:602:PHE:CE1	2.13	0.83
1:F:517:VAL:HB	1:F:525:LYS:HZ1	1.43	0.83
1:F:615:ILE:HG23	1:F:619:ILE:HD12	1.59	0.83
1:F:360:VAL:HG22	1:F:360:VAL:O	1.76	0.83
1:B:165:GLN:HE21	1:B:251:PRO:CG	1.89	0.83
2:S:12:PHE:CE1	2:S:72:MET:HE3	2.11	0.83
1:D:302:LEU:HD22	1:D:602:PHE:HE1	1.43	0.83
2:Q:49:GLN:N	2:Q:49:GLN:HE21	1.74	0.83
1:F:657:ILE:HG13	1:F:756:ILE:HD13	1.60	0.83
2:T:49:GLN:HE21	2:T:49:GLN:H	1.23	0.83
1:E:270:LYS:O	1:E:273:LYS:HB2	1.78	0.83
1:E:517:VAL:HB	1:E:525:LYS:HZ1	1.42	0.83
2:O:49:GLN:HE21	2:O:49:GLN:H	1.24	0.83
1:C:302:LEU:HD22	1:C:602:PHE:HE1	1.42	0.83
2:O:12:PHE:CE1	2:O:72:MET:HE3	2.13	0.83
2:O:37:ARG:HA	2:O:41:GLN:O	1.79	0.83
1:B:270:LYS:O	1:B:273:LYS:HB2	1.79	0.83
1:D:729:TYR:HB2	1:D:756:ILE:HG21	1.61	0.83
1:A:270:LYS:O	1:A:273:LYS:HB2	1.79	0.82
1:D:355:SER:HB2	1:D:371:SER:HA	1.60	0.82
2:Q:49:GLN:HE21	2:Q:49:GLN:H	1.23	0.82
1:A:165:GLN:HE21	1:A:251:PRO:CG	1.89	0.82
1:B:302:LEU:HD22	1:B:602:PHE:HE1	1.43	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:729:TYR:HB2	1:A:756:ILE:HG21	1.61	0.82
1:B:354:SER:O	1:B:371:SER:HB2	1.79	0.82
1:C:153:ILE:O	1:C:154:ILE:HD13	1.79	0.82
1:C:615:ILE:HG23	1:C:619:ILE:HD12	1.58	0.82
1:C:728:ALA:O	1:C:732:ILE:HG12	1.79	0.82
1:D:270:LYS:O	1:D:273:LYS:HB2	1.79	0.82
1:D:776:LEU:O	1:D:776:LEU:HD23	1.79	0.82
1:E:153:ILE:O	1:E:154:ILE:HD13	1.79	0.82
2:S:49:GLN:HE21	2:S:49:GLN:H	1.23	0.82
2:P:37:ARG:HA	2:P:41:GLN:O	1.80	0.82
1:B:657:ILE:HG13	1:B:756:ILE:HD13	1.60	0.82
1:C:462:ILE:HG12	1:C:463:THR:H	1.44	0.82
1:A:776:LEU:O	1:A:776:LEU:HD23	1.79	0.82
1:E:165:GLN:HE21	1:E:251:PRO:CG	1.90	0.82
1:B:236:GLU:HA	1:B:239:HIS:HD2	1.45	0.82
1:C:236:GLU:HA	1:C:239:HIS:HD2	1.45	0.82
1:E:236:GLU:HA	1:E:239:HIS:HD2	1.45	0.82
1:E:462:ILE:HG12	1:E:463:THR:H	1.45	0.82
1:F:728:ALA:O	1:F:732:ILE:HG12	1.80	0.82
2:R:49:GLN:N	2:R:49:GLN:HE21	1.75	0.82
1:C:776:LEU:HD23	1:C:776:LEU:O	1.80	0.82
1:F:305:SER:OG	1:F:307:LEU:HD13	1.80	0.82
1:E:354:SER:O	1:E:371:SER:HB2	1.80	0.82
1:E:657:ILE:HG13	1:E:756:ILE:HD13	1.62	0.82
2:T:37:ARG:HA	2:T:41:GLN:O	1.80	0.82
1:F:270:LYS:O	1:F:273:LYS:HB2	1.79	0.81
1:F:302:LEU:HD22	1:F:602:PHE:HE1	1.43	0.81
2:S:37:ARG:HA	2:S:41:GLN:O	1.80	0.81
1:A:236:GLU:HA	1:A:239:HIS:HD2	1.45	0.81
1:B:179:LEU:HD23	1:B:179:LEU:H	1.45	0.81
1:C:729:TYR:HB2	1:C:756:ILE:HG21	1.61	0.81
1:A:497:LEU:HD13	1:A:556:MET:HG2	1.62	0.81
1:A:728:ALA:O	1:A:732:ILE:HG12	1.78	0.81
1:D:153:ILE:O	1:D:154:ILE:HD13	1.79	0.81
1:D:462:ILE:HG12	1:D:463:THR:H	1.45	0.81
1:E:302:LEU:HB2	1:E:602:PHE:HD1	1.46	0.81
2:R:49:GLN:HE21	2:R:49:GLN:H	1.24	0.81
1:B:497:LEU:HD13	1:B:556:MET:HG2	1.63	0.81
1:B:776:LEU:O	1:B:776:LEU:HD23	1.79	0.81
1:C:540:ARG:NH2	2:Q:87:GLU:OE1	2.12	0.81
2:R:102:ALA:HB2	2:R:125:ILE:HG13	1.63	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:123:GLU:CG	1:C:124:GLU:H	1.94	0.81
1:A:462:ILE:HG12	1:A:463:THR:H	1.45	0.81
1:B:302:LEU:HB2	1:B:602:PHE:HD1	1.46	0.81
1:C:354:SER:O	1:C:371:SER:HB2	1.79	0.81
2:O:102:ALA:HB2	2:O:125:ILE:HG13	1.63	0.81
1:B:462:ILE:HG12	1:B:463:THR:H	1.44	0.81
1:C:462:ILE:HG12	1:C:463:THR:N	1.96	0.81
2:R:48:LEU:HA	2:R:51:MET:HE2	1.63	0.81
1:A:354:SER:O	1:A:371:SER:HB2	1.80	0.81
1:B:196:ILE:HA	1:B:199:LEU:HD12	1.62	0.81
1:E:302:LEU:HD22	1:E:602:PHE:HE1	1.42	0.81
1:A:517:VAL:HB	1:A:525:LYS:HZ1	1.43	0.81
1:A:550:SER:HB3	1:A:553:GLN:CG	2.02	0.81
1:E:179:LEU:HD23	1:E:179:LEU:H	1.45	0.81
1:E:305:SER:OG	1:E:307:LEU:HD13	1.81	0.81
1:B:93:VAL:HA	1:B:96:ILE:HD12	1.63	0.81
1:B:153:ILE:O	1:B:154:ILE:HD13	1.79	0.81
1:B:305:SER:OG	1:B:307:LEU:HD13	1.81	0.81
1:B:729:TYR:HB2	1:B:756:ILE:HG21	1.61	0.81
1:D:354:SER:O	1:D:371:SER:HB2	1.80	0.81
1:E:462:ILE:HG12	1:E:463:THR:N	1.96	0.81
1:E:729:TYR:HB2	1:E:756:ILE:HG21	1.62	0.81
1:A:657:ILE:HG13	1:A:756:ILE:HD13	1.62	0.80
1:D:236:GLU:HA	1:D:239:HIS:HD2	1.45	0.80
1:D:597:ASN:HB2	1:D:598:PRO:HD2	1.62	0.80
1:E:123:GLU:CG	1:E:124:GLU:H	1.94	0.80
1:E:776:LEU:HD23	1:E:776:LEU:O	1.80	0.80
1:F:236:GLU:HA	1:F:239:HIS:HD2	1.46	0.80
1:F:354:SER:O	1:F:371:SER:HB2	1.79	0.80
1:C:360:VAL:HG22	1:C:360:VAL:O	1.82	0.80
1:D:728:ALA:O	1:D:732:ILE:HG12	1.80	0.80
1:F:462:ILE:HG12	1:F:463:THR:N	1.96	0.80
1:F:497:LEU:HD13	1:F:556:MET:HG2	1.63	0.80
1:F:540:ARG:NH2	2:T:87:GLU:OE1	2.14	0.80
2:Q:37:ARG:HA	2:Q:41:GLN:O	1.81	0.80
1:E:629:ASN:HB3	1:E:632:TYR:CE1	2.16	0.80
1:F:318:ILE:H	1:F:318:ILE:HD12	1.46	0.80
2:S:48:LEU:HA	2:S:51:MET:HE2	1.63	0.80
1:A:462:ILE:HG12	1:A:463:THR:N	1.96	0.80
1:D:462:ILE:HG12	1:D:463:THR:N	1.97	0.80
1:E:318:ILE:H	1:E:318:ILE:HD12	1.46	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:302:LEU:HB2	1:F:602:PHE:HD1	1.46	0.80
1:A:345:THR:HB	1:A:491:ASP:HB3	1.64	0.80
1:E:728:ALA:O	1:E:732:ILE:HG12	1.80	0.80
1:F:776:LEU:HD23	1:F:776:LEU:O	1.79	0.80
2:P:102:ALA:HB2	2:P:125:ILE:HG13	1.63	0.80
1:A:179:LEU:HD23	1:A:179:LEU:H	1.45	0.80
1:C:657:ILE:HG13	1:C:756:ILE:HD13	1.61	0.80
1:E:387:ASN:HD22	1:E:387:ASN:N	1.80	0.80
1:F:196:ILE:HA	1:F:199:LEU:HD12	1.63	0.80
1:A:93:VAL:HA	1:A:96:ILE:HD12	1.64	0.80
1:A:122:GLU:HG3	1:A:147:ARG:CB	2.11	0.80
1:A:196:ILE:HA	1:A:199:LEU:HD12	1.63	0.80
1:D:345:THR:HB	1:D:491:ASP:HB3	1.64	0.80
1:D:657:ILE:HG13	1:D:756:ILE:HD13	1.62	0.80
1:E:93:VAL:HA	1:E:96:ILE:HD12	1.63	0.80
2:O:13:LYS:NZ	2:O:65:PHE:HB3	1.97	0.80
2:R:13:LYS:NZ	2:R:65:PHE:HB3	1.96	0.80
1:A:302:LEU:HB2	1:A:602:PHE:HD1	1.46	0.80
1:D:179:LEU:HD23	1:D:179:LEU:H	1.45	0.80
1:D:93:VAL:HA	1:D:96:ILE:HD12	1.63	0.80
1:B:345:THR:HB	1:B:491:ASP:HB3	1.64	0.80
1:C:497:LEU:HD13	1:C:556:MET:HG2	1.64	0.80
1:C:597:ASN:HB2	1:C:598:PRO:HD2	1.63	0.80
1:F:123:GLU:HG2	1:F:124:GLU:N	1.97	0.80
1:B:728:ALA:O	1:B:732:ILE:HG12	1.80	0.79
1:E:196:ILE:HA	1:E:199:LEU:HD12	1.62	0.79
1:F:413:LEU:HB2	1:F:419:ILE:HG12	1.64	0.79
2:P:13:LYS:NZ	2:P:65:PHE:HB3	1.97	0.79
2:P:48:LEU:HA	2:P:51:MET:HE2	1.64	0.79
2:Q:102:ALA:HB2	2:Q:125:ILE:HG13	1.64	0.79
1:C:408:LEU:H	1:C:408:LEU:CD1	1.96	0.79
1:C:179:LEU:HD23	1:C:179:LEU:H	1.46	0.79
1:C:302:LEU:HB2	1:C:602:PHE:HD1	1.46	0.79
1:F:345:THR:HB	1:F:491:ASP:HB3	1.64	0.79
1:A:413:LEU:HB2	1:A:419:ILE:HG12	1.64	0.79
1:D:134:LYS:O	1:D:135:VAL:HG12	1.83	0.79
1:D:302:LEU:HB2	1:D:602:PHE:HD1	1.46	0.79
1:D:413:LEU:HB2	1:D:419:ILE:HG12	1.65	0.79
1:A:540:ARG:NH2	2:O:87:GLU:OE1	2.15	0.79
1:C:318:ILE:H	1:C:318:ILE:HD12	1.47	0.79
1:C:345:THR:HB	1:C:491:ASP:HB3	1.64	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:413:LEU:HB2	1:C:419:ILE:HG12	1.64	0.79
1:D:694:VAL:CG2	2:R:18:LEU:HD21	2.13	0.79
1:F:179:LEU:HD23	1:F:179:LEU:H	1.46	0.79
1:F:301:ALA:C	1:F:303:LYS:H	1.90	0.79
2:Q:48:LEU:HA	2:Q:51:MET:HE2	1.64	0.79
1:B:123:GLU:HG2	1:B:124:GLU:N	1.96	0.79
1:B:318:ILE:H	1:B:318:ILE:HD12	1.47	0.79
1:C:305:SER:OG	1:C:307:LEU:HD13	1.83	0.79
1:F:629:ASN:HB3	1:F:632:TYR:CE1	2.17	0.79
1:A:305:SER:OG	1:A:307:LEU:HD13	1.83	0.79
1:B:134:LYS:O	1:B:135:VAL:HG12	1.83	0.79
1:C:79:ILE:C	1:C:81:GLN:H	1.90	0.79
1:E:497:LEU:HD13	1:E:556:MET:HG2	1.63	0.79
2:R:37:ARG:HA	2:R:41:GLN:O	1.81	0.79
2:S:13:LYS:NZ	2:S:65:PHE:HB3	1.97	0.79
2:T:13:LYS:NZ	2:T:65:PHE:HB3	1.98	0.79
2:T:102:ALA:HB2	2:T:125:ILE:HG13	1.64	0.79
1:D:122:GLU:HG3	1:D:147:ARG:CB	2.11	0.79
2:T:92:PHE:O	2:T:94:LYS:N	2.16	0.79
1:A:79:ILE:C	1:A:81:GLN:H	1.90	0.78
1:A:123:GLU:HG2	1:A:124:GLU:N	1.97	0.78
1:B:79:ILE:C	1:B:81:GLN:H	1.91	0.78
1:D:295:VAL:C	1:D:296:LEU:HD23	2.09	0.78
1:B:462:ILE:HG12	1:B:463:THR:N	1.96	0.78
1:C:93:VAL:HA	1:C:96:ILE:HD12	1.63	0.78
1:C:134:LYS:O	1:C:135:VAL:HG12	1.83	0.78
1:C:196:ILE:HA	1:C:199:LEU:HD12	1.64	0.78
1:D:497:LEU:HD13	1:D:556:MET:HG2	1.63	0.78
1:D:710:HIS:C	1:D:712:PHE:H	1.91	0.78
1:F:123:GLU:CG	1:F:124:GLU:H	1.95	0.78
1:C:295:VAL:C	1:C:296:LEU:HD23	2.08	0.78
1:E:413:LEU:HB2	1:E:419:ILE:HG12	1.64	0.78
2:T:48:LEU:HA	2:T:51:MET:HE2	1.64	0.78
1:B:112:VAL:HG12	1:B:113:GLU:H	1.49	0.78
1:B:597:ASN:HB2	1:B:598:PRO:HD2	1.65	0.78
1:B:629:ASN:HB3	1:B:632:TYR:CE1	2.19	0.78
1:C:301:ALA:C	1:C:303:LYS:H	1.90	0.78
1:F:93:VAL:HA	1:F:96:ILE:HD12	1.64	0.78
1:F:597:ASN:HB2	1:F:598:PRO:HD2	1.64	0.78
1:F:597:ASN:HD21	1:F:601:GLU:H	1.30	0.78
2:O:48:LEU:HA	2:O:51:MET:HE2	1.63	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:318:ILE:HD12	1:D:318:ILE:H	1.47	0.78
1:E:79:ILE:C	1:E:81:GLN:H	1.91	0.78
1:F:397:GLU:HA	1:F:480:ASN:HB2	1.65	0.78
2:S:102:ALA:HB2	2:S:125:ILE:HG13	1.64	0.78
1:D:301:ALA:C	1:D:303:LYS:H	1.90	0.78
1:E:597:ASN:HB2	1:E:598:PRO:HD2	1.65	0.78
1:F:462:ILE:HG12	1:F:463:THR:H	1.45	0.78
1:A:134:LYS:O	1:A:135:VAL:HG12	1.83	0.78
1:A:478:ALA:HB1	1:A:486:LYS:O	1.84	0.78
1:B:122:GLU:HG3	1:B:147:ARG:CB	2.11	0.78
1:B:387:ASN:HD22	1:B:387:ASN:N	1.81	0.78
1:D:597:ASN:HD21	1:D:601:GLU:H	1.31	0.78
1:E:345:THR:HB	1:E:491:ASP:HB3	1.65	0.78
1:B:295:VAL:C	1:B:296:LEU:HD23	2.08	0.78
1:D:196:ILE:HA	1:D:199:LEU:HD12	1.63	0.78
1:E:301:ALA:C	1:E:303:LYS:H	1.90	0.78
1:F:387:ASN:HD22	1:F:387:ASN:N	1.81	0.78
1:A:282:SER:HA	1:A:285:LYS:HD3	1.65	0.78
1:B:581:GLN:HE21	1:B:629:ASN:H	1.31	0.78
1:C:710:HIS:C	1:C:712:PHE:H	1.91	0.78
1:D:123:GLU:HG2	1:D:124:GLU:N	1.97	0.78
1:B:540:ARG:NH2	2:P:87:GLU:OE1	2.16	0.78
1:C:122:GLU:HG3	1:C:147:ARG:CB	2.11	0.78
1:D:305:SER:OG	1:D:307:LEU:HD13	1.84	0.78
1:D:716:LYS:O	1:D:720:ILE:HG22	1.84	0.78
1:E:397:GLU:HA	1:E:480:ASN:HB2	1.65	0.78
2:Q:13:LYS:NZ	2:Q:65:PHE:HB3	1.98	0.78
1:A:141:PHE:HD1	1:A:141:PHE:H	1.32	0.77
1:B:413:LEU:HB2	1:B:419:ILE:HG12	1.64	0.77
1:C:123:GLU:HG2	1:C:124:GLU:N	1.96	0.77
1:F:729:TYR:HB2	1:F:756:ILE:HG21	1.64	0.77
1:B:397:GLU:HA	1:B:480:ASN:HB2	1.65	0.77
1:B:746:LYS:O	1:B:750:GLN:HG2	1.85	0.77
1:D:282:SER:HA	1:D:285:LYS:HD3	1.65	0.77
1:D:629:ASN:HB3	1:D:632:TYR:CE1	2.18	0.77
1:E:540:ARG:NH2	2:S:87:GLU:OE1	2.16	0.77
1:F:295:VAL:C	1:F:296:LEU:HD23	2.09	0.77
1:A:318:ILE:HD12	1:A:318:ILE:H	1.47	0.77
1:A:710:HIS:C	1:A:712:PHE:H	1.91	0.77
1:C:629:ASN:HB3	1:C:632:TYR:CE1	2.18	0.77
1:A:597:ASN:HD21	1:A:601:GLU:H	1.30	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:257:LEU:O	1:D:257:LEU:HG	1.83	0.77
1:E:134:LYS:O	1:E:135:VAL:HG12	1.83	0.77
1:E:295:VAL:C	1:E:296:LEU:HD23	2.09	0.77
1:F:134:LYS:O	1:F:135:VAL:HG12	1.83	0.77
1:F:282:SER:HA	1:F:285:LYS:HD3	1.64	0.77
1:B:282:SER:HA	1:B:285:LYS:HD3	1.65	0.77
1:B:301:ALA:C	1:B:303:LYS:H	1.91	0.77
1:C:694:VAL:CG2	2:Q:18:LEU:HD21	2.15	0.77
1:E:581:GLN:HE21	1:E:629:ASN:H	1.32	0.77
1:A:387:ASN:HD22	1:A:387:ASN:N	1.81	0.77
1:A:397:GLU:HA	1:A:480:ASN:HB2	1.66	0.77
1:A:629:ASN:HB3	1:A:632:TYR:CE1	2.19	0.77
1:A:722:ILE:HG23	1:A:760:VAL:CG1	2.15	0.77
1:D:397:GLU:HA	1:D:480:ASN:HB2	1.66	0.77
1:A:694:VAL:CG2	2:O:18:LEU:HD21	2.15	0.77
1:D:540:ARG:NH2	2:R:87:GLU:OE1	2.17	0.77
1:E:112:VAL:HG12	1:E:113:GLU:H	1.50	0.77
1:E:122:GLU:HG3	1:E:147:ARG:CB	2.11	0.77
1:F:79:ILE:C	1:F:81:GLN:H	1.91	0.77
1:A:472:ARG:HB3	1:A:472:ARG:NH1	1.99	0.77
1:C:282:SER:HA	1:C:285:LYS:HD3	1.65	0.77
1:E:141:PHE:HD1	1:E:141:PHE:H	1.32	0.77
1:E:257:LEU:HG	1:E:257:LEU:O	1.83	0.77
2:O:92:PHE:O	2:O:94:LYS:N	2.18	0.77
1:B:597:ASN:HD21	1:B:601:GLU:H	1.31	0.77
1:B:694:VAL:CG2	2:P:18:LEU:HD21	2.14	0.77
1:C:184:LYS:HE3	1:C:191:GLU:HB2	1.67	0.77
1:D:77:ASP:O	1:D:81:GLN:HB2	1.85	0.77
1:D:79:ILE:C	1:D:81:GLN:H	1.91	0.77
1:D:387:ASN:N	1:D:387:ASN:HD22	1.80	0.77
1:F:694:VAL:CG2	2:T:18:LEU:HD21	2.15	0.77
1:F:710:HIS:C	1:F:712:PHE:H	1.92	0.77
1:A:301:ALA:C	1:A:303:LYS:H	1.91	0.77
1:A:597:ASN:HB2	1:A:598:PRO:HD2	1.64	0.77
1:B:472:ARG:HB3	1:B:472:ARG:NH1	2.00	0.77
1:C:112:VAL:HG12	1:C:113:GLU:H	1.49	0.77
1:C:746:LYS:O	1:C:750:GLN:HG2	1.85	0.77
1:D:746:LYS:O	1:D:750:GLN:HG2	1.85	0.77
1:F:122:GLU:HG3	1:F:147:ARG:CB	2.11	0.77
1:F:372:LYS:HG3	1:F:373:LYS:N	2.00	0.77
1:C:397:GLU:HA	1:C:480:ASN:HB2	1.65	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:478:ALA:HB1	1:C:486:LYS:O	1.85	0.76
1:E:282:SER:HA	1:E:285:LYS:HD3	1.65	0.76
1:E:597:ASN:HD21	1:E:601:GLU:H	1.32	0.76
1:A:295:VAL:C	1:A:296:LEU:HD23	2.09	0.76
1:D:112:VAL:HG12	1:D:113:GLU:H	1.50	0.76
1:E:710:HIS:C	1:E:712:PHE:H	1.92	0.76
1:F:112:VAL:HG12	1:F:113:GLU:H	1.50	0.76
1:F:184:LYS:HE3	1:F:191:GLU:HB2	1.67	0.76
1:B:141:PHE:H	1:B:141:PHE:HD1	1.34	0.76
1:F:257:LEU:O	1:F:257:LEU:HG	1.84	0.76
1:B:710:HIS:C	1:B:712:PHE:H	1.92	0.76
1:D:372:LYS:HG3	1:D:373:LYS:N	2.00	0.76
1:F:581:GLN:HE21	1:F:629:ASN:H	1.33	0.76
1:A:112:VAL:HG12	1:A:113:GLU:H	1.51	0.76
1:A:746:LYS:O	1:A:750:GLN:HG2	1.85	0.76
1:C:372:LYS:HG3	1:C:373:LYS:N	2.00	0.76
1:C:597:ASN:HD21	1:C:601:GLU:H	1.33	0.76
1:E:123:GLU:HG2	1:E:124:GLU:N	1.97	0.76
1:E:478:ALA:HB1	1:E:486:LYS:O	1.86	0.76
1:B:257:LEU:HG	1:B:257:LEU:O	1.84	0.76
1:C:716:LYS:O	1:C:720:ILE:HG22	1.86	0.76
1:D:478:ALA:HB1	1:D:486:LYS:O	1.85	0.76
1:D:499:PRO:HD3	1:D:552:TRP:CH2	2.21	0.76
1:E:77:ASP:O	1:E:81:GLN:HB2	1.85	0.76
2:P:36:MET:CE	2:P:43:PRO:HG3	2.16	0.76
1:D:472:ARG:HB3	1:D:472:ARG:NH1	2.00	0.76
1:E:90:PRO:O	1:E:93:VAL:HG12	1.86	0.76
1:F:90:PRO:O	1:F:93:VAL:HG12	1.86	0.76
1:A:257:LEU:O	1:A:257:LEU:HG	1.83	0.76
1:D:90:PRO:O	1:D:93:VAL:HG12	1.86	0.76
1:D:184:LYS:HE3	1:D:191:GLU:HB2	1.67	0.76
1:E:327:LEU:HG	1:E:595:ILE:HG23	1.68	0.76
1:F:327:LEU:HG	1:F:595:ILE:HG23	1.68	0.76
1:F:472:ARG:HB3	1:F:472:ARG:NH1	2.00	0.76
1:F:722:ILE:HG23	1:F:760:VAL:CG1	2.16	0.76
1:B:345:THR:HG22	1:B:490:ALA:O	1.86	0.76
1:C:180:ASP:CG	1:C:181:ILE:H	1.94	0.76
1:E:345:THR:HG22	1:E:490:ALA:O	1.86	0.76
1:D:141:PHE:H	1:D:141:PHE:HD1	1.32	0.75
1:E:180:ASP:CG	1:E:181:ILE:H	1.94	0.75
1:E:372:LYS:HG3	1:E:373:LYS:N	2.00	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:716:LYS:O	1:F:720:ILE:HG22	1.86	0.75
1:A:716:LYS:O	1:A:720:ILE:HG22	1.86	0.75
1:C:257:LEU:HG	1:C:257:LEU:O	1.84	0.75
1:C:472:ARG:HB3	1:C:472:ARG:NH1	2.00	0.75
1:D:338:LEU:HD21	1:D:409:ARG:CZ	2.17	0.75
1:E:746:LYS:O	1:E:750:GLN:HG2	1.86	0.75
1:A:180:ASP:CG	1:A:181:ILE:H	1.94	0.75
1:A:581:GLN:HE21	1:A:629:ASN:H	1.31	0.75
1:B:77:ASP:O	1:B:81:GLN:HB2	1.86	0.75
1:C:387:ASN:HD22	1:C:387:ASN:N	1.81	0.75
1:E:184:LYS:HE3	1:E:191:GLU:HB2	1.67	0.75
1:F:338:LEU:HD21	1:F:409:ARG:CZ	2.16	0.75
1:F:746:LYS:O	1:F:750:GLN:HG2	1.86	0.75
1:D:180:ASP:CG	1:D:181:ILE:H	1.94	0.75
1:E:236:GLU:HA	1:E:239:HIS:CD2	2.22	0.75
1:E:408:LEU:H	1:E:408:LEU:CD1	1.94	0.75
1:E:472:ARG:HB3	1:E:472:ARG:NH1	2.00	0.75
1:F:141:PHE:HD1	1:F:141:PHE:H	1.33	0.75
1:B:184:LYS:HE3	1:B:191:GLU:HB2	1.67	0.75
1:C:499:PRO:HD3	1:C:552:TRP:CH2	2.21	0.75
1:B:478:ALA:HB1	1:B:486:LYS:O	1.86	0.75
1:B:722:ILE:HG23	1:B:760:VAL:CG1	2.16	0.75
1:E:99:GLU:C	1:E:101:GLY:H	1.95	0.75
1:A:184:LYS:HE3	1:A:191:GLU:HB2	1.67	0.75
1:A:345:THR:HG22	1:A:490:ALA:O	1.86	0.75
1:C:338:LEU:HD21	1:C:409:ARG:CZ	2.17	0.75
1:A:90:PRO:O	1:A:93:VAL:HG12	1.87	0.75
1:E:142:VAL:HG13	1:E:154:ILE:HD12	1.69	0.75
1:B:408:LEU:H	1:B:408:LEU:CD1	1.95	0.75
1:E:722:ILE:HG23	1:E:760:VAL:CG1	2.16	0.75
2:O:36:MET:CE	2:O:43:PRO:HG3	2.17	0.75
2:S:22:ASP:O	2:S:24:ASP:N	2.20	0.75
1:A:499:PRO:HD3	1:A:552:TRP:CH2	2.22	0.74
1:F:180:ASP:CG	1:F:181:ILE:H	1.94	0.74
2:R:22:ASP:O	2:R:24:ASP:N	2.20	0.74
1:B:90:PRO:O	1:B:93:VAL:HG12	1.86	0.74
1:B:123:GLU:CG	1:B:124:GLU:H	1.94	0.74
1:C:236:GLU:HA	1:C:239:HIS:CD2	2.22	0.74
1:D:324:THR:CB	1:D:499:PRO:HA	2.16	0.74
1:F:478:ALA:HB1	1:F:486:LYS:O	1.87	0.74
1:F:499:PRO:HD3	1:F:552:TRP:CH2	2.21	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:T:36:MET:CE	2:T:43:PRO:HG3	2.17	0.74
1:A:77:ASP:O	1:A:81:GLN:HB2	1.86	0.74
1:B:180:ASP:CG	1:B:181:ILE:H	1.94	0.74
1:B:236:GLU:HA	1:B:239:HIS:CD2	2.22	0.74
1:B:716:LYS:O	1:B:720:ILE:HG22	1.87	0.74
1:C:90:PRO:O	1:C:93:VAL:HG12	1.86	0.74
1:E:694:VAL:CG2	2:S:18:LEU:HD21	2.16	0.74
1:B:372:LYS:HG3	1:B:373:LYS:N	2.00	0.74
1:D:345:THR:HG22	1:D:490:ALA:O	1.87	0.74
1:D:722:ILE:HG23	1:D:760:VAL:CG1	2.18	0.74
1:E:338:LEU:HD21	1:E:409:ARG:CZ	2.16	0.74
1:E:716:LYS:O	1:E:720:ILE:HG22	1.86	0.74
1:A:372:LYS:HG3	1:A:373:LYS:N	2.00	0.74
1:F:345:THR:HG22	1:F:490:ALA:O	1.87	0.74
2:Q:92:PHE:O	2:Q:94:LYS:N	2.19	0.74
1:B:338:LEU:HD21	1:B:409:ARG:CZ	2.17	0.74
1:C:722:ILE:HG23	1:C:760:VAL:CG1	2.18	0.74
2:Q:36:MET:CE	2:Q:43:PRO:HG3	2.17	0.74
1:D:142:VAL:HG13	1:D:154:ILE:HD12	1.70	0.74
1:D:236:GLU:HA	1:D:239:HIS:CD2	2.22	0.74
2:P:22:ASP:O	2:P:24:ASP:N	2.20	0.74
1:A:327:LEU:HG	1:A:595:ILE:HG23	1.70	0.74
1:B:499:PRO:HD3	1:B:552:TRP:CH2	2.23	0.74
1:D:581:GLN:HE21	1:D:629:ASN:H	1.33	0.74
1:E:315:PHE:HA	1:E:318:ILE:HD13	1.70	0.74
1:F:77:ASP:O	1:F:81:GLN:HB2	1.88	0.74
1:A:236:GLU:HA	1:A:239:HIS:CD2	2.22	0.73
1:B:99:GLU:C	1:B:101:GLY:H	1.96	0.73
1:B:142:VAL:HG22	1:B:154:ILE:HG23	1.70	0.73
1:E:499:PRO:HD3	1:E:552:TRP:CH2	2.23	0.73
1:F:99:GLU:C	1:F:101:GLY:H	1.95	0.73
1:C:767:GLN:HG2	1:C:768:LYS:N	2.02	0.73
1:D:767:GLN:HG2	1:D:768:LYS:N	2.02	0.73
1:C:77:ASP:O	1:C:81:GLN:HB2	1.87	0.73
1:C:142:VAL:HG22	1:C:154:ILE:HG23	1.70	0.73
1:C:315:PHE:HA	1:C:318:ILE:HD13	1.70	0.73
1:D:218:LEU:HD11	1:D:225:ILE:HD11	1.69	0.73
1:F:315:PHE:HA	1:F:318:ILE:HD13	1.71	0.73
2:S:94:LYS:NZ	2:S:94:LYS:HB3	2.03	0.73
1:C:141:PHE:HD1	1:C:141:PHE:H	1.33	0.73
1:C:581:GLN:HE21	1:C:629:ASN:H	1.33	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:327:LEU:HG	1:D:595:ILE:HG23	1.69	0.73
1:E:668:SER:HA	2:S:14:GLU:HG3	1.71	0.73
1:F:236:GLU:HA	1:F:239:HIS:CD2	2.23	0.73
1:F:780:LEU:HD12	1:F:789:ASN:OD1	1.89	0.73
2:S:12:PHE:CE1	2:S:72:MET:HG3	2.24	0.73
1:A:338:LEU:HD21	1:A:409:ARG:CZ	2.19	0.73
1:C:780:LEU:HD12	1:C:789:ASN:OD1	1.88	0.73
1:E:324:THR:CB	1:E:499:PRO:HA	2.14	0.73
2:P:63:ILE:HG13	2:P:67:GLU:CB	2.18	0.73
2:R:63:ILE:HG13	2:R:67:GLU:CB	2.18	0.73
2:T:12:PHE:CE1	2:T:72:MET:HG3	2.24	0.73
1:A:142:VAL:HG13	1:A:154:ILE:CD1	2.19	0.73
1:A:315:PHE:HA	1:A:318:ILE:HD13	1.71	0.73
1:C:345:THR:HG22	1:C:490:ALA:O	1.87	0.73
2:O:63:ILE:HG13	2:O:67:GLU:CB	2.18	0.73
2:R:36:MET:CE	2:R:43:PRO:HG3	2.17	0.73
1:A:123:GLU:CG	1:A:124:GLU:H	1.94	0.73
1:A:767:GLN:HG2	1:A:768:LYS:N	2.02	0.73
1:A:780:LEU:HD12	1:A:789:ASN:OD1	1.88	0.73
1:B:315:PHE:HA	1:B:318:ILE:HD13	1.71	0.73
1:B:327:LEU:HG	1:B:595:ILE:HG23	1.71	0.73
1:F:142:VAL:HG22	1:F:154:ILE:HG23	1.70	0.73
1:F:337:ASN:ND2	1:F:412:GLU:OE2	2.21	0.73
2:Q:22:ASP:O	2:Q:24:ASP:N	2.20	0.73
2:T:6:GLU:HG3	2:T:7:GLU:N	2.04	0.73
2:T:22:ASP:O	2:T:24:ASP:N	2.19	0.73
1:B:337:ASN:ND2	1:B:412:GLU:OE2	2.22	0.73
1:C:142:VAL:HG13	1:C:154:ILE:HD12	1.71	0.73
1:D:99:GLU:C	1:D:101:GLY:H	1.95	0.73
1:D:315:PHE:HA	1:D:318:ILE:HD13	1.71	0.73
1:E:142:VAL:HG22	1:E:154:ILE:HG23	1.70	0.73
1:E:780:LEU:HD12	1:E:789:ASN:OD1	1.88	0.73
1:A:142:VAL:HG13	1:A:154:ILE:HD12	1.69	0.73
1:A:148:GLU:HG3	1:A:149:THR:N	2.04	0.73
1:C:324:THR:CB	1:C:499:PRO:HA	2.16	0.73
1:C:409:ARG:CZ	1:C:413:LEU:HD21	2.19	0.73
1:E:142:VAL:HG13	1:E:154:ILE:CD1	2.18	0.73
2:Q:63:ILE:HG13	2:Q:67:GLU:CB	2.18	0.73
2:R:109:MET:HG3	2:R:116:LEU:HD11	1.71	0.73
2:S:63:ILE:HG13	2:S:67:GLU:CB	2.18	0.73
2:T:63:ILE:HG13	2:T:67:GLU:CB	2.18	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:99:GLU:C	1:A:101:GLY:H	1.96	0.72
1:A:337:ASN:ND2	1:A:412:GLU:OE2	2.22	0.72
1:C:99:GLU:C	1:C:101:GLY:H	1.95	0.72
1:A:409:ARG:CZ	1:A:413:LEU:HD21	2.19	0.72
1:B:142:VAL:HG13	1:B:154:ILE:HD12	1.71	0.72
1:F:324:THR:CB	1:F:499:PRO:HA	2.15	0.72
1:F:767:GLN:HG2	1:F:768:LYS:N	2.02	0.72
2:O:12:PHE:CE1	2:O:72:MET:HG3	2.23	0.72
1:B:668:SER:HA	2:P:14:GLU:HG3	1.71	0.72
1:B:767:GLN:HG2	1:B:768:LYS:N	2.02	0.72
1:C:327:LEU:HG	1:C:595:ILE:HG23	1.70	0.72
1:E:337:ASN:ND2	1:E:412:GLU:OE2	2.21	0.72
1:E:767:GLN:HG2	1:E:768:LYS:N	2.02	0.72
1:C:337:ASN:ND2	1:C:412:GLU:OE2	2.21	0.72
1:D:275:GLY:CA	1:D:278:LYS:HE2	2.20	0.72
1:E:409:ARG:CZ	1:E:413:LEU:HD21	2.19	0.72
2:P:9:ILE:HD12	2:P:69:LEU:HD11	1.72	0.72
1:B:148:GLU:HG3	1:B:149:THR:N	2.04	0.72
1:B:409:ARG:CZ	1:B:413:LEU:HD21	2.19	0.72
1:C:360:VAL:O	1:C:360:VAL:CG2	2.38	0.72
1:D:780:LEU:HD12	1:D:789:ASN:OD1	1.89	0.72
1:E:275:GLY:CA	1:E:278:LYS:HE2	2.19	0.72
1:E:372:LYS:HG3	1:E:373:LYS:H	1.54	0.72
1:D:142:VAL:HG13	1:D:154:ILE:CD1	2.20	0.72
2:Q:12:PHE:CE1	2:Q:72:MET:HG3	2.25	0.72
1:B:372:LYS:HG3	1:B:373:LYS:H	1.54	0.72
1:C:148:GLU:HG3	1:C:149:THR:N	2.04	0.72
1:A:142:VAL:HG22	1:A:154:ILE:HG23	1.72	0.72
1:A:668:SER:HA	2:O:14:GLU:HG3	1.72	0.72
1:B:275:GLY:CA	1:B:278:LYS:HE2	2.20	0.72
1:B:780:LEU:HD12	1:B:789:ASN:OD1	1.88	0.72
1:D:123:GLU:CG	1:D:124:GLU:H	1.94	0.72
1:D:142:VAL:HG22	1:D:154:ILE:HG23	1.72	0.72
1:D:337:ASN:ND2	1:D:412:GLU:OE2	2.21	0.72
1:F:360:VAL:O	1:F:360:VAL:CG2	2.37	0.72
2:R:12:PHE:CE1	2:R:72:MET:HG3	2.24	0.72
2:S:6:GLU:HG3	2:S:7:GLU:N	2.05	0.72
1:B:142:VAL:HG13	1:B:154:ILE:CD1	2.20	0.72
2:O:22:ASP:O	2:O:24:ASP:N	2.21	0.72
2:P:12:PHE:CE1	2:P:72:MET:HG3	2.24	0.72
2:R:58:ASP:C	2:R:60:ASN:H	1.98	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:T:58:ASP:C	2:T:60:ASN:H	1.98	0.72
1:A:715:GLU:HA	1:A:718:ARG:CZ	2.20	0.72
1:F:148:GLU:HG3	1:F:149:THR:N	2.04	0.72
1:A:227:ILE:O	1:A:228:ASN:C	2.33	0.71
1:C:218:LEU:HD11	1:C:225:ILE:HD11	1.72	0.71
1:F:227:ILE:O	1:F:228:ASN:C	2.33	0.71
1:F:306:GLY:O	1:F:336:THR:HG23	1.90	0.71
1:F:372:LYS:HG3	1:F:373:LYS:H	1.55	0.71
2:O:9:ILE:HD12	2:O:69:LEU:HD11	1.72	0.71
2:S:36:MET:CE	2:S:43:PRO:HG3	2.18	0.71
1:D:334:LEU:H	1:D:334:LEU:HD12	1.55	0.71
1:F:409:ARG:CZ	1:F:413:LEU:HD21	2.20	0.71
1:F:789:ASN:O	1:F:792:VAL:HB	1.90	0.71
2:Q:58:ASP:C	2:Q:60:ASN:H	1.99	0.71
1:A:349:ASN:HD22	1:A:350:VAL:N	1.87	0.71
1:F:142:VAL:HG13	1:F:154:ILE:HD12	1.70	0.71
1:F:360:VAL:CG1	1:F:370:LEU:HD22	2.20	0.71
1:D:409:ARG:CZ	1:D:413:LEU:HD21	2.19	0.71
1:F:275:GLY:CA	1:F:278:LYS:HE2	2.20	0.71
2:O:6:GLU:HG3	2:O:7:GLU:N	2.05	0.71
2:O:58:ASP:C	2:O:60:ASN:H	1.98	0.71
2:S:9:ILE:HD12	2:S:69:LEU:HD11	1.72	0.71
1:A:408:LEU:H	1:A:408:LEU:CD1	1.95	0.71
1:F:611:THR:HG22	1:F:615:ILE:HD11	1.73	0.71
2:Q:109:MET:HG3	2:Q:116:LEU:HD11	1.72	0.71
1:B:377:GLN:O	1:B:381:GLU:HB2	1.91	0.71
1:C:234:LEU:HD23	1:C:235:THR:H	1.56	0.71
2:P:6:GLU:HG3	2:P:7:GLU:N	2.05	0.71
1:B:324:THR:CB	1:B:499:PRO:HA	2.16	0.71
1:B:789:ASN:O	1:B:792:VAL:HB	1.91	0.71
1:C:79:ILE:O	1:C:81:GLN:N	2.24	0.71
1:D:227:ILE:O	1:D:228:ASN:C	2.33	0.71
1:E:148:GLU:HG3	1:E:149:THR:N	2.04	0.71
1:F:218:LEU:HD11	1:F:225:ILE:HD11	1.72	0.71
1:F:270:LYS:HD3	1:F:273:LYS:HD2	1.73	0.71
2:T:109:MET:HG3	2:T:116:LEU:HD11	1.73	0.71
1:A:197:LYS:NZ	1:A:197:LYS:HB3	2.06	0.71
1:C:275:GLY:CA	1:C:278:LYS:HE2	2.20	0.71
1:F:234:LEU:HD23	1:F:235:THR:H	1.56	0.71
2:R:6:GLU:HG3	2:R:7:GLU:N	2.04	0.71
1:C:142:VAL:HG13	1:C:154:ILE:CD1	2.21	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:789:ASN:O	1:D:792:VAL:HB	1.91	0.71
2:S:109:MET:HG3	2:S:116:LEU:HD11	1.73	0.71
2:T:9:ILE:HD12	2:T:69:LEU:HD11	1.72	0.71
1:A:115:LYS:CE	1:A:116:GLU:HG2	2.21	0.71
1:A:275:GLY:CA	1:A:278:LYS:HE2	2.20	0.71
1:B:197:LYS:HZ2	1:B:197:LYS:HB3	1.55	0.71
1:D:148:GLU:HG3	1:D:149:THR:N	2.04	0.71
1:E:739:LYS:HG2	1:E:740:GLN:H	1.56	0.71
2:S:58:ASP:C	2:S:60:ASN:H	1.98	0.71
1:A:89:ILE:HD13	1:A:175:LYS:HE2	1.73	0.70
1:D:372:LYS:HG3	1:D:373:LYS:H	1.55	0.70
1:D:668:SER:HA	2:R:14:GLU:HG3	1.73	0.70
1:D:739:LYS:HG2	1:D:740:GLN:H	1.56	0.70
1:F:142:VAL:HG13	1:F:154:ILE:CD1	2.20	0.70
1:B:218:LEU:HD11	1:B:225:ILE:HD11	1.73	0.70
1:B:227:ILE:O	1:B:228:ASN:C	2.33	0.70
1:D:115:LYS:CE	1:D:116:GLU:HG2	2.21	0.70
1:E:218:LEU:HD11	1:E:225:ILE:HD11	1.73	0.70
1:E:592:GLU:HB3	1:E:604:LEU:HD11	1.72	0.70
1:A:739:LYS:HG2	1:A:740:GLN:H	1.56	0.70
1:E:88:LYS:HZ3	1:E:172:GLU:CD	1.97	0.70
1:E:227:ILE:O	1:E:228:ASN:C	2.33	0.70
2:P:58:ASP:C	2:P:60:ASN:H	1.98	0.70
1:A:592:GLU:HB3	1:A:604:LEU:HD11	1.74	0.70
1:C:270:LYS:HD3	1:C:273:LYS:HD2	1.73	0.70
1:D:349:ASN:HD22	1:D:350:VAL:N	1.88	0.70
1:D:715:GLU:HA	1:D:718:ARG:CZ	2.20	0.70
1:E:270:LYS:HD3	1:E:273:LYS:HD2	1.73	0.70
1:F:739:LYS:HG2	1:F:740:GLN:H	1.56	0.70
2:Q:9:ILE:HD12	2:Q:69:LEU:HD11	1.71	0.70
1:B:349:ASN:HD22	1:B:350:VAL:N	1.88	0.70
1:B:715:GLU:HA	1:B:718:ARG:CZ	2.21	0.70
1:B:739:LYS:HG2	1:B:740:GLN:H	1.56	0.70
1:C:739:LYS:HG2	1:C:740:GLN:H	1.56	0.70
1:E:377:GLN:O	1:E:381:GLU:HB2	1.92	0.70
2:Q:6:GLU:HG3	2:Q:7:GLU:N	2.06	0.70
2:T:13:LYS:C	2:T:15:ALA:H	1.97	0.70
1:A:620:THR:HG22	1:A:621:GLY:N	2.07	0.70
1:A:789:ASN:O	1:A:792:VAL:HB	1.91	0.70
1:B:389:LYS:HA	1:B:392:THR:HB	1.73	0.70
1:C:789:ASN:O	1:C:792:VAL:HB	1.91	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:789:ASN:O	1:E:792:VAL:HB	1.91	0.70
1:F:90:PRO:HG2	1:F:93:VAL:HB	1.74	0.70
1:F:715:GLU:HA	1:F:718:ARG:CZ	2.21	0.70
2:O:13:LYS:C	2:O:15:ALA:H	2.00	0.70
1:A:218:LEU:HD11	1:A:225:ILE:HD11	1.72	0.70
1:A:372:LYS:HG3	1:A:373:LYS:H	1.56	0.70
1:C:668:SER:HA	2:Q:14:GLU:HG3	1.73	0.70
1:E:90:PRO:HG2	1:E:93:VAL:HB	1.73	0.70
1:E:165:GLN:NE2	1:E:252:ASP:HB3	2.07	0.70
1:F:334:LEU:HD12	1:F:334:LEU:H	1.56	0.70
1:F:412:GLU:HA	1:F:415:GLU:OE2	1.92	0.70
2:R:9:ILE:HD12	2:R:69:LEU:HD11	1.72	0.70
1:A:412:GLU:HA	1:A:415:GLU:OE2	1.92	0.70
1:B:234:LEU:HD23	1:B:235:THR:H	1.56	0.70
1:C:115:LYS:CE	1:C:116:GLU:HG2	2.21	0.70
1:C:349:ASN:HD22	1:C:350:VAL:N	1.88	0.70
1:D:89:ILE:HD13	1:D:175:LYS:HE2	1.74	0.70
1:E:197:LYS:HB3	1:E:197:LYS:NZ	2.07	0.70
1:E:464:VAL:HG23	1:E:465:LEU:HD12	1.74	0.70
1:E:715:GLU:HA	1:E:718:ARG:CZ	2.20	0.70
1:F:668:SER:HA	2:T:14:GLU:HG3	1.73	0.70
1:B:301:ALA:O	1:B:303:LYS:N	2.25	0.70
1:C:377:GLN:O	1:C:381:GLU:HB2	1.92	0.70
1:D:189:ASP:O	1:D:191:GLU:N	2.23	0.70
1:D:270:LYS:HD3	1:D:273:LYS:HD2	1.73	0.70
1:D:389:LYS:HA	1:D:392:THR:HB	1.73	0.70
1:D:792:VAL:O	1:D:796:ILE:HG12	1.92	0.70
1:F:349:ASN:HD22	1:F:350:VAL:N	1.88	0.70
2:P:109:MET:HG3	2:P:116:LEU:HD11	1.74	0.70
2:Q:31:GLU:O	2:Q:35:VAL:HG23	1.92	0.70
1:D:270:LYS:HA	1:D:273:LYS:HD2	1.74	0.70
1:E:412:GLU:HA	1:E:415:GLU:OE2	1.92	0.70
2:T:72:MET:O	2:T:76:MET:CB	2.40	0.70
1:A:234:LEU:HD23	1:A:235:THR:H	1.56	0.69
1:A:464:VAL:HG23	1:A:465:LEU:HD12	1.74	0.69
1:B:697:ILE:HD13	1:B:732:ILE:HD13	1.74	0.69
1:C:227:ILE:O	1:C:228:ASN:C	2.33	0.69
1:C:270:LYS:HA	1:C:273:LYS:HD2	1.74	0.69
1:E:357:TRP:CZ3	1:E:439:ASN:HB2	2.27	0.69
1:F:377:GLN:O	1:F:381:GLU:HB2	1.91	0.69
1:A:270:LYS:HA	1:A:273:LYS:HD2	1.74	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:306:GLY:O	1:A:336:THR:HG23	1.92	0.69
1:B:464:VAL:HG23	1:B:465:LEU:HD12	1.74	0.69
1:C:334:LEU:HD12	1:C:334:LEU:H	1.57	0.69
1:C:611:THR:HG22	1:C:615:ILE:HD11	1.74	0.69
1:C:697:ILE:HD13	1:C:732:ILE:HD13	1.75	0.69
1:D:79:ILE:O	1:D:81:GLN:N	2.25	0.69
1:D:301:ALA:O	1:D:303:LYS:N	2.25	0.69
1:E:389:LYS:HA	1:E:392:THR:HB	1.74	0.69
1:F:592:GLU:HB3	1:F:604:LEU:HD11	1.73	0.69
2:S:13:LYS:C	2:S:15:ALA:H	2.00	0.69
1:A:165:GLN:CD	1:A:252:ASP:HB3	2.17	0.69
1:C:715:GLU:HA	1:C:718:ARG:CZ	2.22	0.69
1:D:234:LEU:HD23	1:D:235:THR:H	1.56	0.69
1:D:377:GLN:O	1:D:381:GLU:HB2	1.92	0.69
1:E:334:LEU:H	1:E:334:LEU:HD12	1.56	0.69
1:F:567:THR:HG23	1:F:568:GLY:N	2.07	0.69
1:B:357:TRP:CZ3	1:B:439:ASN:HB2	2.28	0.69
1:D:412:GLU:HA	1:D:415:GLU:OE2	1.92	0.69
1:D:611:THR:HG22	1:D:615:ILE:HD11	1.74	0.69
1:F:301:ALA:O	1:F:303:LYS:N	2.25	0.69
1:A:165:GLN:NE2	1:A:252:ASP:HB3	2.07	0.69
1:A:357:TRP:CZ3	1:A:439:ASN:HB2	2.28	0.69
1:C:81:GLN:OE1	1:C:156:ILE:HG21	1.92	0.69
1:C:301:ALA:O	1:C:303:LYS:N	2.25	0.69
1:D:592:GLU:HB3	1:D:604:LEU:HD11	1.75	0.69
1:F:197:LYS:NZ	1:F:197:LYS:HB3	2.07	0.69
1:B:165:GLN:NE2	1:B:252:ASP:HB3	2.08	0.69
1:B:270:LYS:HA	1:B:273:LYS:HD2	1.74	0.69
1:B:620:THR:HG22	1:B:621:GLY:N	2.08	0.69
1:C:620:THR:HG22	1:C:621:GLY:N	2.08	0.69
1:E:234:LEU:HD23	1:E:235:THR:H	1.56	0.69
1:F:165:GLN:CD	1:F:252:ASP:HB3	2.18	0.69
1:F:357:TRP:CZ3	1:F:439:ASN:HB2	2.27	0.69
1:F:464:VAL:HG23	1:F:465:LEU:HD12	1.74	0.69
2:O:109:MET:HG3	2:O:116:LEU:HD11	1.74	0.69
2:R:13:LYS:C	2:R:15:ALA:H	2.00	0.69
1:A:188:LEU:CD2	1:A:188:LEU:N	2.47	0.69
1:A:189:ASP:O	1:A:191:GLU:N	2.24	0.69
1:A:611:THR:HG22	1:A:615:ILE:HD11	1.75	0.69
1:A:792:VAL:O	1:A:796:ILE:HG12	1.93	0.69
1:C:389:LYS:HA	1:C:392:THR:HB	1.73	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:165:GLN:CD	1:D:252:ASP:HB3	2.17	0.69
1:D:357:TRP:CZ3	1:D:439:ASN:HB2	2.28	0.69
1:E:270:LYS:HA	1:E:273:LYS:HD2	1.73	0.69
1:F:389:LYS:HA	1:F:392:THR:HB	1.73	0.69
1:A:301:ALA:O	1:A:303:LYS:N	2.26	0.69
1:A:377:GLN:O	1:A:381:GLU:HB2	1.92	0.69
1:C:90:PRO:HG2	1:C:93:VAL:HB	1.75	0.69
1:C:792:VAL:O	1:C:796:ILE:HG12	1.92	0.69
1:D:306:GLY:O	1:D:336:THR:HG23	1.92	0.69
1:D:481:VAL:O	1:D:484:VAL:HG23	1.93	0.69
1:E:115:LYS:CE	1:E:116:GLU:HG2	2.21	0.69
1:E:120:LEU:C	1:E:120:LEU:HD13	2.18	0.69
1:E:349:ASN:HD22	1:E:350:VAL:N	1.89	0.69
1:E:792:VAL:O	1:E:796:ILE:HG12	1.92	0.69
2:R:31:GLU:O	2:R:35:VAL:HG23	1.92	0.69
2:T:31:GLU:O	2:T:35:VAL:HG23	1.93	0.69
1:A:79:ILE:O	1:A:81:GLN:N	2.25	0.69
1:A:324:THR:CB	1:A:499:PRO:HA	2.16	0.69
1:D:408:LEU:H	1:D:408:LEU:CD1	1.95	0.69
1:D:611:THR:O	1:D:615:ILE:HG13	1.93	0.69
1:E:301:ALA:O	1:E:303:LYS:N	2.25	0.69
1:E:620:THR:HG22	1:E:621:GLY:N	2.06	0.69
1:F:115:LYS:CE	1:F:116:GLU:HG2	2.22	0.69
1:A:90:PRO:HG2	1:A:93:VAL:HB	1.75	0.69
1:B:165:GLN:CD	1:B:252:ASP:HB3	2.18	0.69
1:B:306:GLY:O	1:B:336:THR:HG23	1.92	0.69
1:D:90:PRO:HG2	1:D:93:VAL:HB	1.75	0.69
1:D:620:THR:HG22	1:D:621:GLY:N	2.08	0.69
1:F:792:VAL:O	1:F:796:ILE:HG12	1.92	0.69
1:A:389:LYS:HA	1:A:392:THR:HB	1.74	0.68
1:B:115:LYS:C	1:B:117:LEU:H	2.01	0.68
1:B:293:ILE:HG22	1:B:613:ARG:HH21	1.57	0.68
1:B:412:GLU:HA	1:B:415:GLU:OE2	1.92	0.68
1:C:357:TRP:CZ3	1:C:439:ASN:HB2	2.28	0.68
1:D:464:VAL:HG23	1:D:465:LEU:HD12	1.74	0.68
1:F:270:LYS:HA	1:F:273:LYS:HD2	1.74	0.68
2:S:30:LYS:H	2:S:30:LYS:CD	2.03	0.68
1:A:270:LYS:HD3	1:A:273:LYS:HD2	1.74	0.68
1:A:457:THR:HG21	1:A:468:LYS:HA	1.75	0.68
1:B:120:LEU:HD13	1:B:120:LEU:C	2.18	0.68
1:B:197:LYS:HB3	1:B:197:LYS:NZ	2.07	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:165:GLN:CD	1:C:252:ASP:HB3	2.18	0.68
1:C:464:VAL:HG23	1:C:465:LEU:HD12	1.73	0.68
1:E:165:GLN:CD	1:E:252:ASP:HB3	2.17	0.68
2:S:106:ARG:O	2:S:110:THR:HG23	1.93	0.68
1:B:115:LYS:CE	1:B:116:GLU:HG2	2.22	0.68
1:C:306:GLY:O	1:C:336:THR:HG23	1.92	0.68
1:C:412:GLU:HA	1:C:415:GLU:OE2	1.93	0.68
1:E:567:THR:HG23	1:E:568:GLY:N	2.09	0.68
1:F:190:PRO:HB2	1:F:195:LEU:HD13	1.75	0.68
2:P:31:GLU:O	2:P:35:VAL:HG23	1.93	0.68
1:B:792:VAL:O	1:B:796:ILE:HG12	1.94	0.68
1:C:372:LYS:HG3	1:C:373:LYS:H	1.56	0.68
1:F:165:GLN:NE2	1:F:252:ASP:HB3	2.08	0.68
1:B:481:VAL:O	1:B:484:VAL:HG23	1.93	0.68
1:B:592:GLU:HB3	1:B:604:LEU:HD11	1.74	0.68
1:B:611:THR:HG22	1:B:615:ILE:HD11	1.74	0.68
1:D:472:ARG:HH11	1:D:472:ARG:CB	2.07	0.68
1:F:120:LEU:C	1:F:120:LEU:HD13	2.18	0.68
1:C:165:GLN:NE2	1:C:252:ASP:HB3	2.07	0.68
1:D:81:GLN:OE1	1:D:156:ILE:HG21	1.93	0.68
1:E:457:THR:HG21	1:E:468:LYS:HA	1.76	0.68
1:F:79:ILE:O	1:F:81:GLN:N	2.27	0.68
1:F:697:ILE:HD13	1:F:732:ILE:HD13	1.76	0.68
2:O:30:LYS:H	2:O:30:LYS:CD	2.03	0.68
2:R:106:ARG:O	2:R:110:THR:HG23	1.94	0.68
1:A:120:LEU:HD13	1:A:120:LEU:C	2.18	0.68
1:B:457:THR:HG21	1:B:468:LYS:HA	1.76	0.68
1:E:190:PRO:HB2	1:E:195:LEU:HD13	1.76	0.68
2:P:106:ARG:O	2:P:110:THR:HG23	1.93	0.68
1:C:115:LYS:C	1:C:117:LEU:H	2.02	0.68
1:C:190:PRO:HB2	1:C:195:LEU:HD13	1.75	0.68
2:O:31:GLU:O	2:O:35:VAL:HG23	1.94	0.68
2:S:12:PHE:HD1	2:S:72:MET:HG3	1.59	0.68
2:S:100:ILE:HB	2:S:136:VAL:CG2	2.21	0.68
1:A:334:LEU:H	1:A:334:LEU:HD12	1.58	0.68
1:B:472:ARG:HH11	1:B:472:ARG:CB	2.07	0.68
1:D:293:ILE:HG22	1:D:613:ARG:HH21	1.59	0.68
1:E:306:GLY:O	1:E:336:THR:HG23	1.93	0.68
1:F:89:ILE:HD13	1:F:175:LYS:HE2	1.76	0.68
2:Q:13:LYS:C	2:Q:15:ALA:H	2.00	0.68
2:Q:30:LYS:H	2:Q:30:LYS:CD	2.03	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:R:30:LYS:H	2:R:30:LYS:CD	2.03	0.68
1:C:89:ILE:HD13	1:C:175:LYS:HE2	1.76	0.68
1:D:197:LYS:NZ	1:D:197:LYS:HB3	2.07	0.68
1:D:457:THR:HG21	1:D:468:LYS:HA	1.75	0.68
1:E:115:LYS:C	1:E:117:LEU:H	2.01	0.68
1:E:611:THR:HG22	1:E:615:ILE:HD11	1.75	0.68
1:E:697:ILE:HD13	1:E:732:ILE:HD13	1.76	0.68
2:P:30:LYS:H	2:P:30:LYS:CD	2.04	0.68
2:R:94:LYS:NZ	2:R:94:LYS:HB3	2.08	0.68
1:B:190:PRO:HB2	1:B:195:LEU:HD13	1.75	0.67
1:B:270:LYS:HD3	1:B:273:LYS:HD2	1.74	0.67
1:B:567:THR:HG23	1:B:568:GLY:N	2.07	0.67
1:C:611:THR:O	1:C:615:ILE:HG13	1.93	0.67
1:C:697:ILE:CD1	1:C:732:ILE:HD13	2.24	0.67
1:D:120:LEU:HD13	1:D:120:LEU:C	2.18	0.67
1:F:81:GLN:OE1	1:F:156:ILE:HG21	1.93	0.67
1:B:79:ILE:O	1:B:81:GLN:N	2.26	0.67
1:C:478:ALA:HA	1:C:488:LEU:HG	1.74	0.67
1:C:592:GLU:HB3	1:C:604:LEU:HD11	1.75	0.67
1:D:697:ILE:HD13	1:D:732:ILE:HD13	1.76	0.67
1:B:478:ALA:HA	1:B:488:LEU:HG	1.76	0.67
1:C:197:LYS:NZ	1:C:197:LYS:HB3	2.08	0.67
1:D:165:GLN:NE2	1:D:252:ASP:HB3	2.08	0.67
1:E:254:ARG:H	1:E:254:ARG:HD2	1.59	0.67
1:F:472:ARG:HH11	1:F:472:ARG:CB	2.07	0.67
1:F:481:VAL:O	1:F:484:VAL:HG23	1.94	0.67
1:B:90:PRO:HG2	1:B:93:VAL:HB	1.76	0.67
1:B:159:TYR:HD1	1:B:159:TYR:H	1.42	0.67
1:B:697:ILE:CD1	1:B:732:ILE:HD13	2.24	0.67
1:C:120:LEU:C	1:C:120:LEU:HD13	2.18	0.67
1:F:357:TRP:HZ3	1:F:439:ASN:HB2	1.60	0.67
2:P:13:LYS:C	2:P:15:ALA:H	2.01	0.67
1:A:385:LEU:O	1:A:385:LEU:HD13	1.95	0.67
1:A:481:VAL:O	1:A:484:VAL:HG23	1.94	0.67
1:B:334:LEU:H	1:B:334:LEU:HD12	1.59	0.67
1:C:481:VAL:O	1:C:484:VAL:HG23	1.95	0.67
1:D:478:ALA:HA	1:D:488:LEU:HG	1.76	0.67
1:F:254:ARG:H	1:F:254:ARG:HD2	1.59	0.67
1:F:293:ILE:HG22	1:F:613:ARG:HH21	1.59	0.67
1:F:620:THR:HG22	1:F:621:GLY:N	2.09	0.67
2:T:12:PHE:HD1	2:T:72:MET:HG3	1.60	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:T:106:ARG:O	2:T:110:THR:HG23	1.93	0.67
1:A:719:LYS:HG2	1:A:797:ILE:CD1	2.24	0.67
2:S:31:GLU:O	2:S:35:VAL:HG23	1.94	0.67
1:A:159:TYR:H	1:A:159:TYR:HD1	1.42	0.67
1:A:293:ILE:HG22	1:A:613:ARG:HH21	1.60	0.67
1:A:478:ALA:HA	1:A:488:LEU:HG	1.75	0.67
1:B:189:ASP:O	1:B:191:GLU:N	2.24	0.67
1:B:301:ALA:C	1:B:303:LYS:N	2.53	0.67
1:C:335:ALA:O	1:C:339:ILE:HG13	1.95	0.67
1:D:435:LEU:HG	1:D:446:ILE:HG22	1.76	0.67
1:E:81:GLN:OE1	1:E:156:ILE:HG21	1.94	0.67
1:E:357:TRP:HZ3	1:E:439:ASN:HB2	1.60	0.67
1:F:611:THR:O	1:F:615:ILE:HG13	1.94	0.67
1:F:719:LYS:HG2	1:F:797:ILE:CD1	2.24	0.67
2:O:106:ARG:O	2:O:110:THR:HG23	1.94	0.67
1:A:719:LYS:HG2	1:A:797:ILE:HD11	1.77	0.67
1:C:435:LEU:HG	1:C:446:ILE:HG22	1.75	0.67
1:D:115:LYS:C	1:D:117:LEU:H	2.02	0.67
1:D:719:LYS:HG2	1:D:797:ILE:CD1	2.24	0.67
1:E:719:LYS:HG2	1:E:797:ILE:CD1	2.24	0.67
2:P:100:ILE:HB	2:P:136:VAL:CG2	2.23	0.67
2:Q:106:ARG:O	2:Q:110:THR:HG23	1.93	0.67
1:D:159:TYR:HD1	1:D:159:TYR:H	1.42	0.67
1:D:186:LYS:C	1:D:188:LEU:O	2.38	0.67
1:D:190:PRO:HB2	1:D:195:LEU:HD13	1.75	0.67
1:F:657:ILE:HG13	1:F:756:ILE:CD1	2.25	0.67
1:B:81:GLN:OE1	1:B:156:ILE:HG21	1.94	0.67
1:C:254:ARG:H	1:C:254:ARG:HD2	1.60	0.67
1:D:335:ALA:O	1:D:339:ILE:HG13	1.95	0.67
1:F:385:LEU:HD13	1:F:385:LEU:O	1.95	0.67
1:F:457:THR:HG21	1:F:468:LYS:HA	1.76	0.67
1:F:550:SER:HG	1:F:552:TRP:CD1	2.13	0.67
1:C:457:THR:HG21	1:C:468:LYS:HA	1.75	0.66
1:E:122:GLU:HB2	1:E:147:ARG:HG3	1.77	0.66
2:T:30:LYS:H	2:T:30:LYS:CD	2.04	0.66
1:A:81:GLN:OE1	1:A:156:ILE:HG21	1.94	0.66
1:A:567:THR:HG23	1:A:568:GLY:N	2.08	0.66
1:B:235:THR:O	1:B:238:GLN:HB2	1.96	0.66
1:B:611:THR:O	1:B:615:ILE:HG13	1.95	0.66
1:E:481:VAL:O	1:E:484:VAL:HG23	1.95	0.66
2:P:33:GLY:O	2:P:37:ARG:HG3	1.96	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:Q:100:ILE:HB	2:Q:136:VAL:CG2	2.21	0.66
1:A:190:PRO:HB2	1:A:195:LEU:HD13	1.75	0.66
1:C:235:THR:O	1:C:238:GLN:HB2	1.96	0.66
2:O:72:MET:O	2:O:76:MET:CB	2.40	0.66
2:Q:72:MET:O	2:Q:76:MET:CB	2.40	0.66
1:B:719:LYS:HG2	1:B:797:ILE:CD1	2.25	0.66
1:C:293:ILE:HG22	1:C:613:ARG:HH21	1.60	0.66
1:D:697:ILE:CD1	1:D:732:ILE:HD13	2.25	0.66
1:E:326:ILE:HG22	1:E:328:PHE:CE1	2.30	0.66
2:T:58:ASP:CB	2:T:62:THR:HG23	2.26	0.66
1:A:435:LEU:HG	1:A:446:ILE:HG22	1.76	0.66
1:A:678:VAL:HG13	1:A:745:TYR:CD2	2.31	0.66
1:B:254:ARG:H	1:B:254:ARG:HD2	1.60	0.66
1:C:189:ASP:O	1:C:191:GLU:N	2.26	0.66
1:C:688:PHE:C	1:C:688:PHE:CD2	2.74	0.66
1:D:719:LYS:HG2	1:D:797:ILE:HD11	1.77	0.66
1:E:293:ILE:HG22	1:E:613:ARG:HH21	1.60	0.66
1:F:478:ALA:HA	1:F:488:LEU:HG	1.76	0.66
1:B:122:GLU:HB2	1:B:147:ARG:HG3	1.77	0.66
1:C:567:THR:HG23	1:C:568:GLY:N	2.07	0.66
1:D:567:THR:HG23	1:D:568:GLY:N	2.08	0.66
1:D:688:PHE:C	1:D:688:PHE:CD2	2.73	0.66
1:F:326:ILE:HG22	1:F:328:PHE:CE1	2.31	0.66
1:A:335:ALA:O	1:A:339:ILE:HG13	1.95	0.66
1:A:697:ILE:HD13	1:A:732:ILE:HD13	1.77	0.66
1:B:210:PHE:C	1:B:212:GLN:H	2.04	0.66
1:D:385:LEU:O	1:D:385:LEU:HD13	1.96	0.66
1:D:550:SER:HG	1:D:552:TRP:CD1	2.14	0.66
1:E:210:PHE:C	1:E:212:GLN:H	2.04	0.66
1:E:238:GLN:C	1:E:240:ALA:H	2.04	0.66
1:E:435:LEU:HG	1:E:446:ILE:HG22	1.77	0.66
1:F:186:LYS:C	1:F:188:LEU:O	2.38	0.66
1:F:235:THR:O	1:F:238:GLN:HB2	1.96	0.66
2:T:100:ILE:HB	2:T:136:VAL:CG2	2.23	0.66
1:A:187:SER:O	1:A:188:LEU:C	2.37	0.66
1:E:235:THR:O	1:E:238:GLN:HB2	1.96	0.66
1:E:478:ALA:HA	1:E:488:LEU:HG	1.76	0.66
1:F:567:THR:CG2	1:F:568:GLY:N	2.58	0.66
1:A:115:LYS:C	1:A:117:LEU:H	2.02	0.66
1:A:657:ILE:HG13	1:A:756:ILE:CD1	2.25	0.66
1:A:697:ILE:CD1	1:A:732:ILE:HD13	2.26	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:188:LEU:CD2	1:B:188:LEU:N	2.47	0.66
1:B:357:TRP:HZ3	1:B:439:ASN:HB2	1.61	0.66
1:C:122:GLU:HB2	1:C:147:ARG:HG3	1.77	0.66
1:D:73:ASN:O	1:D:74:GLU:O	2.14	0.66
1:D:254:ARG:H	1:D:254:ARG:HD2	1.60	0.66
1:E:89:ILE:HD13	1:E:175:LYS:HE2	1.78	0.66
2:R:72:MET:O	2:R:76:MET:CB	2.40	0.66
2:S:13:LYS:HZ3	2:S:65:PHE:HB3	1.61	0.66
1:A:357:TRP:HZ3	1:A:439:ASN:HB2	1.61	0.66
1:A:472:ARG:HH11	1:A:472:ARG:CB	2.07	0.66
1:C:657:ILE:HG13	1:C:756:ILE:CD1	2.26	0.66
1:F:397:GLU:HG3	1:F:480:ASN:HB3	1.77	0.66
1:B:73:ASN:O	1:B:74:GLU:O	2.14	0.65
1:B:435:LEU:HG	1:B:446:ILE:HG22	1.77	0.65
1:D:718:ARG:NH1	1:D:767:GLN:HE21	1.94	0.65
2:R:126:ARG:HH21	2:R:126:ARG:HG3	1.61	0.65
1:A:88:LYS:HZ3	1:A:172:GLU:CD	2.02	0.65
1:B:326:ILE:HG22	1:B:328:PHE:CE1	2.32	0.65
1:B:385:LEU:O	1:B:385:LEU:HD13	1.97	0.65
1:C:246:SER:O	1:C:250:ALA:HB2	1.97	0.65
1:C:350:VAL:HG12	1:C:352:GLY:H	1.62	0.65
1:C:550:SER:HG	1:C:552:TRP:CD1	2.15	0.65
1:D:187:SER:O	1:D:188:LEU:C	2.37	0.65
1:F:115:LYS:C	1:F:117:LEU:H	2.02	0.65
1:A:301:ALA:C	1:A:303:LYS:N	2.53	0.65
1:C:159:TYR:HD1	1:C:159:TYR:H	1.42	0.65
1:C:567:THR:CG2	1:C:568:GLY:N	2.58	0.65
1:C:719:LYS:HG2	1:C:797:ILE:CD1	2.25	0.65
1:D:130:SER:HB2	1:D:170:TYR:CE2	2.31	0.65
1:D:301:ALA:C	1:D:303:LYS:N	2.52	0.65
1:E:79:ILE:O	1:E:81:GLN:N	2.26	0.65
2:P:12:PHE:HD1	2:P:72:MET:HG3	1.60	0.65
2:T:21:LYS:C	2:T:21:LYS:HD3	2.22	0.65
1:A:186:LYS:C	1:A:188:LEU:O	2.40	0.65
1:A:243:LEU:HA	1:A:246:SER:OG	1.97	0.65
1:E:335:ALA:O	1:E:339:ILE:HG13	1.96	0.65
1:F:73:ASN:O	1:F:74:GLU:O	2.14	0.65
1:F:435:LEU:HG	1:F:446:ILE:HG22	1.77	0.65
1:B:719:LYS:HG2	1:B:797:ILE:HD11	1.77	0.65
1:C:186:LYS:C	1:C:188:LEU:O	2.39	0.65
1:E:73:ASN:O	1:E:74:GLU:O	2.15	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:159:TYR:HD1	1:E:159:TYR:H	1.42	0.65
1:E:678:VAL:HG13	1:E:745:TYR:CD2	2.32	0.65
1:E:719:LYS:HG2	1:E:797:ILE:HD11	1.77	0.65
1:F:159:TYR:H	1:F:159:TYR:HD1	1.42	0.65
1:F:697:ILE:CD1	1:F:732:ILE:HD13	2.26	0.65
1:F:735:VAL:HA	1:F:738:SER:HB2	1.79	0.65
2:P:126:ARG:HH21	2:P:126:ARG:HG3	1.61	0.65
2:R:100:ILE:HB	2:R:136:VAL:CG2	2.21	0.65
2:S:58:ASP:CB	2:S:62:THR:HG23	2.26	0.65
1:B:186:LYS:C	1:B:188:LEU:O	2.40	0.65
1:B:688:PHE:C	1:B:688:PHE:CD2	2.74	0.65
1:F:122:GLU:HB2	1:F:147:ARG:HG3	1.77	0.65
1:F:305:SER:HB2	1:F:594:PHE:CD1	2.32	0.65
1:F:692:GLU:HA	1:F:692:GLU:OE2	1.95	0.65
1:B:711:ILE:HG13	1:B:712:PHE:CD2	2.32	0.65
1:C:199:LEU:C	1:C:201:ASP:H	2.05	0.65
1:C:301:ALA:C	1:C:303:LYS:N	2.52	0.65
1:C:735:VAL:HA	1:C:738:SER:HB2	1.79	0.65
1:D:122:GLU:HB2	1:D:147:ARG:HG3	1.77	0.65
1:D:199:LEU:C	1:D:201:ASP:H	2.05	0.65
1:F:243:LEU:HA	1:F:246:SER:OG	1.96	0.65
1:F:688:PHE:C	1:F:688:PHE:CD2	2.74	0.65
1:F:718:ARG:NH1	1:F:767:GLN:HE21	1.94	0.65
2:T:33:GLY:O	2:T:37:ARG:HG3	1.97	0.65
1:A:235:THR:O	1:A:238:GLN:HB2	1.97	0.65
1:A:305:SER:HB2	1:A:594:PHE:CD1	2.32	0.65
1:B:238:GLN:C	1:B:240:ALA:H	2.05	0.65
1:B:567:THR:CG2	1:B:568:GLY:N	2.59	0.65
1:B:692:GLU:HA	1:B:692:GLU:OE2	1.96	0.65
1:B:697:ILE:CG2	1:B:732:ILE:HD11	2.27	0.65
1:D:735:VAL:HA	1:D:738:SER:HB2	1.79	0.65
1:E:611:THR:O	1:E:615:ILE:HG13	1.97	0.65
1:F:408:LEU:H	1:F:408:LEU:CD1	1.94	0.65
2:S:33:GLY:O	2:S:37:ARG:HG3	1.97	0.65
1:A:73:ASN:O	1:A:74:GLU:O	2.14	0.65
1:A:210:PHE:C	1:A:212:GLN:H	2.05	0.65
1:C:718:ARG:NH1	1:C:767:GLN:HE21	1.95	0.65
1:E:246:SER:O	1:E:250:ALA:HB2	1.97	0.65
1:F:678:VAL:HG13	1:F:745:TYR:CD2	2.31	0.65
2:O:21:LYS:HD3	2:O:21:LYS:C	2.22	0.65
2:P:117:THR:HG23	2:P:120:GLU:HB2	1.79	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:R:21:LYS:HD3	2:R:21:LYS:C	2.21	0.65
1:A:254:ARG:H	1:A:254:ARG:HD2	1.60	0.65
1:A:567:THR:CG2	1:A:568:GLY:N	2.59	0.65
1:A:735:VAL:HA	1:A:738:SER:HB2	1.79	0.65
1:B:305:SER:HB2	1:B:594:PHE:CD1	2.32	0.65
1:B:657:ILE:HG13	1:B:756:ILE:CD1	2.27	0.65
1:C:305:SER:HB2	1:C:594:PHE:CD1	2.32	0.65
1:C:692:GLU:HA	1:C:692:GLU:OE2	1.96	0.65
1:D:567:THR:CG2	1:D:568:GLY:N	2.59	0.65
1:E:189:ASP:O	1:E:191:GLU:N	2.25	0.65
1:E:688:PHE:C	1:E:688:PHE:CD2	2.74	0.65
1:E:711:ILE:HG13	1:E:712:PHE:CD2	2.32	0.65
2:P:65:PHE:HB2	2:P:66:PRO:HD3	1.79	0.65
2:T:13:LYS:C	2:T:15:ALA:N	2.54	0.65
1:A:326:ILE:HG22	1:A:328:PHE:CE1	2.32	0.64
1:A:688:PHE:C	1:A:688:PHE:CD2	2.75	0.64
1:C:243:LEU:HA	1:C:246:SER:OG	1.97	0.64
1:C:472:ARG:HH11	1:C:472:ARG:CB	2.07	0.64
1:E:692:GLU:HA	1:E:692:GLU:OE2	1.96	0.64
1:F:189:ASP:O	1:F:191:GLU:N	2.25	0.64
1:F:297:LYS:HZ3	1:F:601:GLU:HB3	1.62	0.64
1:F:350:VAL:HG12	1:F:352:GLY:H	1.62	0.64
1:F:719:LYS:HG2	1:F:797:ILE:HD11	1.78	0.64
2:Q:65:PHE:HB2	2:Q:66:PRO:HD3	1.79	0.64
1:A:130:SER:HB2	1:A:170:TYR:CE2	2.32	0.64
1:A:246:SER:O	1:A:250:ALA:HB2	1.97	0.64
1:A:711:ILE:HG13	1:A:712:PHE:CD2	2.32	0.64
1:B:130:SER:HB2	1:B:170:TYR:CE2	2.32	0.64
1:C:385:LEU:O	1:C:385:LEU:HD13	1.98	0.64
1:C:719:LYS:HG2	1:C:797:ILE:HD11	1.78	0.64
1:E:657:ILE:HG13	1:E:756:ILE:CD1	2.27	0.64
1:F:246:SER:O	1:F:250:ALA:HB2	1.98	0.64
1:F:711:ILE:HG13	1:F:712:PHE:CD2	2.32	0.64
2:P:72:MET:O	2:P:76:MET:CB	2.40	0.64
2:Q:11:GLU:O	2:Q:13:LYS:N	2.30	0.64
1:A:397:GLU:HG3	1:A:480:ASN:HB3	1.78	0.64
1:D:326:ILE:HG22	1:D:328:PHE:CE1	2.32	0.64
1:D:657:ILE:HG13	1:D:756:ILE:CD1	2.27	0.64
1:E:629:ASN:HB3	1:E:632:TYR:CD1	2.33	0.64
2:S:55:VAL:HG21	2:S:67:GLU:OE1	1.98	0.64
2:T:117:THR:HG23	2:T:120:GLU:HB2	1.79	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:199:LEU:C	1:A:201:ASP:H	2.05	0.64
1:A:478:ALA:HB1	1:A:486:LYS:C	2.23	0.64
1:B:335:ALA:O	1:B:339:ILE:HG13	1.98	0.64
1:B:678:VAL:HG13	1:B:745:TYR:CD2	2.33	0.64
1:E:275:GLY:HA2	1:E:278:LYS:CG	2.28	0.64
1:E:697:ILE:CD1	1:E:732:ILE:HD13	2.27	0.64
1:E:788:ASP:O	1:E:792:VAL:HG23	1.98	0.64
2:O:126:ARG:HH21	2:O:126:ARG:HG3	1.62	0.64
2:Q:12:PHE:HD1	2:Q:72:MET:HG3	1.59	0.64
2:Q:33:GLY:O	2:Q:37:ARG:HG3	1.98	0.64
2:T:110:THR:O	2:T:113:GLY:N	2.30	0.64
1:A:122:GLU:HB2	1:A:147:ARG:HG3	1.77	0.64
1:A:238:GLN:C	1:A:240:ALA:H	2.04	0.64
1:B:89:ILE:HD13	1:B:175:LYS:HE2	1.79	0.64
1:B:246:SER:O	1:B:250:ALA:HB2	1.96	0.64
1:B:355:SER:CB	1:B:371:SER:HA	2.27	0.64
1:C:296:LEU:O	1:C:301:ALA:HB2	1.98	0.64
1:D:357:TRP:HZ3	1:D:439:ASN:HB2	1.62	0.64
1:E:115:LYS:HB2	1:E:118:GLN:CG	2.28	0.64
1:E:567:THR:CG2	1:E:568:GLY:N	2.60	0.64
1:E:579:THR:O	1:E:581:GLN:N	2.31	0.64
1:F:199:LEU:C	1:F:201:ASP:H	2.05	0.64
1:F:736:LEU:HD11	1:F:750:GLN:NE2	2.13	0.64
2:O:58:ASP:CB	2:O:62:THR:HG23	2.26	0.64
1:C:736:LEU:HD11	1:C:750:GLN:NE2	2.13	0.64
1:D:235:THR:O	1:D:238:GLN:HB2	1.97	0.64
1:E:472:ARG:HH11	1:E:472:ARG:CB	2.07	0.64
2:S:11:GLU:O	2:S:13:LYS:N	2.30	0.64
1:A:297:LYS:NZ	1:A:601:GLU:HB3	2.13	0.64
1:A:350:VAL:HG12	1:A:352:GLY:H	1.62	0.64
1:A:550:SER:HG	1:A:552:TRP:CD1	2.16	0.64
1:A:611:THR:O	1:A:615:ILE:HG13	1.97	0.64
1:C:360:VAL:CG1	1:C:370:LEU:HD22	2.26	0.64
1:D:350:VAL:HG12	1:D:352:GLY:H	1.63	0.64
1:F:210:PHE:C	1:F:212:GLN:H	2.06	0.64
2:P:21:LYS:C	2:P:21:LYS:HD3	2.22	0.64
2:S:137:ASN:OD1	2:S:140:GLU:HG2	1.98	0.64
2:T:13:LYS:HZ3	2:T:65:PHE:HB3	1.61	0.64
1:A:127:SER:O	1:A:133:GLU:OE2	2.16	0.64
1:A:275:GLY:HA2	1:A:278:LYS:CG	2.28	0.64
1:A:579:THR:O	1:A:581:GLN:N	2.31	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:92:ASP:O	1:B:96:ILE:HG13	1.97	0.64
1:B:296:LEU:O	1:B:301:ALA:HB2	1.97	0.64
1:B:736:LEU:HD11	1:B:750:GLN:NE2	2.13	0.64
1:C:238:GLN:C	1:C:240:ALA:H	2.04	0.64
1:C:357:TRP:HZ3	1:C:439:ASN:HB2	1.62	0.64
1:C:711:ILE:HG13	1:C:712:PHE:CD2	2.32	0.64
1:D:127:SER:O	1:D:133:GLU:OE2	2.16	0.64
1:D:238:GLN:C	1:D:240:ALA:H	2.04	0.64
1:D:243:LEU:HA	1:D:246:SER:OG	1.97	0.64
1:D:297:LYS:NZ	1:D:601:GLU:HB3	2.13	0.64
1:D:397:GLU:HG3	1:D:480:ASN:HB3	1.79	0.64
1:D:711:ILE:HG13	1:D:712:PHE:CD2	2.32	0.64
2:R:49:GLN:N	2:R:49:GLN:NE2	2.46	0.64
1:B:127:SER:O	1:B:133:GLU:OE2	2.16	0.64
1:B:550:SER:HG	1:B:552:TRP:CD1	2.16	0.64
1:B:735:VAL:HA	1:B:738:SER:HB2	1.80	0.64
1:C:397:GLU:HG3	1:C:480:ASN:HB3	1.78	0.64
1:E:243:LEU:HA	1:E:246:SER:OG	1.98	0.64
1:E:305:SER:HB2	1:E:594:PHE:CD1	2.32	0.64
1:E:385:LEU:O	1:E:385:LEU:HD13	1.97	0.64
1:E:718:ARG:NH1	1:E:767:GLN:HE21	1.96	0.64
1:F:97:TYR:OH	1:F:150:PRO:HB2	1.98	0.64
1:F:297:LYS:NZ	1:F:601:GLU:HB3	2.12	0.64
1:F:335:ALA:O	1:F:339:ILE:HG13	1.97	0.64
2:P:58:ASP:CB	2:P:62:THR:HG23	2.26	0.64
2:T:126:ARG:HG3	2:T:126:ARG:HH21	1.63	0.64
1:B:187:SER:O	1:B:188:LEU:C	2.37	0.64
1:B:397:GLU:HG3	1:B:480:ASN:HB3	1.78	0.64
1:C:130:SER:HB2	1:C:170:TYR:CE2	2.32	0.64
1:C:297:LYS:NZ	1:C:601:GLU:HB3	2.13	0.64
1:C:355:SER:CB	1:C:371:SER:HA	2.28	0.64
1:C:478:ALA:HB1	1:C:486:LYS:C	2.23	0.64
1:D:478:ALA:HB1	1:D:486:LYS:C	2.23	0.64
1:D:597:ASN:ND2	1:D:601:GLU:N	2.46	0.64
1:E:92:ASP:O	1:E:96:ILE:HG13	1.98	0.64
1:E:127:SER:O	1:E:133:GLU:OE2	2.16	0.64
1:E:735:VAL:HA	1:E:738:SER:HB2	1.80	0.64
2:Q:49:GLN:N	2:Q:49:GLN:NE2	2.45	0.64
1:A:718:ARG:NH1	1:A:767:GLN:HE21	1.96	0.63
1:B:630:ARG:NE	2:P:83:GLU:HG2	2.13	0.63
1:B:788:ASP:O	1:B:792:VAL:HG23	1.98	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:108:ASP:O	1:C:155:ASN:HB2	1.98	0.63
1:C:210:PHE:C	1:C:212:GLN:H	2.06	0.63
1:C:607:ASN:HB3	1:C:609:GLU:OE2	1.98	0.63
1:D:597:ASN:HD21	1:D:601:GLU:N	1.96	0.63
1:D:697:ILE:CG2	1:D:732:ILE:HD11	2.28	0.63
1:E:199:LEU:C	1:E:201:ASP:H	2.06	0.63
2:O:33:GLY:O	2:O:37:ARG:HG3	1.98	0.63
2:Q:21:LYS:C	2:Q:21:LYS:HD3	2.22	0.63
2:S:126:ARG:HG3	2:S:126:ARG:HH21	1.63	0.63
2:T:65:PHE:HB2	2:T:66:PRO:HD3	1.79	0.63
1:A:355:SER:CB	1:A:371:SER:HA	2.28	0.63
1:A:597:ASN:HD21	1:A:601:GLU:N	1.96	0.63
1:A:692:GLU:OE2	1:A:692:GLU:HA	1.97	0.63
1:B:243:LEU:HA	1:B:246:SER:OG	1.97	0.63
1:B:350:VAL:HG12	1:B:352:GLY:H	1.62	0.63
1:B:597:ASN:HD21	1:B:601:GLU:N	1.96	0.63
1:C:275:GLY:HA2	1:C:278:LYS:CG	2.27	0.63
1:C:326:ILE:HG22	1:C:328:PHE:CE1	2.33	0.63
1:D:275:GLY:HA2	1:D:278:LYS:CG	2.28	0.63
1:D:355:SER:CB	1:D:371:SER:HA	2.29	0.63
1:E:397:GLU:HG3	1:E:480:ASN:HB3	1.79	0.63
1:E:478:ALA:HB1	1:E:486:LYS:C	2.24	0.63
1:F:127:SER:O	1:F:133:GLU:OE2	2.16	0.63
1:F:275:GLY:HA2	1:F:278:LYS:CG	2.27	0.63
1:F:405:LEU:HD13	1:F:453:VAL:HG21	1.81	0.63
2:Q:117:THR:HG23	2:Q:120:GLU:HB2	1.79	0.63
2:T:49:GLN:N	2:T:49:GLN:NE2	2.45	0.63
1:C:788:ASP:O	1:C:792:VAL:HG23	1.97	0.63
1:E:130:SER:HB2	1:E:170:TYR:CE2	2.33	0.63
1:E:186:LYS:C	1:E:188:LEU:O	2.41	0.63
1:F:187:SER:O	1:F:188:LEU:C	2.37	0.63
2:P:11:GLU:O	2:P:13:LYS:N	2.30	0.63
2:P:55:VAL:HG21	2:P:67:GLU:OE1	1.98	0.63
1:A:296:LEU:O	1:A:301:ALA:HB2	1.99	0.63
1:A:405:LEU:HD13	1:A:453:VAL:HG21	1.81	0.63
1:A:597:ASN:ND2	1:A:601:GLU:N	2.46	0.63
1:A:697:ILE:CG2	1:A:732:ILE:HD11	2.29	0.63
1:B:97:TYR:OH	1:B:150:PRO:HB2	1.99	0.63
1:F:355:SER:CB	1:F:371:SER:HA	2.28	0.63
2:Q:13:LYS:C	2:Q:15:ALA:N	2.56	0.63
2:R:58:ASP:CB	2:R:62:THR:HG23	2.26	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:692:GLU:OE2	1:D:692:GLU:HA	1.98	0.63
1:D:788:ASP:O	1:D:792:VAL:HG23	1.97	0.63
1:E:187:SER:O	1:E:188:LEU:C	2.37	0.63
1:E:225:ILE:HA	1:E:229:PHE:HE2	1.64	0.63
1:F:296:LEU:O	1:F:301:ALA:HB2	1.98	0.63
1:F:478:ALA:HB1	1:F:486:LYS:C	2.24	0.63
1:F:788:ASP:O	1:F:792:VAL:HG23	1.98	0.63
2:R:13:LYS:HZ3	2:R:65:PHE:HB3	1.63	0.63
2:S:110:THR:O	2:S:113:GLY:N	2.31	0.63
1:A:88:LYS:NZ	1:A:172:GLU:OE2	2.31	0.63
1:B:199:LEU:C	1:B:201:ASP:H	2.06	0.63
1:B:225:ILE:HA	1:B:229:PHE:HE2	1.64	0.63
1:C:179:LEU:O	1:C:182:ILE:HG22	1.98	0.63
1:D:296:LEU:O	1:D:301:ALA:HB2	1.99	0.63
1:D:376:GLN:HB3	1:D:379:ALA:HB3	1.81	0.63
1:D:405:LEU:HD13	1:D:453:VAL:HG21	1.80	0.63
1:D:697:ILE:C	1:D:699:GLY:H	2.07	0.63
1:F:301:ALA:C	1:F:303:LYS:N	2.53	0.63
2:O:11:GLU:O	2:O:13:LYS:N	2.31	0.63
2:Q:126:ARG:HH21	2:Q:126:ARG:HG3	1.63	0.63
2:R:33:GLY:O	2:R:37:ARG:HG3	1.99	0.63
1:A:788:ASP:O	1:A:792:VAL:HG23	1.98	0.63
1:B:275:GLY:HA2	1:B:278:LYS:CG	2.28	0.63
1:B:368:GLN:HG3	1:B:383:GLY:C	2.24	0.63
1:C:127:SER:O	1:C:133:GLU:OE2	2.16	0.63
1:C:697:ILE:CG2	1:C:732:ILE:HD11	2.29	0.63
1:D:129:ASN:HD22	1:D:129:ASN:N	1.97	0.63
1:D:678:VAL:HG13	1:D:745:TYR:CD2	2.33	0.63
1:E:350:VAL:HG12	1:E:352:GLY:H	1.63	0.63
1:F:92:ASP:O	1:F:96:ILE:HG13	1.97	0.63
1:F:130:SER:HB2	1:F:170:TYR:CE2	2.33	0.63
1:F:697:ILE:CG2	1:F:732:ILE:HD11	2.29	0.63
2:O:65:PHE:HB2	2:O:66:PRO:HD3	1.80	0.63
2:S:49:GLN:N	2:S:49:GLN:NE2	2.46	0.63
1:B:115:LYS:HB2	1:B:118:GLN:CG	2.29	0.63
1:C:115:LYS:HB2	1:C:118:GLN:CG	2.29	0.63
1:D:246:SER:O	1:D:250:ALA:HB2	1.98	0.63
1:E:297:LYS:NZ	1:E:601:GLU:HB3	2.13	0.63
1:F:368:GLN:HG3	1:F:383:GLY:C	2.24	0.63
1:F:630:ARG:NE	2:T:83:GLU:HG2	2.14	0.63
2:S:65:PHE:HB2	2:S:66:PRO:HD3	1.80	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:97:TYR:OH	1:A:150:PRO:HB2	1.99	0.63
1:A:597:ASN:HD22	1:A:601:GLU:HB2	1.64	0.63
1:B:297:LYS:NZ	1:B:601:GLU:HB3	2.13	0.63
1:C:97:TYR:OH	1:C:150:PRO:HB2	1.99	0.63
1:C:129:ASN:HD22	1:C:129:ASN:N	1.97	0.63
1:C:597:ASN:HD22	1:C:601:GLU:HB2	1.63	0.63
1:D:210:PHE:C	1:D:212:GLN:H	2.07	0.63
1:E:355:SER:CB	1:E:371:SER:HA	2.27	0.63
1:E:368:GLN:HG3	1:E:383:GLY:C	2.24	0.63
1:F:225:ILE:HA	1:F:229:PHE:HE2	1.64	0.63
1:A:92:ASP:O	1:A:96:ILE:HG13	1.99	0.62
1:A:97:TYR:CD2	1:A:102:GLY:HA3	2.34	0.62
1:B:478:ALA:HB1	1:B:486:LYS:C	2.24	0.62
1:E:105:TYR:N	1:E:152:LEU:O	2.30	0.62
1:F:238:GLN:C	1:F:240:ALA:H	2.06	0.62
1:F:288:VAL:HG23	1:F:289:GLU:H	1.64	0.62
1:F:607:ASN:HB3	1:F:609:GLU:OE2	1.99	0.62
1:F:700:TYR:CD1	1:F:727:GLN:HB3	2.34	0.62
2:Q:55:VAL:HG21	2:Q:67:GLU:OE1	1.98	0.62
2:R:11:GLU:O	2:R:13:LYS:N	2.32	0.62
1:A:630:ARG:NE	2:O:83:GLU:HG2	2.14	0.62
1:B:93:VAL:HA	1:B:96:ILE:CD1	2.29	0.62
1:B:97:TYR:C	1:B:99:GLU:H	2.08	0.62
1:B:607:ASN:HB3	1:B:609:GLU:OE2	1.99	0.62
1:C:225:ILE:HA	1:C:229:PHE:HE2	1.64	0.62
1:C:368:GLN:HG3	1:C:383:GLY:C	2.24	0.62
1:C:463:THR:HB	1:C:467:GLU:H	1.64	0.62
1:C:678:VAL:HG13	1:C:745:TYR:CD2	2.33	0.62
1:D:97:TYR:CD2	1:D:102:GLY:HA3	2.34	0.62
1:D:179:LEU:O	1:D:182:ILE:HG22	1.99	0.62
1:E:607:ASN:HB3	1:E:609:GLU:OE2	1.98	0.62
1:A:296:LEU:HD22	1:A:606:LYS:HE2	1.81	0.62
1:A:413:LEU:N	1:A:413:LEU:HD23	2.14	0.62
1:A:607:ASN:HB3	1:A:609:GLU:OE2	1.98	0.62
1:A:736:LEU:HD11	1:A:750:GLN:NE2	2.13	0.62
1:B:115:LYS:CB	1:B:118:GLN:HG2	2.29	0.62
1:D:105:TYR:N	1:D:152:LEU:O	2.30	0.62
1:D:630:ARG:NE	2:R:83:GLU:HG2	2.14	0.62
1:E:424:LYS:HZ2	1:E:424:LYS:HB3	1.64	0.62
1:E:597:ASN:ND2	1:E:601:GLU:N	2.47	0.62
1:E:630:ARG:NE	2:S:83:GLU:HG2	2.13	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:108:ASP:O	1:F:155:ASN:HB2	1.99	0.62
2:R:13:LYS:CE	2:R:65:PHE:HB3	2.29	0.62
2:R:55:VAL:HG21	2:R:67:GLU:OE1	1.99	0.62
2:R:65:PHE:HB2	2:R:66:PRO:HD3	1.79	0.62
2:S:117:THR:HG23	2:S:120:GLU:HB2	1.81	0.62
2:T:137:ASN:OD1	2:T:140:GLU:HG2	1.99	0.62
1:A:427:ASP:O	1:A:428:ASN:HB2	1.98	0.62
1:B:405:LEU:HD13	1:B:453:VAL:HG21	1.80	0.62
1:D:92:ASP:O	1:D:96:ILE:HG13	1.99	0.62
1:D:607:ASN:HB3	1:D:609:GLU:OE2	1.99	0.62
1:E:115:LYS:CB	1:E:118:GLN:HG2	2.29	0.62
1:E:296:LEU:O	1:E:301:ALA:HB2	1.99	0.62
1:E:446:ILE:HG13	1:E:452:GLU:O	1.99	0.62
1:E:522:SER:O	2:S:124:MET:HE2	1.99	0.62
1:F:115:LYS:HB2	1:F:118:GLN:CG	2.30	0.62
1:F:296:LEU:HD22	1:F:606:LYS:HE2	1.80	0.62
1:F:463:THR:HB	1:F:467:GLU:H	1.65	0.62
2:S:72:MET:O	2:S:76:MET:CB	2.40	0.62
1:A:129:ASN:N	1:A:129:ASN:HD22	1.97	0.62
1:B:108:ASP:O	1:B:155:ASN:HB2	2.00	0.62
1:C:296:LEU:HD22	1:C:606:LYS:HE2	1.82	0.62
1:E:97:TYR:C	1:E:99:GLU:H	2.07	0.62
1:E:108:ASP:O	1:E:155:ASN:HB2	2.00	0.62
1:E:597:ASN:HD21	1:E:601:GLU:N	1.97	0.62
1:E:697:ILE:CG2	1:E:732:ILE:HD11	2.29	0.62
1:E:736:LEU:HD11	1:E:750:GLN:NE2	2.14	0.62
1:F:97:TYR:CD2	1:F:102:GLY:HA3	2.33	0.62
1:F:376:GLN:HB3	1:F:379:ALA:HB3	1.81	0.62
1:F:407:HIS:HB2	1:F:408:LEU:HD12	1.81	0.62
2:R:110:THR:O	2:R:113:GLY:N	2.30	0.62
2:S:21:LYS:C	2:S:21:LYS:HD3	2.22	0.62
2:T:11:GLU:O	2:T:13:LYS:N	2.31	0.62
1:A:108:ASP:O	1:A:155:ASN:HB2	1.99	0.62
1:B:296:LEU:HD22	1:B:606:LYS:HE2	1.81	0.62
1:B:700:TYR:CD1	1:B:727:GLN:HB3	2.35	0.62
1:C:697:ILE:C	1:C:699:GLY:H	2.08	0.62
1:D:97:TYR:C	1:D:99:GLU:H	2.08	0.62
1:D:115:LYS:HB2	1:D:118:GLN:CG	2.29	0.62
1:E:405:LEU:HD13	1:E:453:VAL:HG21	1.82	0.62
1:E:736:LEU:HD21	1:E:750:GLN:NE2	2.15	0.62
2:O:13:LYS:CE	2:O:65:PHE:HB3	2.29	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:P:49:GLN:N	2:P:49:GLN:NE2	2.45	0.62
1:A:376:GLN:HB3	1:A:379:ALA:HB3	1.81	0.62
1:C:97:TYR:CD2	1:C:102:GLY:HA3	2.34	0.62
1:D:597:ASN:HD22	1:D:601:GLU:HB2	1.64	0.62
1:D:629:ASN:HB3	1:D:632:TYR:CD1	2.35	0.62
1:E:93:VAL:HA	1:E:96:ILE:CD1	2.29	0.62
1:F:597:ASN:ND2	1:F:601:GLU:N	2.47	0.62
2:S:13:LYS:CE	2:S:65:PHE:HB3	2.30	0.62
1:A:115:LYS:CB	1:A:118:GLN:HG2	2.30	0.62
1:A:161:ILE:HG23	1:A:168:GLU:HB2	1.81	0.62
1:A:172:GLU:HB3	1:A:246:SER:HA	1.82	0.62
1:A:238:GLN:C	1:A:240:ALA:N	2.58	0.62
1:A:368:GLN:HG3	1:A:383:GLY:C	2.24	0.62
1:A:446:ILE:HG13	1:A:452:GLU:O	1.99	0.62
1:A:700:TYR:CD1	1:A:727:GLN:HB3	2.35	0.62
1:B:597:ASN:ND2	1:B:601:GLU:N	2.47	0.62
1:B:697:ILE:C	1:B:699:GLY:H	2.07	0.62
1:C:376:GLN:HB3	1:C:379:ALA:HB3	1.82	0.62
1:D:225:ILE:HA	1:D:229:PHE:HE2	1.64	0.62
1:D:305:SER:HB2	1:D:594:PHE:CD1	2.34	0.62
1:E:288:VAL:HG23	1:E:289:GLU:H	1.64	0.62
2:O:13:LYS:HZ3	2:O:65:PHE:HB3	1.63	0.62
2:O:117:THR:HG23	2:O:120:GLU:HB2	1.81	0.62
2:R:94:LYS:HE3	2:R:107:HIS:CD2	2.34	0.62
1:A:225:ILE:HA	1:A:229:PHE:HE2	1.64	0.62
1:A:332:ASN:HD21	1:A:334:LEU:HB2	1.65	0.62
1:D:165:GLN:C	1:D:167:LYS:H	2.08	0.62
1:F:172:GLU:HB3	1:F:246:SER:HA	1.82	0.62
1:F:597:ASN:HD21	1:F:601:GLU:N	1.96	0.62
1:F:629:ASN:HB3	1:F:632:TYR:CD1	2.34	0.62
2:Q:110:THR:O	2:Q:113:GLY:N	2.32	0.62
2:T:55:VAL:HG21	2:T:67:GLU:OE1	1.99	0.62
1:A:93:VAL:HA	1:A:96:ILE:CD1	2.30	0.62
1:C:97:TYR:C	1:C:99:GLU:H	2.07	0.62
1:C:718:ARG:O	1:C:722:ILE:HG13	2.00	0.62
1:D:115:LYS:CB	1:D:118:GLN:HG2	2.30	0.62
1:D:296:LEU:HD22	1:D:606:LYS:HE2	1.82	0.62
1:D:636:ALA:O	1:D:640:LYS:HA	2.00	0.62
1:E:318:ILE:HG23	1:E:322:LEU:HD12	1.82	0.62
1:E:413:LEU:N	1:E:413:LEU:HD23	2.15	0.62
1:E:463:THR:HB	1:E:467:GLU:H	1.65	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:413:LEU:N	1:F:413:LEU:HD23	2.15	0.62
1:F:597:ASN:HD22	1:F:601:GLU:HB2	1.65	0.62
2:R:117:THR:HG23	2:R:120:GLU:HB2	1.80	0.62
1:A:115:LYS:HB2	1:A:118:GLN:CG	2.29	0.61
1:B:165:GLN:C	1:B:167:LYS:H	2.07	0.61
1:B:463:THR:HB	1:B:467:GLU:H	1.65	0.61
1:D:446:ILE:HG13	1:D:452:GLU:O	1.99	0.61
1:D:700:TYR:CD1	1:D:727:GLN:HB3	2.35	0.61
1:E:161:ILE:HG23	1:E:168:GLU:HB2	1.80	0.61
2:P:110:THR:O	2:P:113:GLY:N	2.31	0.61
1:B:446:ILE:HG13	1:B:452:GLU:O	2.01	0.61
1:B:718:ARG:O	1:B:722:ILE:HG13	2.00	0.61
1:C:187:SER:O	1:C:188:LEU:C	2.37	0.61
1:C:288:VAL:HG23	1:C:289:GLU:H	1.65	0.61
1:C:529:VAL:HG21	2:Q:109:MET:HE1	1.82	0.61
1:C:629:ASN:HB3	1:C:632:TYR:CD1	2.35	0.61
1:D:718:ARG:O	1:D:722:ILE:HG13	2.00	0.61
1:E:97:TYR:CD2	1:E:102:GLY:HA3	2.34	0.61
1:E:700:TYR:CD1	1:E:727:GLN:HB3	2.35	0.61
1:F:97:TYR:C	1:F:99:GLU:H	2.07	0.61
1:F:129:ASN:HD22	1:F:129:ASN:N	1.97	0.61
2:Q:58:ASP:CB	2:Q:62:THR:HG23	2.26	0.61
2:Q:138:TYR:O	2:Q:142:VAL:HG23	2.00	0.61
2:R:137:ASN:OD1	2:R:140:GLU:HG2	2.01	0.61
1:A:179:LEU:O	1:A:182:ILE:HG22	1.99	0.61
1:A:407:HIS:HB2	1:A:408:LEU:HD12	1.82	0.61
1:B:413:LEU:N	1:B:413:LEU:HD23	2.16	0.61
1:B:718:ARG:NH1	1:B:767:GLN:HE21	1.98	0.61
1:C:405:LEU:HD13	1:C:453:VAL:HG21	1.81	0.61
1:C:630:ARG:NE	2:Q:83:GLU:HG2	2.15	0.61
1:C:700:TYR:CD1	1:C:727:GLN:HB3	2.35	0.61
1:D:238:GLN:C	1:D:240:ALA:N	2.58	0.61
1:D:736:LEU:HD11	1:D:750:GLN:NE2	2.14	0.61
1:E:298:GLY:C	1:E:300:LYS:H	2.08	0.61
1:E:407:HIS:HB2	1:E:408:LEU:HD12	1.81	0.61
1:F:298:GLY:C	1:F:300:LYS:H	2.09	0.61
2:P:13:LYS:HZ3	2:P:65:PHE:HB3	1.64	0.61
1:B:97:TYR:CD2	1:B:102:GLY:HA3	2.36	0.61
1:B:161:ILE:HG23	1:B:168:GLU:HB2	1.81	0.61
1:B:179:LEU:O	1:B:182:ILE:HG22	1.99	0.61
1:B:407:HIS:HB2	1:B:408:LEU:HD12	1.81	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:92:ASP:O	1:C:96:ILE:HG13	2.00	0.61
1:C:298:GLY:C	1:C:300:LYS:H	2.08	0.61
1:C:636:ALA:O	1:C:640:LYS:HA	2.00	0.61
1:D:88:LYS:HZ3	1:D:172:GLU:CD	2.08	0.61
1:E:97:TYR:OH	1:E:150:PRO:HB2	2.00	0.61
2:P:13:LYS:CE	2:P:65:PHE:HB3	2.30	0.61
2:R:97:ASN:ND2	2:R:99:TYR:H	1.99	0.61
1:C:115:LYS:C	1:C:117:LEU:N	2.58	0.61
1:C:332:ASN:HD21	1:C:334:LEU:HB2	1.65	0.61
1:C:579:THR:O	1:C:581:GLN:N	2.33	0.61
1:D:318:ILE:HG23	1:D:322:LEU:HD12	1.82	0.61
1:D:413:LEU:HD23	1:D:413:LEU:N	2.14	0.61
1:E:88:LYS:NZ	1:E:172:GLU:OE2	2.31	0.61
1:E:179:LEU:O	1:E:182:ILE:HG22	2.00	0.61
1:E:332:ASN:HD21	1:E:334:LEU:HB2	1.66	0.61
1:B:217:LYS:HB3	1:B:217:LYS:NZ	2.12	0.61
1:C:115:LYS:CB	1:C:118:GLN:HG2	2.30	0.61
1:C:710:HIS:C	1:C:712:PHE:N	2.59	0.61
1:D:108:ASP:O	1:D:155:ASN:HB2	2.00	0.61
1:D:463:THR:HB	1:D:467:GLU:H	1.65	0.61
1:E:298:GLY:O	1:E:300:LYS:N	2.34	0.61
1:F:636:ALA:O	1:F:640:LYS:HA	2.01	0.61
1:A:97:TYR:C	1:A:99:GLU:H	2.08	0.61
1:A:401:ILE:HD13	1:A:485:LEU:O	2.01	0.61
1:A:597:ASN:OD1	1:A:599:GLU:N	2.31	0.61
1:B:172:GLU:HB3	1:B:246:SER:HA	1.82	0.61
1:C:424:LYS:HZ2	1:C:424:LYS:HB3	1.65	0.61
1:D:368:GLN:HG3	1:D:383:GLY:C	2.25	0.61
1:D:609:GLU:OE2	1:D:609:GLU:N	2.31	0.61
1:E:376:GLN:HB3	1:E:379:ALA:HB3	1.81	0.61
1:F:275:GLY:O	1:F:278:LYS:HB2	2.01	0.61
2:T:13:LYS:CE	2:T:65:PHE:HB3	2.29	0.61
1:A:298:GLY:C	1:A:300:LYS:H	2.09	0.61
1:A:424:LYS:HZ2	1:A:424:LYS:HB3	1.65	0.61
1:B:579:THR:O	1:B:581:GLN:N	2.33	0.61
1:C:597:ASN:HD21	1:C:601:GLU:N	1.98	0.61
1:D:141:PHE:HD1	1:D:141:PHE:N	1.99	0.61
1:E:172:GLU:HB3	1:E:246:SER:HA	1.83	0.61
1:E:597:ASN:HD22	1:E:601:GLU:HB2	1.65	0.61
1:F:446:ILE:HG13	1:F:452:GLU:O	2.00	0.61
1:F:579:THR:O	1:F:581:GLN:N	2.34	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:O:13:LYS:C	2:O:15:ALA:N	2.56	0.61
2:S:13:LYS:C	2:S:15:ALA:N	2.57	0.61
1:A:109:ILE:HD13	1:A:157:LYS:NZ	2.16	0.61
1:A:275:GLY:CA	1:A:278:LYS:HG3	2.31	0.61
1:A:463:THR:HB	1:A:467:GLU:H	1.65	0.61
1:B:275:GLY:O	1:B:278:LYS:HB2	2.01	0.61
1:C:238:GLN:C	1:C:240:ALA:N	2.58	0.61
1:C:479:LYS:CG	1:C:488:LEU:HD21	2.30	0.61
1:C:597:ASN:ND2	1:C:601:GLU:N	2.48	0.61
1:D:522:SER:O	2:R:124:MET:HE2	2.00	0.61
1:D:579:THR:O	1:D:581:GLN:N	2.34	0.61
1:E:179:LEU:HG	1:E:180:ASP:N	2.16	0.61
1:F:88:LYS:NZ	1:F:172:GLU:OE2	2.32	0.61
1:A:122:GLU:HG2	1:A:122:GLU:O	2.00	0.61
1:A:318:ILE:HG23	1:A:322:LEU:HD12	1.83	0.61
1:A:718:ARG:O	1:A:722:ILE:HG13	2.00	0.61
1:B:112:VAL:C	1:B:114:HIS:H	2.09	0.61
1:B:343:VAL:HG13	1:B:487:PRO:HG2	1.83	0.61
1:D:161:ILE:HG23	1:D:168:GLU:HB2	1.81	0.61
1:D:759:GLN:HA	1:D:759:GLN:HE21	1.66	0.61
1:E:141:PHE:HD1	1:E:141:PHE:N	1.98	0.61
1:E:275:GLY:O	1:E:278:LYS:HB2	2.01	0.61
1:E:301:ALA:C	1:E:303:LYS:N	2.52	0.61
1:F:115:LYS:CB	1:F:118:GLN:HG2	2.30	0.61
1:F:122:GLU:CD	1:F:147:ARG:HB2	2.26	0.61
1:F:298:GLY:O	1:F:300:LYS:N	2.34	0.61
1:F:332:ASN:HD21	1:F:334:LEU:HB2	1.66	0.61
1:F:424:LYS:HZ2	1:F:424:LYS:HB3	1.65	0.61
2:Q:13:LYS:CE	2:Q:65:PHE:HB3	2.30	0.61
2:S:83:GLU:O	2:S:87:GLU:HG3	2.01	0.61
1:A:141:PHE:HD1	1:A:141:PHE:N	1.98	0.60
1:A:636:ALA:O	1:A:640:LYS:HA	2.01	0.60
1:A:671:ARG:HD2	2:O:14:GLU:CG	2.31	0.60
1:B:88:LYS:NZ	1:B:172:GLU:OE2	2.33	0.60
1:B:115:LYS:C	1:B:117:LEU:N	2.58	0.60
1:B:122:GLU:CD	1:B:147:ARG:HB2	2.26	0.60
1:C:161:ILE:HG23	1:C:168:GLU:HB2	1.83	0.60
1:C:413:LEU:N	1:C:413:LEU:HD23	2.15	0.60
1:D:179:LEU:HG	1:D:180:ASP:N	2.16	0.60
1:D:298:GLY:C	1:D:300:LYS:H	2.09	0.60
1:D:343:VAL:HG13	1:D:487:PRO:HG2	1.82	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:479:LYS:CG	1:D:488:LEU:HD21	2.30	0.60
1:E:122:GLU:O	1:E:122:GLU:HG2	2.00	0.60
2:R:12:PHE:HD1	2:R:72:MET:HG3	1.60	0.60
1:B:129:ASN:HD22	1:B:129:ASN:N	1.97	0.60
1:B:360:VAL:CG1	1:B:370:LEU:HD22	2.27	0.60
1:B:636:ALA:O	1:B:640:LYS:HA	2.01	0.60
1:D:597:ASN:OD1	1:D:599:GLU:N	2.31	0.60
1:E:115:LYS:C	1:E:117:LEU:N	2.57	0.60
1:E:165:GLN:C	1:E:167:LYS:H	2.08	0.60
1:E:293:ILE:O	1:E:295:VAL:HG13	2.01	0.60
1:E:296:LEU:HD22	1:E:606:LYS:HE2	1.82	0.60
1:F:161:ILE:HG23	1:F:168:GLU:HB2	1.82	0.60
1:A:122:GLU:CD	1:A:147:ARG:HB2	2.26	0.60
1:C:122:GLU:CD	1:C:147:ARG:HB2	2.26	0.60
1:C:165:GLN:C	1:C:167:LYS:H	2.08	0.60
1:C:279:ILE:O	1:C:283:LEU:HD13	2.01	0.60
1:D:275:GLY:O	1:D:278:LYS:HB2	2.02	0.60
1:E:759:GLN:HA	1:E:759:GLN:HE21	1.66	0.60
1:F:349:ASN:ND2	1:F:350:VAL:HG23	2.16	0.60
1:F:718:ARG:O	1:F:722:ILE:HG13	2.01	0.60
2:O:12:PHE:HD1	2:O:72:MET:HG3	1.60	0.60
2:P:11:GLU:C	2:P:13:LYS:H	2.09	0.60
1:A:165:GLN:C	1:A:167:LYS:H	2.07	0.60
1:B:179:LEU:HG	1:B:180:ASP:N	2.17	0.60
1:B:629:ASN:HB3	1:B:632:TYR:CD1	2.36	0.60
1:C:217:LYS:CB	1:C:236:GLU:HG3	2.31	0.60
1:D:141:PHE:N	1:D:141:PHE:CD1	2.69	0.60
1:D:710:HIS:O	1:D:712:PHE:N	2.35	0.60
1:F:159:TYR:N	1:F:159:TYR:CD1	2.70	0.60
2:O:100:ILE:HB	2:O:136:VAL:CG2	2.23	0.60
2:O:110:THR:O	2:O:113:GLY:N	2.32	0.60
1:A:141:PHE:N	1:A:141:PHE:CD1	2.69	0.60
1:A:343:VAL:HG13	1:A:487:PRO:HG2	1.82	0.60
1:A:349:ASN:HD22	1:A:350:VAL:H	1.50	0.60
1:A:759:GLN:HA	1:A:759:GLN:HE21	1.66	0.60
1:B:122:GLU:O	1:B:122:GLU:HG2	2.00	0.60
1:B:298:GLY:C	1:B:300:LYS:H	2.09	0.60
1:C:141:PHE:N	1:C:141:PHE:CD1	2.70	0.60
1:C:159:TYR:N	1:C:159:TYR:CD1	2.70	0.60
1:C:736:LEU:HD21	1:C:750:GLN:NE2	2.16	0.60
1:E:112:VAL:C	1:E:114:HIS:H	2.10	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:636:ALA:O	1:E:640:LYS:HA	2.00	0.60
1:E:697:ILE:C	1:E:699:GLY:H	2.07	0.60
1:F:105:TYR:N	1:F:152:LEU:O	2.30	0.60
1:F:115:LYS:C	1:F:117:LEU:N	2.58	0.60
1:F:122:GLU:O	1:F:122:GLU:HG2	2.00	0.60
1:F:179:LEU:O	1:F:182:ILE:HG22	2.00	0.60
1:F:736:LEU:HD21	1:F:750:GLN:NE2	2.17	0.60
2:T:37:ARG:HG2	2:T:37:ARG:HH11	1.67	0.60
1:A:217:LYS:HB3	1:A:217:LYS:NZ	2.12	0.60
1:A:349:ASN:ND2	1:A:350:VAL:HG23	2.17	0.60
1:C:275:GLY:CA	1:C:278:LYS:HG3	2.29	0.60
1:C:759:GLN:HE21	1:C:759:GLN:HA	1.66	0.60
1:D:407:HIS:HB2	1:D:408:LEU:HD12	1.82	0.60
1:F:759:GLN:HE21	1:F:759:GLN:HA	1.66	0.60
2:P:83:GLU:O	2:P:87:GLU:HG3	2.02	0.60
2:Q:37:ARG:HG2	2:Q:37:ARG:HH11	1.67	0.60
2:R:37:ARG:HG2	2:R:37:ARG:HH11	1.67	0.60
1:A:529:VAL:HG21	2:O:109:MET:HE1	1.84	0.60
1:A:629:ASN:HB3	1:A:632:TYR:CD1	2.36	0.60
1:B:88:LYS:HZ3	1:B:172:GLU:CD	2.09	0.60
1:B:376:GLN:HB3	1:B:379:ALA:HB3	1.82	0.60
1:C:141:PHE:HD1	1:C:141:PHE:N	1.99	0.60
1:C:179:LEU:HG	1:C:180:ASP:N	2.17	0.60
1:C:407:HIS:HB2	1:C:408:LEU:HD12	1.82	0.60
1:E:275:GLY:CA	1:E:278:LYS:HG3	2.31	0.60
1:F:710:HIS:O	1:F:712:PHE:N	2.35	0.60
2:O:11:GLU:C	2:O:13:LYS:H	2.08	0.60
2:O:97:ASN:ND2	2:O:99:TYR:H	1.99	0.60
2:P:12:PHE:HE1	2:P:72:MET:CE	2.12	0.60
2:P:49:GLN:O	2:P:51:MET:N	2.35	0.60
2:P:137:ASN:OD1	2:P:140:GLU:HG2	2.01	0.60
1:A:179:LEU:HG	1:A:180:ASP:N	2.17	0.60
1:B:318:ILE:HG23	1:B:322:LEU:HD12	1.82	0.60
1:C:109:ILE:C	1:C:111:LEU:H	2.10	0.60
1:C:173:ILE:HG23	1:C:174:GLY:H	1.66	0.60
1:C:293:ILE:O	1:C:295:VAL:HG13	2.02	0.60
1:D:517:VAL:HG23	1:D:518:ASN:ND2	2.17	0.60
1:E:180:ASP:CG	1:E:181:ILE:N	2.59	0.60
1:E:427:ASP:O	1:E:428:ASN:HB2	2.00	0.60
1:E:529:VAL:HG21	2:S:109:MET:HE1	1.83	0.60
1:E:718:ARG:O	1:E:722:ILE:HG13	2.01	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:165:GLN:C	1:F:167:LYS:H	2.09	0.60
1:B:288:VAL:HG23	1:B:289:GLU:H	1.65	0.60
1:D:597:ASN:ND2	1:D:601:GLU:HB2	2.17	0.60
1:E:343:VAL:HG13	1:E:487:PRO:HG2	1.82	0.60
1:F:176:GLY:C	1:F:178:SER:H	2.10	0.60
1:F:318:ILE:HG23	1:F:322:LEU:HD12	1.83	0.60
2:O:137:ASN:OD1	2:O:140:GLU:HG2	2.02	0.60
1:A:479:LYS:CG	1:A:488:LEU:HD21	2.30	0.60
1:A:710:HIS:O	1:A:712:PHE:N	2.35	0.60
1:A:752:LEU:O	1:A:756:ILE:HG12	2.02	0.60
1:B:79:ILE:C	1:B:81:GLN:N	2.60	0.60
1:B:349:ASN:ND2	1:B:350:VAL:HG23	2.17	0.60
1:B:517:VAL:HG23	1:B:518:ASN:ND2	2.17	0.60
1:B:759:GLN:HA	1:B:759:GLN:HE21	1.67	0.60
1:C:105:TYR:N	1:C:152:LEU:O	2.31	0.60
1:D:97:TYR:OH	1:D:150:PRO:HB2	2.02	0.60
1:D:122:GLU:HG2	1:D:122:GLU:O	2.00	0.60
1:D:275:GLY:CA	1:D:278:LYS:HG3	2.31	0.60
1:D:298:GLY:O	1:D:300:LYS:N	2.35	0.60
1:E:165:GLN:C	1:E:167:LYS:N	2.59	0.60
1:F:179:LEU:HG	1:F:180:ASP:N	2.17	0.60
2:Q:11:GLU:C	2:Q:13:LYS:H	2.08	0.60
2:R:11:GLU:C	2:R:13:LYS:H	2.09	0.60
1:A:112:VAL:C	1:A:114:HIS:H	2.09	0.59
1:B:332:ASN:HD21	1:B:334:LEU:HB2	1.66	0.59
1:B:522:SER:O	2:P:124:MET:HE2	2.02	0.59
1:C:88:LYS:NZ	1:C:172:GLU:OE2	2.34	0.59
1:C:122:GLU:O	1:C:122:GLU:HG2	2.00	0.59
1:C:446:ILE:HG13	1:C:452:GLU:O	2.00	0.59
1:C:469:PHE:O	1:C:469:PHE:HD1	1.84	0.59
1:C:597:ASN:ND2	1:C:601:GLU:HB2	2.17	0.59
1:D:122:GLU:CD	1:D:147:ARG:HB2	2.26	0.59
1:E:349:ASN:ND2	1:E:350:VAL:HG23	2.16	0.59
1:E:469:PHE:O	1:E:469:PHE:HD1	1.85	0.59
1:F:109:ILE:C	1:F:111:LEU:H	2.10	0.59
1:F:368:GLN:C	1:F:370:LEU:H	2.10	0.59
1:F:529:VAL:HG21	2:T:109:MET:HE1	1.83	0.59
2:O:49:GLN:N	2:O:49:GLN:NE2	2.46	0.59
2:P:138:TYR:O	2:P:142:VAL:HG23	2.02	0.59
2:S:11:GLU:C	2:S:13:LYS:H	2.09	0.59
2:S:97:ASN:ND2	2:S:99:TYR:H	1.99	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:T:49:GLN:O	2:T:51:MET:N	2.35	0.59
1:A:79:ILE:C	1:A:81:GLN:N	2.60	0.59
1:A:710:HIS:C	1:A:712:PHE:N	2.59	0.59
1:B:275:GLY:CA	1:B:278:LYS:HG3	2.31	0.59
1:B:597:ASN:HD22	1:B:601:GLU:HB2	1.67	0.59
1:B:710:HIS:O	1:B:712:PHE:N	2.35	0.59
1:C:275:GLY:O	1:C:278:LYS:HB2	2.01	0.59
1:C:298:GLY:O	1:C:300:LYS:N	2.35	0.59
1:D:172:GLU:HB3	1:D:246:SER:HA	1.83	0.59
1:D:671:ARG:HD2	2:R:14:GLU:CG	2.32	0.59
1:E:79:ILE:C	1:E:81:GLN:N	2.61	0.59
1:E:179:LEU:HG	1:E:180:ASP:H	1.67	0.59
1:F:112:VAL:C	1:F:114:HIS:H	2.10	0.59
2:O:83:GLU:O	2:O:87:GLU:HG3	2.02	0.59
2:R:13:LYS:C	2:R:15:ALA:N	2.56	0.59
1:A:697:ILE:C	1:A:699:GLY:H	2.08	0.59
1:B:279:ILE:O	1:B:283:LEU:HD13	2.02	0.59
1:B:401:ILE:HD13	1:B:485:LEU:O	2.03	0.59
1:C:308:VAL:HB	1:C:311:HIS:ND1	2.17	0.59
1:D:112:VAL:C	1:D:114:HIS:H	2.09	0.59
1:D:334:LEU:HD12	1:D:334:LEU:N	2.17	0.59
1:D:387:ASN:N	1:D:387:ASN:ND2	2.49	0.59
1:E:122:GLU:CD	1:E:147:ARG:HB2	2.26	0.59
1:E:129:ASN:HD22	1:E:129:ASN:N	1.97	0.59
1:E:308:VAL:HB	1:E:311:HIS:ND1	2.18	0.59
1:E:517:VAL:HG23	1:E:518:ASN:ND2	2.17	0.59
1:E:622:LYS:O	1:E:623:ASP:HB2	2.02	0.59
1:F:141:PHE:HD1	1:F:141:PHE:N	1.99	0.59
2:P:37:ARG:HG2	2:P:37:ARG:HH11	1.67	0.59
2:R:49:GLN:O	2:R:51:MET:N	2.35	0.59
2:T:138:TYR:O	2:T:142:VAL:HG23	2.02	0.59
1:A:165:GLN:C	1:A:167:LYS:N	2.58	0.59
1:A:298:GLY:O	1:A:300:LYS:N	2.35	0.59
1:B:298:GLY:O	1:B:300:LYS:N	2.36	0.59
1:B:671:ARG:HD2	2:P:14:GLU:CG	2.32	0.59
1:C:343:VAL:HG13	1:C:487:PRO:HG2	1.83	0.59
1:D:288:VAL:HG23	1:D:289:GLU:H	1.66	0.59
1:D:752:LEU:O	1:D:756:ILE:HG12	2.02	0.59
1:E:159:TYR:N	1:E:159:TYR:CD1	2.70	0.59
1:E:279:ILE:O	1:E:283:LEU:HD13	2.03	0.59
1:E:401:ILE:HD13	1:E:485:LEU:O	2.01	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:93:VAL:HA	1:F:96:ILE:CD1	2.31	0.59
2:O:49:GLN:O	2:O:51:MET:N	2.35	0.59
2:S:37:ARG:HH11	2:S:37:ARG:HG2	1.67	0.59
1:A:279:ILE:O	1:A:283:LEU:HD13	2.03	0.59
1:A:318:ILE:HD12	1:A:318:ILE:N	2.18	0.59
1:A:515:LYS:HG2	1:A:515:LYS:O	2.03	0.59
1:B:173:ILE:HG23	1:B:174:GLY:H	1.67	0.59
1:B:308:VAL:HB	1:B:311:HIS:ND1	2.18	0.59
1:B:736:LEU:HD21	1:B:750:GLN:NE2	2.17	0.59
1:C:112:VAL:C	1:C:114:HIS:H	2.10	0.59
1:C:217:LYS:HB3	1:C:217:LYS:NZ	2.14	0.59
1:C:279:ILE:HG22	1:C:283:LEU:HD13	1.85	0.59
1:E:173:ILE:HG23	1:E:174:GLY:H	1.66	0.59
1:E:217:LYS:CB	1:E:236:GLU:HG3	2.32	0.59
1:E:257:LEU:O	1:E:265:PHE:HB2	2.02	0.59
1:E:550:SER:HG	1:E:552:TRP:CD1	2.21	0.59
1:F:180:ASP:CG	1:F:181:ILE:N	2.59	0.59
1:F:190:PRO:O	1:F:195:LEU:HB2	2.03	0.59
1:F:343:VAL:HG13	1:F:487:PRO:HG2	1.83	0.59
2:Q:49:GLN:O	2:Q:51:MET:N	2.35	0.59
2:S:36:MET:HE1	2:S:51:MET:HE1	1.85	0.59
1:B:165:GLN:C	1:B:167:LYS:N	2.58	0.59
1:B:217:LYS:CB	1:B:236:GLU:HG3	2.32	0.59
1:B:469:PHE:HD1	1:B:469:PHE:O	1.85	0.59
1:B:622:LYS:O	1:B:623:ASP:HB2	2.02	0.59
1:C:710:HIS:O	1:C:712:PHE:N	2.35	0.59
1:E:109:ILE:C	1:E:111:LEU:H	2.10	0.59
1:E:165:GLN:HG2	1:E:251:PRO:HG2	1.85	0.59
1:E:671:ARG:HD2	2:S:14:GLU:CG	2.32	0.59
1:F:275:GLY:CA	1:F:278:LYS:HG3	2.30	0.59
1:F:469:PHE:O	1:F:469:PHE:HD1	1.85	0.59
1:B:493:ASP:OD2	1:B:577:HIS:CE1	2.55	0.59
1:B:581:GLN:HE21	1:B:629:ASN:N	1.99	0.59
1:B:752:LEU:O	1:B:756:ILE:HG12	2.03	0.59
1:C:93:VAL:HA	1:C:96:ILE:CD1	2.30	0.59
1:C:172:GLU:HB3	1:C:246:SER:HA	1.83	0.59
1:C:268:MET:O	1:C:271:LEU:HB2	2.03	0.59
1:C:517:VAL:HG23	1:C:518:ASN:ND2	2.17	0.59
1:D:529:VAL:HG21	2:R:109:MET:HE1	1.85	0.59
1:E:752:LEU:O	1:E:756:ILE:HG12	2.03	0.59
1:F:597:ASN:OD1	1:F:599:GLU:N	2.31	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:O:37:ARG:HG2	2:O:37:ARG:HH11	1.68	0.59
2:P:97:ASN:ND2	2:P:99:TYR:H	2.00	0.59
2:Q:137:ASN:OD1	2:Q:140:GLU:HG2	2.03	0.59
1:A:469:PHE:O	1:A:469:PHE:HD1	1.85	0.59
1:B:293:ILE:O	1:B:295:VAL:HG13	2.03	0.59
1:B:609:GLU:OE2	1:B:609:GLU:N	2.31	0.59
1:C:176:GLY:C	1:C:178:SER:H	2.10	0.59
1:C:318:ILE:HG23	1:C:322:LEU:HD12	1.82	0.59
1:E:318:ILE:HD12	1:E:318:ILE:N	2.17	0.59
1:F:401:ILE:HD13	1:F:485:LEU:O	2.02	0.59
1:A:217:LYS:CB	1:A:236:GLU:HG3	2.32	0.59
1:B:529:VAL:HG21	2:P:109:MET:HE1	1.85	0.59
1:C:165:GLN:C	1:C:167:LYS:N	2.59	0.59
1:C:493:ASP:OD2	1:C:577:HIS:CE1	2.56	0.59
1:C:522:SER:O	2:Q:124:MET:HE2	2.03	0.59
1:C:671:ARG:HD2	2:Q:14:GLU:CG	2.33	0.59
1:D:217:LYS:CB	1:D:236:GLU:HG3	2.32	0.59
1:D:332:ASN:HD21	1:D:334:LEU:HB2	1.67	0.59
1:D:401:ILE:HD13	1:D:485:LEU:O	2.02	0.59
1:D:424:LYS:HZ2	1:D:424:LYS:HB3	1.66	0.59
1:D:515:LYS:HG2	1:D:515:LYS:O	2.03	0.59
1:E:238:GLN:C	1:E:240:ALA:N	2.58	0.59
1:F:217:LYS:CB	1:F:236:GLU:HG3	2.31	0.59
2:P:13:LYS:C	2:P:15:ALA:N	2.57	0.59
2:S:49:GLN:O	2:S:51:MET:N	2.35	0.59
1:A:190:PRO:O	1:A:195:LEU:HB2	2.03	0.59
1:B:141:PHE:HD1	1:B:141:PHE:N	1.99	0.59
1:D:88:LYS:NZ	1:D:172:GLU:OE2	2.34	0.59
1:D:190:PRO:O	1:D:195:LEU:HB2	2.03	0.59
1:D:356:ASP:O	1:D:357:TRP:HB3	2.03	0.59
1:D:724:ARG:HG3	1:D:724:ARG:HH11	1.68	0.59
1:E:176:GLY:C	1:E:178:SER:H	2.11	0.59
1:F:173:ILE:HG23	1:F:174:GLY:H	1.68	0.59
1:F:268:MET:O	1:F:271:LEU:HB2	2.02	0.59
1:F:293:ILE:O	1:F:295:VAL:HG13	2.03	0.59
1:F:308:VAL:HB	1:F:311:HIS:ND1	2.17	0.59
1:F:318:ILE:HD12	1:F:318:ILE:N	2.17	0.59
2:R:94:LYS:HE3	2:R:107:HIS:HD2	1.66	0.59
2:T:11:GLU:C	2:T:13:LYS:H	2.08	0.59
2:T:124:MET:O	2:T:125:ILE:C	2.46	0.59
1:A:173:ILE:HG23	1:A:174:GLY:H	1.67	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:180:ASP:CG	1:A:181:ILE:N	2.59	0.58
1:A:257:LEU:O	1:A:265:PHE:HB2	2.03	0.58
1:A:275:GLY:O	1:A:278:LYS:HB2	2.02	0.58
1:B:165:GLN:HG2	1:B:251:PRO:HG2	1.85	0.58
1:B:202:ASP:HB2	1:B:208:LEU:HD23	1.85	0.58
1:D:622:LYS:O	1:D:623:ASP:HB2	2.02	0.58
1:F:297:LYS:NZ	1:F:297:LYS:HB3	2.18	0.58
1:A:192:PHE:HA	1:A:195:LEU:HB3	1.85	0.58
1:A:279:ILE:HG22	1:A:283:LEU:HD13	1.85	0.58
1:A:288:VAL:HG23	1:A:289:GLU:H	1.66	0.58
1:B:190:PRO:O	1:B:195:LEU:HB2	2.03	0.58
1:C:190:PRO:O	1:C:195:LEU:HB2	2.03	0.58
1:C:297:LYS:HZ3	1:C:601:GLU:HB3	1.67	0.58
1:F:597:ASN:ND2	1:F:601:GLU:H	2.01	0.58
1:F:622:LYS:O	1:F:623:ASP:HB2	2.02	0.58
1:F:697:ILE:C	1:F:699:GLY:H	2.08	0.58
2:R:138:TYR:O	2:R:142:VAL:HG23	2.03	0.58
1:A:109:ILE:C	1:A:111:LEU:H	2.10	0.58
1:A:268:MET:O	1:A:271:LEU:HB2	2.03	0.58
1:A:457:THR:HG23	1:A:469:PHE:H	1.68	0.58
1:C:94:LEU:O	1:C:97:TYR:N	2.37	0.58
1:C:413:LEU:HB2	1:C:419:ILE:CG1	2.33	0.58
1:D:202:ASP:HB2	1:D:208:LEU:HD23	1.86	0.58
1:D:268:MET:O	1:D:271:LEU:HB2	2.03	0.58
1:D:469:PHE:O	1:D:469:PHE:HD1	1.85	0.58
1:E:217:LYS:HB3	1:E:217:LYS:NZ	2.14	0.58
1:E:334:LEU:HD12	1:E:334:LEU:N	2.18	0.58
1:E:349:ASN:HD22	1:E:350:VAL:HG23	1.68	0.58
1:F:257:LEU:O	1:F:265:PHE:HB2	2.03	0.58
1:F:349:ASN:HD22	1:F:350:VAL:HG23	1.68	0.58
1:F:356:ASP:O	1:F:357:TRP:HB3	2.03	0.58
1:F:517:VAL:HG23	1:F:518:ASN:ND2	2.18	0.58
1:A:308:VAL:HB	1:A:311:HIS:ND1	2.18	0.58
1:B:179:LEU:HG	1:B:180:ASP:H	1.68	0.58
1:B:457:THR:HG23	1:B:469:PHE:H	1.68	0.58
1:C:368:GLN:C	1:C:370:LEU:H	2.11	0.58
1:C:401:ILE:HD13	1:C:485:LEU:O	2.03	0.58
1:C:724:ARG:HH11	1:C:724:ARG:HG3	1.68	0.58
1:D:349:ASN:ND2	1:D:350:VAL:HG23	2.17	0.58
1:D:736:LEU:HD21	1:D:750:GLN:NE2	2.18	0.58
1:E:368:GLN:C	1:E:370:LEU:H	2.11	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:179:LEU:HG	1:F:180:ASP:H	1.68	0.58
1:F:581:GLN:HE21	1:F:629:ASN:N	2.01	0.58
1:A:522:SER:O	2:O:124:MET:HE2	2.03	0.58
1:A:736:LEU:HD21	1:A:750:GLN:NE2	2.18	0.58
1:B:76:LEU:O	1:B:79:ILE:N	2.36	0.58
1:C:192:PHE:HA	1:C:195:LEU:HB3	1.86	0.58
1:C:457:THR:HG23	1:C:469:PHE:H	1.68	0.58
1:D:71:PHE:O	1:D:78:LYS:NZ	2.37	0.58
1:D:165:GLN:HG2	1:D:251:PRO:HG2	1.85	0.58
1:D:318:ILE:HD12	1:D:318:ILE:N	2.17	0.58
1:F:109:ILE:HD13	1:F:157:LYS:NZ	2.18	0.58
1:F:165:GLN:HG2	1:F:251:PRO:HG2	1.86	0.58
1:F:217:LYS:HB3	1:F:217:LYS:NZ	2.13	0.58
1:F:238:GLN:C	1:F:240:ALA:N	2.59	0.58
1:F:279:ILE:O	1:F:283:LEU:HD13	2.03	0.58
1:F:373:LYS:HD3	1:F:376:GLN:NE2	2.19	0.58
1:F:522:SER:O	2:T:124:MET:HE2	2.03	0.58
2:Q:12:PHE:O	2:Q:13:LYS:HD2	2.04	0.58
1:A:71:PHE:O	1:A:78:LYS:NZ	2.37	0.58
1:A:99:GLU:C	1:A:101:GLY:N	2.61	0.58
1:A:159:TYR:CD1	1:A:159:TYR:N	2.70	0.58
1:A:187:SER:O	1:A:188:LEU:O	2.22	0.58
1:C:188:LEU:CD2	1:C:188:LEU:N	2.47	0.58
1:C:257:LEU:O	1:C:265:PHE:HB2	2.02	0.58
1:C:297:LYS:NZ	1:C:297:LYS:HB3	2.19	0.58
1:D:93:VAL:HA	1:D:96:ILE:CD1	2.30	0.58
1:D:441:VAL:HG22	1:D:461:LYS:CG	2.31	0.58
1:F:495:PHE:CD1	1:F:495:PHE:C	2.81	0.58
2:O:48:LEU:CA	2:O:51:MET:HE2	2.33	0.58
2:Q:137:ASN:OD1	2:Q:139:GLU:HB2	2.04	0.58
1:A:115:LYS:C	1:A:117:LEU:N	2.58	0.58
1:A:176:GLY:C	1:A:178:SER:H	2.11	0.58
1:A:179:LEU:HG	1:A:180:ASP:H	1.69	0.58
1:A:202:ASP:HB2	1:A:208:LEU:HD23	1.84	0.58
1:B:109:ILE:C	1:B:111:LEU:H	2.10	0.58
1:B:257:LEU:O	1:B:265:PHE:HB2	2.04	0.58
1:B:373:LYS:HD3	1:B:376:GLN:NE2	2.19	0.58
1:C:115:LYS:HG3	1:C:153:ILE:HG21	1.85	0.58
1:C:356:ASP:O	1:C:357:TRP:HB3	2.02	0.58
1:D:176:GLY:C	1:D:178:SER:H	2.11	0.58
1:D:186:LYS:O	1:D:188:LEU:O	2.21	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:293:ILE:O	1:D:295:VAL:HG13	2.03	0.58
1:D:349:ASN:HD22	1:D:350:VAL:HG23	1.69	0.58
1:E:515:LYS:O	1:E:515:LYS:HG2	2.02	0.58
1:F:94:LEU:O	1:F:97:TYR:N	2.36	0.58
1:F:187:SER:O	1:F:188:LEU:O	2.22	0.58
1:F:279:ILE:HG22	1:F:283:LEU:HD13	1.85	0.58
2:T:83:GLU:O	2:T:87:GLU:HG3	2.03	0.58
1:A:105:TYR:N	1:A:152:LEU:O	2.31	0.58
1:A:165:GLN:HG2	1:A:251:PRO:HG2	1.85	0.58
1:A:724:ARG:HG3	1:A:724:ARG:HH11	1.69	0.58
1:A:777:TYR:O	1:A:778:LYS:C	2.46	0.58
1:B:268:MET:O	1:B:271:LEU:HB2	2.03	0.58
1:B:349:ASN:HD22	1:B:350:VAL:H	1.52	0.58
1:C:202:ASP:HB2	1:C:208:LEU:HD23	1.85	0.58
1:C:334:LEU:HD12	1:C:334:LEU:N	2.19	0.58
1:D:581:GLN:HE21	1:D:629:ASN:N	2.01	0.58
1:E:71:PHE:O	1:E:78:LYS:NZ	2.37	0.58
1:E:597:ASN:OD1	1:E:599:GLU:N	2.31	0.58
1:F:192:PHE:HA	1:F:195:LEU:HB3	1.84	0.58
1:F:196:ILE:O	1:F:199:LEU:HB2	2.04	0.58
1:F:333:LYS:H	1:F:333:LYS:HD2	1.69	0.58
1:F:493:ASP:OD2	1:F:577:HIS:CE1	2.57	0.58
1:F:671:ARG:HD2	2:T:14:GLU:CG	2.33	0.58
2:P:124:MET:O	2:P:125:ILE:C	2.47	0.58
2:S:48:LEU:CA	2:S:51:MET:HE2	2.33	0.58
1:A:499:PRO:HG2	1:A:504:ILE:HD11	1.86	0.58
1:B:196:ILE:O	1:B:199:LEU:HB2	2.04	0.58
1:B:515:LYS:O	1:B:515:LYS:HG2	2.03	0.58
1:C:515:LYS:O	1:C:515:LYS:HG2	2.02	0.58
1:D:279:ILE:O	1:D:283:LEU:HD13	2.03	0.58
1:D:349:ASN:HD22	1:D:350:VAL:H	1.52	0.58
1:E:173:ILE:HG23	1:E:174:GLY:N	2.19	0.58
1:E:187:SER:O	1:E:188:LEU:O	2.22	0.58
1:E:710:HIS:O	1:E:712:PHE:N	2.36	0.58
1:E:777:TYR:O	1:E:778:LYS:C	2.47	0.58
1:F:186:LYS:O	1:F:188:LEU:O	2.22	0.58
1:F:441:VAL:HG22	1:F:461:LYS:CG	2.32	0.58
2:Q:55:VAL:CG2	2:Q:67:GLU:OE1	2.52	0.58
2:S:105:LEU:HD21	2:S:124:MET:SD	2.44	0.58
1:B:279:ILE:HG22	1:B:283:LEU:HD13	1.85	0.58
1:B:333:LYS:H	1:B:333:LYS:HD2	1.69	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:106:PHE:HA	1:D:154:ILE:O	2.04	0.58
1:D:279:ILE:HG22	1:D:283:LEU:HD13	1.86	0.58
1:E:493:ASP:OD2	1:E:577:HIS:CE1	2.57	0.58
1:F:88:LYS:HZ3	1:F:172:GLU:CD	2.12	0.58
1:F:752:LEU:O	1:F:756:ILE:HG12	2.04	0.58
1:F:777:TYR:O	1:F:778:LYS:C	2.47	0.58
2:Q:97:ASN:ND2	2:Q:99:TYR:H	2.02	0.58
1:A:293:ILE:O	1:A:295:VAL:HG13	2.04	0.57
1:A:373:LYS:HD3	1:A:376:GLN:NE2	2.19	0.57
1:B:159:TYR:N	1:B:159:TYR:CD1	2.70	0.57
1:B:179:LEU:H	1:B:179:LEU:CD2	2.16	0.57
1:B:187:SER:O	1:B:188:LEU:O	2.22	0.57
1:C:349:ASN:ND2	1:C:350:VAL:HG23	2.18	0.57
1:C:609:GLU:OE2	1:C:609:GLU:N	2.30	0.57
1:C:752:LEU:O	1:C:756:ILE:HG12	2.03	0.57
1:D:109:ILE:C	1:D:111:LEU:H	2.10	0.57
1:D:159:TYR:N	1:D:159:TYR:CD1	2.70	0.57
1:E:297:LYS:NZ	1:E:297:LYS:HB3	2.19	0.57
1:E:499:PRO:HG2	1:E:504:ILE:HD11	1.85	0.57
1:F:165:GLN:C	1:F:167:LYS:N	2.60	0.57
1:F:334:LEU:HD12	1:F:334:LEU:N	2.18	0.57
2:R:55:VAL:CG2	2:R:67:GLU:OE1	2.52	0.57
1:A:413:LEU:HB2	1:A:419:ILE:CG1	2.33	0.57
1:A:581:GLN:HE21	1:A:629:ASN:N	1.99	0.57
1:B:176:GLY:C	1:B:178:SER:H	2.10	0.57
1:C:318:ILE:HD12	1:C:318:ILE:N	2.17	0.57
1:C:441:VAL:HG22	1:C:461:LYS:CG	2.32	0.57
1:D:297:LYS:NZ	1:D:297:LYS:HB3	2.19	0.57
1:D:413:LEU:HB2	1:D:419:ILE:CG1	2.34	0.57
1:E:190:PRO:O	1:E:195:LEU:HB2	2.03	0.57
2:P:48:LEU:CA	2:P:51:MET:HE2	2.34	0.57
2:Q:105:LEU:HD21	2:Q:124:MET:SD	2.43	0.57
1:A:517:VAL:HG23	1:A:518:ASN:ND2	2.18	0.57
1:A:657:ILE:HD11	1:A:701:LEU:HD23	1.86	0.57
1:B:186:LYS:O	1:B:188:LEU:O	2.22	0.57
1:B:318:ILE:HD12	1:B:318:ILE:N	2.17	0.57
1:D:457:THR:HG23	1:D:469:PHE:H	1.68	0.57
1:E:76:LEU:O	1:E:79:ILE:N	2.37	0.57
1:E:186:LYS:O	1:E:188:LEU:O	2.22	0.57
1:E:373:LYS:HD3	1:E:376:GLN:NE2	2.20	0.57
1:E:597:ASN:ND2	1:E:601:GLU:HB2	2.19	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:609:GLU:OE2	1:E:609:GLU:N	2.32	0.57
2:P:55:VAL:CG2	2:P:67:GLU:OE1	2.52	0.57
2:R:12:PHE:O	2:R:13:LYS:HD2	2.04	0.57
1:A:597:ASN:ND2	1:A:601:GLU:HB2	2.18	0.57
1:B:71:PHE:O	1:B:78:LYS:NZ	2.38	0.57
1:B:115:LYS:HG3	1:B:153:ILE:HG21	1.86	0.57
1:B:192:PHE:HA	1:B:195:LEU:HB3	1.86	0.57
1:B:597:ASN:OD1	1:B:599:GLU:N	2.31	0.57
1:B:639:ASN:ND2	1:B:639:ASN:H	2.03	0.57
1:B:697:ILE:HG21	1:B:732:ILE:HD11	1.86	0.57
1:B:732:ILE:HG23	1:B:749:PHE:HD1	1.70	0.57
1:C:499:PRO:HG2	1:C:504:ILE:HD11	1.86	0.57
1:C:694:VAL:HG23	2:Q:18:LEU:HD21	1.86	0.57
1:D:94:LEU:O	1:D:97:TYR:N	2.37	0.57
1:D:179:LEU:HG	1:D:180:ASP:H	1.68	0.57
1:D:187:SER:O	1:D:188:LEU:O	2.22	0.57
1:D:308:VAL:HB	1:D:311:HIS:ND1	2.18	0.57
1:E:268:MET:O	1:E:271:LEU:HB2	2.03	0.57
1:E:305:SER:C	1:E:307:LEU:H	2.12	0.57
1:F:175:LYS:HB2	1:F:175:LYS:NZ	2.19	0.57
1:F:664:ILE:HG21	2:T:15:ALA:HB2	1.87	0.57
1:F:714:GLN:O	1:F:718:ARG:HG3	2.04	0.57
2:P:105:LEU:HD21	2:P:124:MET:SD	2.45	0.57
2:S:48:LEU:HD12	2:S:51:MET:HE1	1.87	0.57
2:S:97:ASN:HD22	2:S:97:ASN:C	2.13	0.57
2:S:138:TYR:O	2:S:142:VAL:HG23	2.04	0.57
1:B:297:LYS:NZ	1:B:297:LYS:HB3	2.19	0.57
1:B:368:GLN:C	1:B:370:LEU:H	2.11	0.57
1:B:710:HIS:C	1:B:712:PHE:N	2.59	0.57
1:C:187:SER:O	1:C:188:LEU:O	2.22	0.57
1:C:521:ASN:HB3	1:C:524:GLU:HB3	1.86	0.57
1:C:639:ASN:ND2	1:C:639:ASN:H	2.03	0.57
1:C:777:TYR:O	1:C:778:LYS:C	2.48	0.57
1:E:333:LYS:H	1:E:333:LYS:HD2	1.69	0.57
1:F:80:GLN:HG3	1:F:80:GLN:O	2.05	0.57
1:F:199:LEU:O	1:F:201:ASP:N	2.32	0.57
2:Q:83:GLU:O	2:Q:87:GLU:HG3	2.05	0.57
2:R:105:LEU:HD21	2:R:124:MET:SD	2.45	0.57
2:S:12:PHE:HE1	2:S:72:MET:CE	2.13	0.57
2:S:55:VAL:CG2	2:S:67:GLU:OE1	2.52	0.57
2:T:97:ASN:ND2	2:T:99:TYR:H	2.02	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:609:GLU:OE2	1:A:609:GLU:N	2.30	0.57
1:B:238:GLN:C	1:B:240:ALA:N	2.59	0.57
1:B:356:ASP:O	1:B:357:TRP:HB3	2.03	0.57
1:B:424:LYS:HB3	1:B:424:LYS:HZ2	1.68	0.57
1:B:519:THR:OG1	1:B:520:PRO:HD2	2.05	0.57
1:C:173:ILE:HG23	1:C:174:GLY:N	2.20	0.57
1:C:622:LYS:O	1:C:623:ASP:HB2	2.03	0.57
1:E:275:GLY:HA2	1:E:278:LYS:HE2	1.86	0.57
1:E:281:GLU:O	1:E:285:LYS:HG2	2.05	0.57
1:F:730:ASN:O	1:F:733:GLU:N	2.38	0.57
2:R:48:LEU:CA	2:R:51:MET:HE2	2.33	0.57
2:S:94:LYS:HB3	2:S:94:LYS:HZ2	1.68	0.57
1:A:199:LEU:O	1:A:201:ASP:N	2.33	0.57
1:A:297:LYS:NZ	1:A:297:LYS:HB3	2.18	0.57
1:A:387:ASN:N	1:A:387:ASN:ND2	2.50	0.57
1:A:403:LEU:HG	1:A:405:LEU:CD1	2.35	0.57
1:A:521:ASN:HB3	1:A:524:GLU:HB3	1.86	0.57
1:B:499:PRO:HG2	1:B:504:ILE:HD11	1.86	0.57
1:B:764:LEU:C	1:B:766:HIS:H	2.11	0.57
1:C:664:ILE:HG21	2:Q:15:ALA:HB2	1.86	0.57
1:D:99:GLU:C	1:D:101:GLY:N	2.60	0.57
1:D:368:GLN:C	1:D:370:LEU:H	2.11	0.57
1:E:279:ILE:HG22	1:E:283:LEU:HD13	1.86	0.57
1:E:714:GLN:O	1:E:718:ARG:HG3	2.05	0.57
1:F:97:TYR:CZ	1:F:150:PRO:HB2	2.40	0.57
2:P:12:PHE:O	2:P:13:LYS:HD2	2.05	0.57
1:A:115:LYS:HB2	1:A:118:GLN:HG2	1.86	0.57
1:A:519:THR:OG1	1:A:520:PRO:HD2	2.05	0.57
1:A:622:LYS:O	1:A:623:ASP:HB2	2.03	0.57
1:B:180:ASP:CG	1:B:181:ILE:N	2.59	0.57
1:B:479:LYS:CG	1:B:488:LEU:HD21	2.30	0.57
1:B:694:VAL:HG23	2:P:18:LEU:HD21	1.86	0.57
1:C:71:PHE:CG	1:C:73:ASN:HB2	2.40	0.57
1:C:99:GLU:C	1:C:101:GLY:N	2.61	0.57
1:C:179:LEU:HG	1:C:180:ASP:H	1.67	0.57
1:C:333:LYS:H	1:C:333:LYS:HD2	1.69	0.57
1:C:349:ASN:HD22	1:C:350:VAL:H	1.52	0.57
1:C:732:ILE:HG23	1:C:749:PHE:HD1	1.69	0.57
1:D:115:LYS:HG3	1:D:153:ILE:HG21	1.86	0.57
1:D:188:LEU:CD2	1:D:188:LEU:N	2.47	0.57
1:D:192:PHE:HA	1:D:195:LEU:HB3	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:333:LYS:HD2	1:D:333:LYS:H	1.68	0.57
1:D:732:ILE:HG23	1:D:749:PHE:HD1	1.69	0.57
1:E:196:ILE:O	1:E:199:LEU:HB2	2.04	0.57
1:E:202:ASP:HB2	1:E:208:LEU:HD23	1.86	0.57
2:T:36:MET:HE1	2:T:51:MET:HE1	1.86	0.57
1:B:205:SER:C	1:B:207:ASP:H	2.13	0.57
1:B:697:ILE:HD13	1:B:732:ILE:CD1	2.35	0.57
1:B:724:ARG:HH11	1:B:724:ARG:HG3	1.70	0.57
1:C:714:GLN:O	1:C:718:ARG:HG3	2.04	0.57
1:D:179:LEU:HD23	1:D:179:LEU:N	2.18	0.57
1:D:305:SER:C	1:D:307:LEU:H	2.12	0.57
1:D:671:ARG:HH12	1:D:677:GLY:HA3	1.70	0.57
1:E:94:LEU:O	1:E:97:TYR:N	2.38	0.57
1:E:115:LYS:HG3	1:E:153:ILE:HG21	1.86	0.57
1:E:639:ASN:H	1:E:639:ASN:ND2	2.03	0.57
1:F:205:SER:C	1:F:207:ASP:H	2.12	0.57
1:F:326:ILE:C	1:F:327:LEU:HD12	2.30	0.57
2:R:109:MET:HG3	2:R:116:LEU:CD1	2.34	0.57
1:A:441:VAL:HG22	1:A:461:LYS:CG	2.31	0.57
1:A:639:ASN:ND2	1:A:639:ASN:H	2.03	0.57
1:C:71:PHE:O	1:C:78:LYS:NZ	2.38	0.57
1:C:667:LEU:O	1:C:668:SER:C	2.48	0.57
1:D:173:ILE:HG23	1:D:174:GLY:H	1.68	0.57
1:D:205:SER:C	1:D:207:ASP:H	2.12	0.57
1:D:493:ASP:OD2	1:D:577:HIS:CE1	2.58	0.57
1:D:639:ASN:ND2	1:D:639:ASN:H	2.03	0.57
1:F:71:PHE:O	1:F:78:LYS:NZ	2.38	0.57
1:F:305:SER:C	1:F:307:LEU:H	2.13	0.57
1:F:597:ASN:ND2	1:F:601:GLU:HB2	2.19	0.57
1:F:625:LEU:HD12	1:F:626:TYR:N	2.19	0.57
2:T:48:LEU:HD12	2:T:51:MET:HE1	1.87	0.57
1:A:196:ILE:O	1:A:199:LEU:HB2	2.05	0.56
1:A:275:GLY:HA2	1:A:278:LYS:HE2	1.86	0.56
1:A:368:GLN:C	1:A:370:LEU:H	2.11	0.56
1:B:76:LEU:O	1:B:80:GLN:N	2.38	0.56
1:B:173:ILE:HG23	1:B:174:GLY:N	2.20	0.56
1:B:597:ASN:ND2	1:B:601:GLU:HB2	2.20	0.56
1:C:165:GLN:HG2	1:C:251:PRO:HG2	1.86	0.56
1:D:360:VAL:CG1	1:D:370:LEU:HD22	2.34	0.56
1:D:597:ASN:CB	1:D:598:PRO:HD2	2.33	0.56
1:D:671:ARG:NH1	1:D:677:GLY:HA3	2.19	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:205:SER:C	1:E:207:ASP:H	2.13	0.56
1:E:441:VAL:HG22	1:E:461:LYS:CG	2.32	0.56
1:E:457:THR:HG23	1:E:469:PHE:H	1.68	0.56
1:E:764:LEU:C	1:E:766:HIS:H	2.13	0.56
1:F:66:LEU:HD12	1:F:103:GLU:HA	1.87	0.56
1:F:521:ASN:HB3	1:F:524:GLU:HB3	1.87	0.56
2:Q:48:LEU:HD12	2:Q:51:MET:HE1	1.87	0.56
2:Q:109:MET:HG3	2:Q:116:LEU:CD1	2.35	0.56
1:A:106:PHE:HA	1:A:154:ILE:O	2.05	0.56
1:A:173:ILE:HG23	1:A:174:GLY:N	2.20	0.56
1:A:360:VAL:CG1	1:A:370:LEU:HD22	2.34	0.56
1:B:66:LEU:HD12	1:B:103:GLU:HA	1.87	0.56
1:B:207:ASP:HB2	1:B:262:PRO:HG3	1.87	0.56
1:B:413:LEU:HB2	1:B:419:ILE:CG1	2.33	0.56
1:D:175:LYS:HB2	1:D:175:LYS:NZ	2.20	0.56
1:D:196:ILE:O	1:D:199:LEU:HB2	2.05	0.56
1:D:199:LEU:O	1:D:201:ASP:N	2.32	0.56
1:D:257:LEU:O	1:D:265:PHE:HB2	2.04	0.56
1:E:217:LYS:HB3	1:E:236:GLU:OE1	2.05	0.56
1:E:495:PHE:C	1:E:495:PHE:CD1	2.82	0.56
1:E:724:ARG:HH11	1:E:724:ARG:HG3	1.70	0.56
1:F:217:LYS:HB3	1:F:236:GLU:OE1	2.05	0.56
1:F:408:LEU:HD12	1:F:408:LEU:N	2.01	0.56
1:F:499:PRO:HG2	1:F:504:ILE:HD11	1.87	0.56
1:F:639:ASN:ND2	1:F:639:ASN:H	2.03	0.56
1:F:724:ARG:HH11	1:F:724:ARG:HG3	1.69	0.56
1:A:94:LEU:O	1:A:97:TYR:N	2.38	0.56
1:A:205:SER:C	1:A:207:ASP:H	2.13	0.56
1:A:305:SER:C	1:A:307:LEU:H	2.13	0.56
1:A:311:HIS:O	1:A:314:ALA:HB3	2.04	0.56
1:A:356:ASP:O	1:A:357:TRP:HB3	2.04	0.56
1:A:495:PHE:CD1	1:A:495:PHE:C	2.83	0.56
1:B:179:LEU:HD23	1:B:179:LEU:N	2.18	0.56
1:B:275:GLY:HA2	1:B:278:LYS:HE2	1.86	0.56
1:B:403:LEU:HG	1:B:405:LEU:CD1	2.36	0.56
1:C:109:ILE:HD13	1:C:157:LYS:NZ	2.21	0.56
1:C:196:ILE:O	1:C:199:LEU:HB2	2.04	0.56
1:C:205:SER:C	1:C:207:ASP:H	2.13	0.56
1:C:275:GLY:HA2	1:C:278:LYS:HE2	1.86	0.56
1:C:495:PHE:CD1	1:C:495:PHE:C	2.83	0.56
1:D:79:ILE:C	1:D:81:GLN:N	2.60	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:109:ILE:HD13	1:D:157:LYS:NZ	2.20	0.56
1:D:115:LYS:C	1:D:117:LEU:N	2.58	0.56
1:E:179:LEU:H	1:E:179:LEU:CD2	2.17	0.56
1:E:192:PHE:HA	1:E:195:LEU:HB3	1.85	0.56
1:E:207:ASP:HB2	1:E:262:PRO:HG3	1.86	0.56
1:E:334:LEU:H	1:E:334:LEU:CD1	2.18	0.56
1:E:356:ASP:O	1:E:357:TRP:HB3	2.03	0.56
1:F:173:ILE:HG23	1:F:174:GLY:N	2.21	0.56
1:F:403:LEU:HG	1:F:405:LEU:CD1	2.34	0.56
1:F:457:THR:HG23	1:F:469:PHE:H	1.69	0.56
1:F:694:VAL:HG23	2:T:18:LEU:HD11	1.87	0.56
1:F:715:GLU:HA	1:F:718:ARG:HH12	1.70	0.56
2:O:105:LEU:HD21	2:O:124:MET:SD	2.45	0.56
2:O:124:MET:O	2:O:125:ILE:C	2.48	0.56
2:R:12:PHE:HE1	2:R:72:MET:CE	2.14	0.56
1:A:179:LEU:H	1:A:179:LEU:CD2	2.17	0.56
1:A:349:ASN:HD22	1:A:350:VAL:HG23	1.70	0.56
1:B:349:ASN:HD22	1:B:350:VAL:HG23	1.69	0.56
1:C:373:LYS:HD3	1:C:376:GLN:NE2	2.20	0.56
1:C:657:ILE:HD11	1:C:701:LEU:HD23	1.88	0.56
1:D:217:LYS:HB3	1:D:217:LYS:NZ	2.13	0.56
1:D:694:VAL:HG22	2:R:18:LEU:HD21	1.87	0.56
1:E:519:THR:OG1	1:E:520:PRO:HD2	2.05	0.56
1:F:202:ASP:HB2	1:F:208:LEU:HD23	1.85	0.56
1:F:515:LYS:HG2	1:F:515:LYS:O	2.04	0.56
2:S:124:MET:O	2:S:125:ILE:C	2.48	0.56
1:A:207:ASP:HB2	1:A:262:PRO:HG3	1.87	0.56
1:A:597:ASN:CB	1:A:598:PRO:HD2	2.35	0.56
1:A:748:TYR:O	1:A:751:TYR:HB3	2.06	0.56
1:B:80:GLN:O	1:B:80:GLN:HG3	2.06	0.56
1:B:199:LEU:O	1:B:201:ASP:N	2.33	0.56
1:B:521:ASN:HB3	1:B:524:GLU:HB3	1.86	0.56
1:D:165:GLN:C	1:D:167:LYS:N	2.59	0.56
1:D:403:LEU:HG	1:D:405:LEU:CD1	2.35	0.56
1:E:109:ILE:HD13	1:E:157:LYS:NZ	2.20	0.56
1:E:175:LYS:NZ	1:E:175:LYS:HB2	2.20	0.56
1:E:360:VAL:CG1	1:E:370:LEU:HD22	2.33	0.56
1:E:671:ARG:NH1	1:E:677:GLY:HA3	2.21	0.56
1:F:671:ARG:NH1	1:F:677:GLY:HA3	2.20	0.56
1:F:697:ILE:HG21	1:F:732:ILE:HD11	1.87	0.56
2:R:48:LEU:HD12	2:R:51:MET:HE1	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:T:12:PHE:HE1	2:T:72:MET:CE	2.13	0.56
1:A:115:LYS:HG3	1:A:153:ILE:HG21	1.87	0.56
1:A:281:GLU:O	1:A:285:LYS:HG2	2.06	0.56
1:B:664:ILE:HG21	2:P:15:ALA:HB2	1.87	0.56
1:C:106:PHE:HA	1:C:154:ILE:O	2.05	0.56
1:C:281:GLU:O	1:C:285:LYS:HG2	2.06	0.56
1:C:597:ASN:OD1	1:C:599:GLU:N	2.32	0.56
1:D:664:ILE:HG21	2:R:15:ALA:HB2	1.86	0.56
1:E:403:LEU:HG	1:E:405:LEU:CD1	2.36	0.56
2:P:36:MET:HE1	2:P:51:MET:HE1	1.88	0.56
2:R:83:GLU:O	2:R:87:GLU:HG3	2.06	0.56
1:A:365:PRO:HB2	1:A:367:ASP:O	2.06	0.56
1:B:115:LYS:HE3	1:B:116:GLU:N	2.21	0.56
1:B:441:VAL:HG22	1:B:461:LYS:CG	2.32	0.56
1:B:777:TYR:O	1:B:778:LYS:C	2.47	0.56
1:C:88:LYS:HZ3	1:C:172:GLU:CD	2.14	0.56
1:C:349:ASN:HD22	1:C:350:VAL:HG23	1.71	0.56
1:C:697:ILE:HG21	1:C:732:ILE:HD11	1.88	0.56
1:E:494:LEU:HB3	1:E:579:THR:HG22	1.88	0.56
2:O:12:PHE:O	2:O:13:LYS:HD2	2.04	0.56
2:R:124:MET:O	2:R:125:ILE:C	2.47	0.56
1:A:175:LYS:HB2	1:A:175:LYS:NZ	2.21	0.56
1:A:334:LEU:HD12	1:A:334:LEU:N	2.19	0.56
1:A:671:ARG:NH1	1:A:677:GLY:HA3	2.20	0.56
1:A:732:ILE:HG23	1:A:749:PHE:HD1	1.70	0.56
1:B:105:TYR:N	1:B:152:LEU:O	2.31	0.56
1:B:141:PHE:N	1:B:141:PHE:CD1	2.70	0.56
1:C:78:LYS:HG3	1:C:79:ILE:N	2.20	0.56
1:C:639:ASN:H	1:C:639:ASN:HD22	1.53	0.56
1:C:671:ARG:NH1	1:C:677:GLY:HA3	2.21	0.56
1:D:495:PHE:CD1	1:D:495:PHE:C	2.84	0.56
1:E:106:PHE:HA	1:E:154:ILE:O	2.05	0.56
1:E:657:ILE:HD11	1:E:701:LEU:HD23	1.88	0.56
1:E:715:GLU:HA	1:E:718:ARG:HH12	1.69	0.56
1:F:106:PHE:HA	1:F:154:ILE:O	2.05	0.56
1:F:134:LYS:C	1:F:136:PRO:HD3	2.31	0.56
1:F:165:GLN:OE1	1:F:252:ASP:HB3	2.06	0.56
1:F:479:LYS:CG	1:F:488:LEU:HD21	2.31	0.56
1:F:667:LEU:O	1:F:668:SER:C	2.48	0.56
2:O:12:PHE:HE1	2:O:72:MET:CE	2.14	0.56
2:R:16:PHE:CE1	2:R:27:ILE:HG12	2.41	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:T:48:LEU:CA	2:T:51:MET:HE2	2.34	0.56
1:A:333:LYS:H	1:A:333:LYS:HD2	1.69	0.56
1:A:387:ASN:HB3	1:A:477:MET:SD	2.46	0.56
1:B:109:ILE:HD13	1:B:157:LYS:NZ	2.21	0.56
1:B:305:SER:C	1:B:307:LEU:H	2.13	0.56
1:B:334:LEU:HD12	1:B:334:LEU:N	2.20	0.56
1:C:115:LYS:HE3	1:C:116:GLU:N	2.20	0.56
1:C:305:SER:C	1:C:307:LEU:H	2.13	0.56
1:D:80:GLN:HG3	1:D:80:GLN:O	2.06	0.56
1:D:308:VAL:O	1:D:311:HIS:HB2	2.06	0.56
1:D:373:LYS:HD3	1:D:376:GLN:NE2	2.20	0.56
1:E:97:TYR:CZ	1:E:150:PRO:HB2	2.41	0.56
1:F:115:LYS:HG3	1:F:153:ILE:HG21	1.86	0.56
1:F:207:ASP:HB2	1:F:262:PRO:HG3	1.87	0.56
2:T:12:PHE:O	2:T:13:LYS:HD2	2.05	0.56
2:T:55:VAL:CG2	2:T:67:GLU:OE1	2.53	0.56
2:T:94:LYS:HB3	2:T:94:LYS:NZ	2.21	0.56
2:T:105:LEU:HD21	2:T:124:MET:SD	2.46	0.56
2:T:109:MET:HG3	2:T:116:LEU:CD1	2.35	0.56
1:A:80:GLN:HG3	1:A:80:GLN:O	2.07	0.56
1:A:186:LYS:O	1:A:188:LEU:O	2.23	0.56
1:B:115:LYS:HB2	1:B:118:GLN:HG2	1.87	0.56
1:C:175:LYS:HB2	1:C:175:LYS:NZ	2.21	0.56
1:C:207:ASP:HB2	1:C:262:PRO:HG3	1.87	0.56
1:D:179:LEU:H	1:D:179:LEU:CD2	2.17	0.56
1:D:275:GLY:HA2	1:D:278:LYS:HE2	1.86	0.56
1:D:311:HIS:O	1:D:314:ALA:HB3	2.06	0.56
1:D:334:LEU:H	1:D:334:LEU:CD1	2.17	0.56
1:E:179:LEU:HD23	1:E:179:LEU:N	2.18	0.56
1:E:579:THR:C	1:E:581:GLN:H	2.14	0.56
1:F:275:GLY:HA2	1:F:278:LYS:HE2	1.87	0.56
1:F:694:VAL:HG23	2:T:18:LEU:HD21	1.87	0.56
2:O:36:MET:HE1	2:O:51:MET:HE1	1.88	0.56
2:P:94:LYS:NZ	2:P:94:LYS:HB3	2.21	0.56
2:Q:36:MET:HE1	2:Q:51:MET:HE1	1.87	0.56
2:R:97:ASN:C	2:R:97:ASN:HD22	2.12	0.56
1:A:165:GLN:O	1:A:167:LYS:N	2.39	0.55
1:B:94:LEU:O	1:B:97:TYR:N	2.38	0.55
1:B:99:GLU:C	1:B:101:GLY:N	2.61	0.55
1:B:165:GLN:O	1:B:167:LYS:N	2.38	0.55
1:B:494:LEU:HB3	1:B:579:THR:HG22	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:495:PHE:C	1:B:495:PHE:CD1	2.84	0.55
1:B:671:ARG:NH1	1:B:677:GLY:HA3	2.22	0.55
1:B:748:TYR:O	1:B:751:TYR:HB3	2.07	0.55
1:D:115:LYS:HB2	1:D:118:GLN:HG2	1.87	0.55
1:D:173:ILE:HG23	1:D:174:GLY:N	2.21	0.55
1:D:207:ASP:HB2	1:D:262:PRO:HG3	1.87	0.55
1:D:521:ASN:HB3	1:D:524:GLU:HB3	1.86	0.55
1:D:694:VAL:HG23	2:R:18:LEU:HD21	1.86	0.55
1:D:748:TYR:O	1:D:751:TYR:HB3	2.06	0.55
1:D:777:TYR:O	1:D:778:LYS:C	2.47	0.55
1:E:66:LEU:HD12	1:E:103:GLU:HA	1.88	0.55
1:F:90:PRO:O	1:F:92:ASP:N	2.39	0.55
2:O:10:ALA:O	2:O:14:GLU:HB2	2.06	0.55
2:P:48:LEU:HD12	2:P:51:MET:HE1	1.88	0.55
2:R:16:PHE:HE1	2:R:27:ILE:HG12	1.71	0.55
2:S:12:PHE:O	2:S:13:LYS:HD2	2.07	0.55
1:A:97:TYR:CZ	1:A:150:PRO:HB2	2.41	0.55
1:A:667:LEU:O	1:A:668:SER:C	2.50	0.55
1:A:671:ARG:HH12	1:A:677:GLY:HA3	1.71	0.55
1:C:165:GLN:O	1:C:167:LYS:N	2.39	0.55
1:C:405:LEU:HD12	1:C:405:LEU:N	2.22	0.55
1:C:671:ARG:HH12	1:C:677:GLY:HA3	1.71	0.55
1:C:730:ASN:O	1:C:733:GLU:N	2.39	0.55
1:D:64:ASN:HD22	1:D:64:ASN:N	2.05	0.55
1:D:375:GLY:O	1:D:377:GLN:N	2.38	0.55
1:D:694:VAL:HG23	2:R:18:LEU:HD11	1.87	0.55
1:E:115:LYS:HE3	1:E:116:GLU:N	2.21	0.55
1:E:694:VAL:HG23	2:S:18:LEU:HD11	1.88	0.55
1:F:519:THR:OG1	1:F:520:PRO:HD2	2.05	0.55
2:O:48:LEU:HD12	2:O:51:MET:HE1	1.87	0.55
2:O:58:ASP:C	2:O:60:ASN:N	2.64	0.55
1:A:115:LYS:HE3	1:A:116:GLU:N	2.21	0.55
1:A:694:VAL:HG23	2:O:18:LEU:HD21	1.88	0.55
1:C:71:PHE:CD2	1:C:73:ASN:HB2	2.41	0.55
1:C:697:ILE:HD13	1:C:732:ILE:CD1	2.36	0.55
1:D:714:GLN:O	1:D:718:ARG:HG3	2.05	0.55
1:E:141:PHE:N	1:E:141:PHE:CD1	2.69	0.55
1:E:165:GLN:OE1	1:E:252:ASP:HB3	2.06	0.55
1:F:78:LYS:HG3	1:F:79:ILE:N	2.21	0.55
1:F:657:ILE:HD11	1:F:701:LEU:HD23	1.89	0.55
1:F:732:ILE:HG23	1:F:749:PHE:HD1	1.70	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:O:138:TYR:O	2:O:142:VAL:HG23	2.06	0.55
1:A:78:LYS:HG3	1:A:79:ILE:N	2.20	0.55
1:A:134:LYS:C	1:A:136:PRO:HD3	2.31	0.55
1:A:493:ASP:OD2	1:A:577:HIS:CE1	2.58	0.55
1:B:106:PHE:HA	1:B:154:ILE:O	2.06	0.55
1:B:512:GLU:O	1:B:516:VAL:HG23	2.07	0.55
1:B:714:GLN:O	1:B:718:ARG:HG3	2.06	0.55
1:C:186:LYS:O	1:C:188:LEU:O	2.23	0.55
1:C:326:ILE:C	1:C:327:LEU:HD12	2.30	0.55
1:C:403:LEU:HG	1:C:405:LEU:CD1	2.35	0.55
1:D:66:LEU:HD12	1:D:103:GLU:HA	1.87	0.55
1:D:165:GLN:O	1:D:167:LYS:N	2.40	0.55
1:E:413:LEU:HB2	1:E:419:ILE:CG1	2.33	0.55
1:E:512:GLU:O	1:E:516:VAL:HG23	2.07	0.55
1:F:122:GLU:CB	1:F:147:ARG:HG3	2.37	0.55
1:F:375:GLY:O	1:F:377:GLN:N	2.38	0.55
2:Q:13:LYS:HZ3	2:Q:65:PHE:HB3	1.70	0.55
1:A:197:LYS:HE2	1:A:263:ASP:HB3	1.89	0.55
1:A:217:LYS:HB2	1:A:236:GLU:HG3	1.88	0.55
1:A:664:ILE:HG21	2:O:15:ALA:HB2	1.87	0.55
1:A:714:GLN:O	1:A:718:ARG:HG3	2.05	0.55
1:B:97:TYR:CZ	1:B:150:PRO:HB2	2.41	0.55
1:B:216:GLU:HG3	1:B:217:LYS:H	1.72	0.55
1:B:694:VAL:HG23	2:P:18:LEU:HD11	1.88	0.55
1:C:97:TYR:CZ	1:C:150:PRO:HB2	2.41	0.55
1:C:179:LEU:HD23	1:C:179:LEU:N	2.19	0.55
1:C:199:LEU:O	1:C:201:ASP:N	2.32	0.55
1:C:375:GLY:O	1:C:377:GLN:N	2.38	0.55
1:C:581:GLN:HE21	1:C:629:ASN:N	2.02	0.55
1:C:699:GLY:O	1:C:703:ASP:OD2	2.25	0.55
1:D:180:ASP:CG	1:D:181:ILE:N	2.58	0.55
1:D:230:ILE:HG13	1:D:237:PHE:CE2	2.42	0.55
1:D:519:THR:OG1	1:D:520:PRO:HD2	2.06	0.55
1:E:122:GLU:CB	1:E:147:ARG:HG3	2.37	0.55
1:E:479:LYS:CG	1:E:488:LEU:HD21	2.30	0.55
1:E:525:LYS:O	1:E:529:VAL:HG23	2.07	0.55
1:E:664:ILE:HG21	2:S:15:ALA:HB2	1.89	0.55
1:E:671:ARG:HD2	2:S:14:GLU:HG2	1.89	0.55
1:E:694:VAL:HG23	2:S:18:LEU:HD21	1.88	0.55
1:F:671:ARG:HH12	1:F:677:GLY:HA3	1.71	0.55
2:S:16:PHE:CE1	2:S:27:ILE:HD13	2.42	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:90:PRO:O	1:A:92:ASP:N	2.39	0.55
1:A:109:ILE:HD13	1:A:157:LYS:HZ3	1.71	0.55
1:A:217:LYS:HB3	1:A:236:GLU:OE1	2.05	0.55
1:A:579:THR:C	1:A:581:GLN:H	2.14	0.55
1:A:730:ASN:O	1:A:733:GLU:N	2.40	0.55
1:A:764:LEU:C	1:A:766:HIS:H	2.13	0.55
1:C:122:GLU:CB	1:C:147:ARG:HG3	2.37	0.55
1:C:365:PRO:HB2	1:C:367:ASP:O	2.07	0.55
1:D:76:LEU:O	1:D:79:ILE:N	2.39	0.55
1:D:134:LYS:C	1:D:136:PRO:HD3	2.32	0.55
1:D:365:PRO:HB2	1:D:367:ASP:O	2.07	0.55
1:E:311:HIS:O	1:E:314:ALA:HB3	2.07	0.55
1:E:521:ASN:HB3	1:E:524:GLU:HB3	1.87	0.55
1:F:413:LEU:HB2	1:F:419:ILE:CG1	2.33	0.55
1:F:748:TYR:O	1:F:751:TYR:HB3	2.06	0.55
2:Q:124:MET:O	2:Q:125:ILE:C	2.47	0.55
1:A:165:GLN:OE1	1:A:252:ASP:HB3	2.06	0.55
1:B:657:ILE:HD11	1:B:701:LEU:HD23	1.88	0.55
1:B:667:LEU:O	1:B:668:SER:C	2.49	0.55
1:D:217:LYS:HB3	1:D:236:GLU:OE1	2.06	0.55
1:D:667:LEU:O	1:D:668:SER:C	2.50	0.55
1:E:78:LYS:HG3	1:E:79:ILE:N	2.21	0.55
1:E:216:GLU:HG3	1:E:217:LYS:H	1.72	0.55
1:E:349:ASN:HD22	1:E:350:VAL:H	1.53	0.55
1:E:667:LEU:O	1:E:668:SER:C	2.50	0.55
1:F:281:GLU:O	1:F:285:LYS:HG2	2.07	0.55
1:F:365:PRO:HB2	1:F:367:ASP:O	2.07	0.55
1:F:579:THR:C	1:F:581:GLN:H	2.15	0.55
1:F:667:LEU:HB3	2:T:14:GLU:OE2	2.06	0.55
2:Q:94:LYS:HB3	2:Q:94:LYS:NZ	2.22	0.55
2:T:16:PHE:CE1	2:T:27:ILE:HD13	2.42	0.55
1:A:189:ASP:C	1:A:191:GLU:H	2.14	0.55
1:B:165:GLN:OE1	1:B:252:ASP:HB3	2.07	0.55
1:B:197:LYS:HE2	1:B:263:ASP:HB3	1.89	0.55
1:B:217:LYS:HB3	1:B:236:GLU:OE1	2.07	0.55
1:B:387:ASN:N	1:B:387:ASN:ND2	2.50	0.55
1:C:115:LYS:HB2	1:C:118:GLN:HG2	1.86	0.55
1:C:201:ASP:O	1:C:210:PHE:CE2	2.60	0.55
1:C:217:LYS:HB3	1:C:236:GLU:OE1	2.06	0.55
1:C:217:LYS:HB2	1:C:236:GLU:HG3	1.88	0.55
1:F:325:TYR:CE1	1:F:598:PRO:HD3	2.42	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:O:94:LYS:NZ	2:O:94:LYS:HB3	2.22	0.55
2:P:16:PHE:CE1	2:P:27:ILE:HD13	2.41	0.55
2:S:58:ASP:C	2:S:60:ASN:N	2.65	0.55
1:A:90:PRO:HD3	1:A:249:PHE:CE2	2.42	0.55
1:A:179:LEU:HD23	1:A:179:LEU:N	2.19	0.55
1:B:122:GLU:CB	1:B:147:ARG:HG3	2.37	0.55
1:B:281:GLU:O	1:B:285:LYS:HG2	2.05	0.55
1:B:297:LYS:HZ3	1:B:601:GLU:HB3	1.72	0.55
1:B:308:VAL:O	1:B:311:HIS:HB2	2.07	0.55
1:C:308:VAL:O	1:C:311:HIS:HB2	2.06	0.55
1:C:519:THR:OG1	1:C:520:PRO:HD2	2.05	0.55
1:C:694:VAL:HG23	2:Q:18:LEU:HD11	1.88	0.55
1:D:90:PRO:HD3	1:D:249:PHE:CE2	2.42	0.55
1:D:281:GLU:O	1:D:285:LYS:HG2	2.06	0.55
1:E:80:GLN:HG3	1:E:80:GLN:O	2.07	0.55
1:E:581:GLN:HE21	1:E:629:ASN:N	2.00	0.55
1:E:732:ILE:HG23	1:E:749:PHE:HD1	1.70	0.55
1:F:179:LEU:H	1:F:179:LEU:CD2	2.18	0.55
1:F:308:VAL:O	1:F:311:HIS:HB2	2.06	0.55
1:F:512:GLU:O	1:F:516:VAL:HG23	2.06	0.55
1:F:794:GLN:NE2	1:F:795:LYS:HG3	2.22	0.55
2:Q:48:LEU:CA	2:Q:51:MET:HE2	2.34	0.55
1:B:217:LYS:HB2	1:B:236:GLU:HG3	1.88	0.55
1:C:93:VAL:HG23	1:C:179:LEU:HD13	1.89	0.55
1:C:625:LEU:HD12	1:C:626:TYR:N	2.22	0.55
1:D:97:TYR:CZ	1:D:150:PRO:HB2	2.42	0.55
1:D:499:PRO:HG2	1:D:504:ILE:HD11	1.88	0.55
1:D:639:ASN:H	1:D:639:ASN:HD22	1.53	0.55
1:D:666:ASN:O	1:D:670:ILE:HG13	2.07	0.55
1:E:325:TYR:CE1	1:E:598:PRO:HD3	2.42	0.55
1:E:387:ASN:N	1:E:387:ASN:ND2	2.49	0.55
1:E:697:ILE:HG21	1:E:732:ILE:HD11	1.88	0.55
1:F:76:LEU:O	1:F:79:ILE:N	2.40	0.55
1:A:66:LEU:HD12	1:A:103:GLU:HA	1.88	0.54
1:A:697:ILE:HD13	1:A:732:ILE:CD1	2.37	0.54
1:B:78:LYS:HG3	1:B:79:ILE:N	2.21	0.54
1:B:115:LYS:HG3	1:B:153:ILE:CG2	2.38	0.54
1:B:671:ARG:HD2	2:P:14:GLU:HG2	1.89	0.54
1:E:76:LEU:O	1:E:80:GLN:N	2.38	0.54
1:E:110:ASP:O	1:E:111:LEU:C	2.51	0.54
1:E:115:LYS:HG3	1:E:153:ILE:CG2	2.38	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:308:VAL:O	1:E:311:HIS:HB2	2.07	0.54
1:F:197:LYS:HB3	1:F:197:LYS:HZ2	1.70	0.54
1:F:609:GLU:O	1:F:613:ARG:N	2.33	0.54
2:Q:10:ALA:O	2:Q:14:GLU:HB2	2.07	0.54
2:R:36:MET:HE1	2:R:51:MET:HE1	1.88	0.54
2:S:16:PHE:HE1	2:S:27:ILE:HG12	1.72	0.54
2:T:16:PHE:HE1	2:T:27:ILE:HG12	1.72	0.54
2:T:137:ASN:OD1	2:T:139:GLU:HB2	2.07	0.54
1:B:110:ASP:O	1:B:111:LEU:C	2.51	0.54
1:B:334:LEU:H	1:B:334:LEU:CD1	2.20	0.54
1:B:375:GLY:O	1:B:377:GLN:N	2.38	0.54
1:B:639:ASN:H	1:B:639:ASN:HD22	1.54	0.54
1:C:80:GLN:HG3	1:C:80:GLN:O	2.06	0.54
1:C:179:LEU:H	1:C:179:LEU:CD2	2.17	0.54
1:C:254:ARG:HD2	1:C:254:ARG:N	2.23	0.54
1:C:579:THR:C	1:C:581:GLN:H	2.15	0.54
1:D:699:GLY:O	1:D:703:ASP:OD2	2.26	0.54
1:E:90:PRO:O	1:E:92:ASP:N	2.40	0.54
1:E:742:ALA:HB1	1:E:744:GLU:OE1	2.07	0.54
1:F:109:ILE:HD13	1:F:157:LYS:HZ3	1.72	0.54
1:F:115:LYS:HE3	1:F:116:GLU:N	2.22	0.54
1:F:712:PHE:HB3	1:F:716:LYS:HG2	1.90	0.54
2:P:97:ASN:HD22	2:P:97:ASN:C	2.15	0.54
1:A:110:ASP:O	1:A:111:LEU:C	2.50	0.54
1:A:715:GLU:HA	1:A:718:ARG:HH12	1.69	0.54
1:B:64:ASN:HD22	1:B:64:ASN:N	2.04	0.54
1:B:513:TRP:CZ3	1:B:517:VAL:HG11	2.43	0.54
1:B:730:ASN:O	1:B:733:GLU:N	2.40	0.54
1:C:597:ASN:CB	1:C:598:PRO:HD2	2.35	0.54
1:C:656:THR:O	1:C:755:ARG:NH1	2.40	0.54
1:C:748:TYR:O	1:C:751:TYR:HB3	2.07	0.54
1:D:90:PRO:O	1:D:92:ASP:N	2.40	0.54
1:D:115:LYS:HE3	1:D:116:GLU:N	2.22	0.54
1:D:657:ILE:HD11	1:D:701:LEU:HD23	1.88	0.54
1:E:671:ARG:O	1:E:673:SER:N	2.40	0.54
1:F:334:LEU:H	1:F:334:LEU:CD1	2.18	0.54
1:F:450:ASN:ND2	1:F:452:GLU:CD	2.66	0.54
2:O:6:GLU:HG3	2:O:7:GLU:H	1.72	0.54
2:R:137:ASN:OD1	2:R:139:GLU:HB2	2.07	0.54
2:T:17:SER:O	2:T:20:ASP:N	2.37	0.54
1:A:76:LEU:O	1:A:79:ILE:N	2.41	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:76:LEU:O	1:A:80:GLN:N	2.40	0.54
1:A:230:ILE:HG13	1:A:237:PHE:CE2	2.43	0.54
1:A:325:TYR:CE1	1:A:598:PRO:HD3	2.41	0.54
1:A:364:ILE:O	1:A:477:MET:HG2	2.08	0.54
1:B:325:TYR:CE1	1:B:598:PRO:HD3	2.42	0.54
1:B:364:ILE:O	1:B:477:MET:HG2	2.08	0.54
1:C:66:LEU:HD12	1:C:103:GLU:HA	1.88	0.54
1:C:216:GLU:HG3	1:C:217:LYS:H	1.72	0.54
1:D:78:LYS:HG3	1:D:79:ILE:N	2.21	0.54
1:D:122:GLU:CB	1:D:147:ARG:HG3	2.37	0.54
1:D:189:ASP:C	1:D:191:GLU:H	2.14	0.54
1:D:730:ASN:O	1:D:733:GLU:N	2.40	0.54
1:E:710:HIS:C	1:E:712:PHE:N	2.59	0.54
1:F:179:LEU:HD23	1:F:179:LEU:N	2.19	0.54
1:F:700:TYR:CD1	1:F:727:GLN:C	2.86	0.54
1:A:326:ILE:C	1:A:327:LEU:HD12	2.32	0.54
1:A:375:GLY:O	1:A:377:GLN:N	2.40	0.54
1:A:405:LEU:HD12	1:A:405:LEU:N	2.22	0.54
1:A:512:GLU:O	1:A:516:VAL:HG23	2.06	0.54
1:A:794:GLN:NE2	1:A:795:LYS:HG3	2.22	0.54
1:B:230:ILE:HG13	1:B:237:PHE:CE2	2.43	0.54
1:B:311:HIS:O	1:B:314:ALA:HB3	2.07	0.54
1:C:132:GLY:O	1:C:133:GLU:HB2	2.06	0.54
1:C:230:ILE:HG13	1:C:237:PHE:CE2	2.42	0.54
1:C:311:HIS:O	1:C:314:ALA:HB3	2.07	0.54
1:C:494:LEU:HB3	1:C:579:THR:HG22	1.89	0.54
1:C:609:GLU:O	1:C:613:ARG:N	2.33	0.54
1:C:671:ARG:O	1:C:673:SER:N	2.40	0.54
1:C:700:TYR:CD1	1:C:727:GLN:C	2.86	0.54
1:D:667:LEU:HB3	2:R:14:GLU:OE2	2.08	0.54
1:D:697:ILE:HD13	1:D:732:ILE:CD1	2.37	0.54
1:E:639:ASN:H	1:E:639:ASN:HD22	1.54	0.54
1:F:217:LYS:HB2	1:F:236:GLU:HG3	1.88	0.54
1:F:327:LEU:HD12	1:F:327:LEU:N	2.22	0.54
2:P:10:ALA:O	2:P:14:GLU:HB2	2.07	0.54
2:S:17:SER:O	2:S:20:ASP:N	2.39	0.54
1:A:615:ILE:CD1	1:A:645:TRP:HH2	2.20	0.54
1:A:713:SER:O	1:A:714:GLN:C	2.51	0.54
1:B:175:LYS:HB2	1:B:175:LYS:NZ	2.22	0.54
1:B:326:ILE:C	1:B:327:LEU:HD12	2.32	0.54
1:B:405:LEU:N	1:B:405:LEU:HD12	2.23	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:712:PHE:HB3	1:B:716:LYS:HG2	1.90	0.54
1:C:79:ILE:C	1:C:81:GLN:N	2.60	0.54
1:C:90:PRO:HD3	1:C:249:PHE:CE2	2.42	0.54
1:C:197:LYS:HB3	1:C:197:LYS:HZ2	1.71	0.54
1:C:513:TRP:CZ3	1:C:517:VAL:HG11	2.43	0.54
1:C:549:LEU:HG	1:C:550:SER:O	2.08	0.54
1:C:666:ASN:O	1:C:670:ILE:HG13	2.07	0.54
1:C:764:LEU:C	1:C:766:HIS:H	2.14	0.54
1:D:132:GLY:O	1:D:133:GLU:HB2	2.07	0.54
1:D:165:GLN:OE1	1:D:252:ASP:HB3	2.06	0.54
1:D:197:LYS:HB3	1:D:197:LYS:HZ2	1.72	0.54
1:E:387:ASN:HB3	1:E:477:MET:SD	2.47	0.54
1:E:713:SER:O	1:E:714:GLN:C	2.51	0.54
1:E:748:TYR:O	1:E:751:TYR:HB3	2.08	0.54
1:F:110:ASP:O	1:F:111:LEU:C	2.50	0.54
1:F:764:LEU:C	1:F:766:HIS:H	2.15	0.54
2:P:16:PHE:HE1	2:P:27:ILE:HG12	1.73	0.54
1:A:115:LYS:HD3	1:A:153:ILE:HD13	1.90	0.54
1:A:513:TRP:CZ3	1:A:517:VAL:HG11	2.43	0.54
1:A:667:LEU:HB3	2:O:14:GLU:OE2	2.08	0.54
1:B:112:VAL:O	1:B:114:HIS:N	2.39	0.54
1:B:365:PRO:HB2	1:B:367:ASP:O	2.07	0.54
1:B:579:THR:C	1:B:581:GLN:H	2.15	0.54
1:C:165:GLN:OE1	1:C:252:ASP:HB3	2.07	0.54
1:C:189:ASP:C	1:C:191:GLU:H	2.14	0.54
1:C:334:LEU:H	1:C:334:LEU:CD1	2.19	0.54
1:D:326:ILE:C	1:D:327:LEU:HD12	2.32	0.54
1:D:494:LEU:HB3	1:D:579:THR:HG22	1.90	0.54
1:E:375:GLY:O	1:E:377:GLN:N	2.38	0.54
1:E:405:LEU:N	1:E:405:LEU:HD12	2.23	0.54
1:F:120:LEU:C	1:F:120:LEU:CD1	2.81	0.54
1:F:349:ASN:HD22	1:F:350:VAL:H	1.53	0.54
1:F:609:GLU:OE2	1:F:609:GLU:N	2.33	0.54
1:F:671:ARG:O	1:F:673:SER:N	2.40	0.54
2:R:10:ALA:O	2:R:14:GLU:HB2	2.07	0.54
2:S:109:MET:HG3	2:S:116:LEU:CD1	2.36	0.54
1:A:122:GLU:CB	1:A:147:ARG:HG3	2.37	0.54
1:A:201:ASP:O	1:A:210:PHE:CE2	2.60	0.54
1:A:216:GLU:HG3	1:A:217:LYS:H	1.72	0.54
1:A:671:ARG:O	1:A:673:SER:N	2.40	0.54
1:B:132:GLY:O	1:B:133:GLU:HB2	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:671:ARG:O	1:B:673:SER:N	2.40	0.54
1:C:134:LYS:C	1:C:136:PRO:HD3	2.33	0.54
1:C:667:LEU:HB3	2:Q:14:GLU:OE2	2.08	0.54
1:C:712:PHE:HB3	1:C:716:LYS:HG2	1.90	0.54
1:D:181:ILE:HD12	1:D:238:GLN:OE1	2.08	0.54
1:D:201:ASP:O	1:D:210:PHE:CE2	2.61	0.54
1:D:254:ARG:HD2	1:D:254:ARG:N	2.23	0.54
1:D:325:TYR:CE1	1:D:598:PRO:HD3	2.43	0.54
1:E:109:ILE:HD13	1:E:157:LYS:HZ3	1.73	0.54
1:E:217:LYS:HB2	1:E:236:GLU:HG3	1.89	0.54
1:E:364:ILE:O	1:E:477:MET:HG2	2.08	0.54
1:E:365:PRO:HB2	1:E:367:ASP:O	2.08	0.54
1:F:141:PHE:N	1:F:141:PHE:CD1	2.69	0.54
1:F:181:ILE:HD12	1:F:238:GLN:OE1	2.08	0.54
1:F:216:GLU:HG3	1:F:217:LYS:H	1.73	0.54
1:F:254:ARG:HD2	1:F:254:ARG:N	2.23	0.54
1:F:639:ASN:H	1:F:639:ASN:HD22	1.55	0.54
1:F:666:ASN:O	1:F:670:ILE:HG13	2.08	0.54
2:Q:97:ASN:ND2	2:Q:97:ASN:H	2.06	0.54
2:S:10:ALA:O	2:S:14:GLU:HB2	2.07	0.54
2:T:16:PHE:CE1	2:T:27:ILE:HG12	2.43	0.54
1:B:90:PRO:O	1:B:92:ASP:N	2.41	0.54
1:B:279:ILE:HG22	1:B:283:LEU:CD1	2.38	0.54
1:B:792:VAL:HG12	1:B:796:ILE:HD11	1.90	0.54
1:C:371:SER:O	1:C:372:LYS:C	2.51	0.54
1:D:217:LYS:HB2	1:D:236:GLU:HG3	1.89	0.54
1:D:500:SER:HG	1:D:502:THR:HG1	1.56	0.54
1:E:165:GLN:O	1:E:167:LYS:N	2.40	0.54
1:E:671:ARG:HH12	1:E:677:GLY:HA3	1.72	0.54
1:E:730:ASN:O	1:E:733:GLU:N	2.39	0.54
1:F:189:ASP:C	1:F:191:GLU:H	2.14	0.54
1:F:659:THR:H	1:F:662:GLU:HB3	1.73	0.54
1:F:694:VAL:HG22	2:T:18:LEU:HD21	1.89	0.54
1:F:713:SER:O	1:F:714:GLN:C	2.51	0.54
2:O:137:ASN:OD1	2:O:139:GLU:HB2	2.08	0.54
2:Q:97:ASN:C	2:Q:97:ASN:HD22	2.15	0.54
1:A:279:ILE:HG22	1:A:283:LEU:CD1	2.38	0.54
1:A:308:VAL:O	1:A:311:HIS:HB2	2.08	0.54
1:A:666:ASN:O	1:A:670:ILE:HG13	2.07	0.54
1:A:694:VAL:HG23	2:O:18:LEU:HD11	1.89	0.54
1:A:697:ILE:HG21	1:A:732:ILE:HD11	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:120:LEU:C	1:B:120:LEU:CD1	2.81	0.54
1:B:201:ASP:O	1:B:210:PHE:CE2	2.61	0.54
1:B:450:ASN:ND2	1:B:452:GLU:CD	2.66	0.54
1:B:666:ASN:O	1:B:670:ILE:HG13	2.08	0.54
1:B:794:GLN:NE2	1:B:795:LYS:HG3	2.22	0.54
1:C:694:VAL:HG22	2:Q:18:LEU:HD21	1.90	0.54
1:C:794:GLN:NE2	1:C:795:LYS:HG3	2.23	0.54
1:D:76:LEU:O	1:D:80:GLN:N	2.39	0.54
1:D:120:LEU:C	1:D:120:LEU:CD1	2.81	0.54
1:D:671:ARG:O	1:D:673:SER:N	2.40	0.54
1:D:697:ILE:HG21	1:D:732:ILE:HD11	1.88	0.54
1:E:700:TYR:CD1	1:E:727:GLN:C	2.86	0.54
1:F:132:GLY:O	1:F:133:GLU:HB2	2.07	0.54
1:F:311:HIS:O	1:F:314:ALA:HB3	2.07	0.54
1:F:597:ASN:CB	1:F:598:PRO:HD2	2.35	0.54
2:P:52:ILE:HG23	2:P:53:ASN:N	2.23	0.54
2:P:109:MET:HG3	2:P:116:LEU:CD1	2.37	0.54
2:S:6:GLU:HG3	2:S:7:GLU:H	1.73	0.54
1:A:334:LEU:H	1:A:334:LEU:CD1	2.20	0.53
1:A:549:LEU:HG	1:A:550:SER:O	2.08	0.53
1:B:90:PRO:HD3	1:B:249:PHE:CE2	2.43	0.53
1:C:110:ASP:O	1:C:111:LEU:C	2.50	0.53
1:C:120:LEU:C	1:C:120:LEU:CD1	2.81	0.53
1:C:181:ILE:HD12	1:C:238:GLN:OE1	2.08	0.53
1:D:93:VAL:HG23	1:D:179:LEU:HD13	1.90	0.53
1:D:197:LYS:HE2	1:D:263:ASP:HB3	1.90	0.53
1:D:794:GLN:NE2	1:D:795:LYS:HG3	2.22	0.53
1:E:181:ILE:HD12	1:E:238:GLN:OE1	2.09	0.53
1:E:615:ILE:CD1	1:E:645:TRP:HH2	2.21	0.53
1:F:549:LEU:HG	1:F:550:SER:O	2.08	0.53
2:Q:12:PHE:HE1	2:Q:72:MET:CE	2.14	0.53
2:R:52:ILE:HG23	2:R:53:ASN:N	2.23	0.53
2:T:64:ASP:OD2	2:T:67:GLU:HG3	2.08	0.53
1:A:120:LEU:C	1:A:120:LEU:CD1	2.81	0.53
1:A:333:LYS:C	1:A:335:ALA:H	2.16	0.53
1:C:76:LEU:O	1:C:79:ILE:N	2.41	0.53
1:C:90:PRO:O	1:C:92:ASP:N	2.41	0.53
1:C:142:VAL:HG22	1:C:154:ILE:HD12	1.90	0.53
1:C:298:GLY:C	1:C:300:LYS:N	2.66	0.53
1:E:85:LEU:O	1:E:88:LYS:HB3	2.08	0.53
1:E:197:LYS:HE2	1:E:263:ASP:HB3	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:678:VAL:HG13	1:E:745:TYR:HD2	1.74	0.53
1:F:115:LYS:HG3	1:F:153:ILE:CG2	2.39	0.53
1:F:197:LYS:NZ	1:F:197:LYS:CB	2.72	0.53
1:F:513:TRP:CZ3	1:F:517:VAL:HG11	2.43	0.53
1:F:656:THR:O	1:F:755:ARG:NH1	2.41	0.53
1:F:709:ASN:O	1:F:717:LYS:HE3	2.08	0.53
1:F:742:ALA:HB1	1:F:744:GLU:OE1	2.08	0.53
2:Q:52:ILE:HG23	2:Q:53:ASN:N	2.23	0.53
1:A:450:ASN:ND2	1:A:452:GLU:CD	2.66	0.53
1:A:671:ARG:HD2	2:O:14:GLU:HG2	1.90	0.53
1:B:625:LEU:HD12	1:B:626:TYR:N	2.23	0.53
1:C:115:LYS:HG3	1:C:153:ILE:CG2	2.38	0.53
1:C:279:ILE:HG22	1:C:283:LEU:CD1	2.38	0.53
1:C:333:LYS:C	1:C:335:ALA:H	2.16	0.53
1:C:659:THR:H	1:C:662:GLU:HB3	1.74	0.53
1:C:742:ALA:HB1	1:C:744:GLU:OE1	2.08	0.53
1:D:615:ILE:CD1	1:D:645:TRP:HH2	2.21	0.53
1:E:120:LEU:C	1:E:120:LEU:CD1	2.81	0.53
1:E:230:ILE:HG13	1:E:237:PHE:CE2	2.43	0.53
1:E:712:PHE:HB3	1:E:716:LYS:HG2	1.91	0.53
1:E:794:GLN:NE2	1:E:795:LYS:HG3	2.23	0.53
1:F:142:VAL:HG22	1:F:154:ILE:HD12	1.90	0.53
1:F:494:LEU:HB3	1:F:579:THR:HG22	1.88	0.53
2:O:97:ASN:C	2:O:97:ASN:HD22	2.14	0.53
1:A:197:LYS:NZ	1:A:197:LYS:CB	2.70	0.53
1:A:597:ASN:ND2	1:A:601:GLU:CB	2.72	0.53
1:B:142:VAL:HG22	1:B:154:ILE:HD12	1.89	0.53
1:B:583:ASN:ND2	1:B:587:PRO:HA	2.23	0.53
1:C:364:ILE:O	1:C:477:MET:HG2	2.08	0.53
1:D:115:LYS:HD3	1:D:153:ILE:HD13	1.90	0.53
1:D:405:LEU:HD12	1:D:405:LEU:N	2.23	0.53
1:D:450:ASN:ND2	1:D:452:GLU:CD	2.66	0.53
1:D:512:GLU:O	1:D:516:VAL:HG23	2.07	0.53
1:D:513:TRP:CZ3	1:D:517:VAL:HG11	2.43	0.53
1:D:656:THR:O	1:D:755:ARG:NH1	2.40	0.53
1:D:671:ARG:HH11	1:D:671:ARG:HG3	1.74	0.53
1:D:712:PHE:HB3	1:D:716:LYS:HG2	1.90	0.53
1:D:764:LEU:C	1:D:766:HIS:H	2.15	0.53
1:E:90:PRO:HD3	1:E:249:PHE:CE2	2.43	0.53
1:E:513:TRP:CZ3	1:E:517:VAL:HG11	2.43	0.53
1:E:671:ARG:HH11	1:E:671:ARG:HG3	1.74	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:188:LEU:CD2	1:F:188:LEU:N	2.47	0.53
1:F:504:ILE:O	1:F:507:GLN:HB3	2.08	0.53
1:F:615:ILE:CD1	1:F:645:TRP:HH2	2.21	0.53
1:F:697:ILE:HD13	1:F:732:ILE:CD1	2.38	0.53
2:O:109:MET:HG3	2:O:116:LEU:CD1	2.37	0.53
2:Q:16:PHE:CE1	2:Q:27:ILE:HD13	2.43	0.53
2:S:36:MET:HE3	2:S:43:PRO:CG	2.32	0.53
2:T:52:ILE:HG23	2:T:53:ASN:N	2.23	0.53
1:A:694:VAL:HG22	2:O:18:LEU:HD21	1.90	0.53
1:B:371:SER:O	1:B:372:LYS:C	2.52	0.53
1:B:699:GLY:O	1:B:703:ASP:OD2	2.25	0.53
1:B:700:TYR:CD1	1:B:727:GLN:C	2.86	0.53
1:C:325:TYR:CE1	1:C:598:PRO:HD3	2.43	0.53
1:C:327:LEU:HD12	1:C:327:LEU:N	2.22	0.53
1:C:504:ILE:O	1:C:507:GLN:HB3	2.09	0.53
1:C:730:ASN:O	1:C:732:ILE:N	2.42	0.53
1:D:175:LYS:O	1:D:178:SER:N	2.42	0.53
1:D:357:TRP:CH2	1:D:439:ASN:ND2	2.77	0.53
1:D:715:GLU:HA	1:D:718:ARG:HH12	1.71	0.53
1:E:625:LEU:HD12	1:E:626:TYR:N	2.23	0.53
2:O:16:PHE:CE1	2:O:27:ILE:HD13	2.43	0.53
2:O:64:ASP:OD2	2:O:67:GLU:HG3	2.09	0.53
2:S:16:PHE:CE1	2:S:27:ILE:HG12	2.43	0.53
2:T:138:TYR:CE1	2:T:142:VAL:HG22	2.43	0.53
1:A:142:VAL:HA	1:A:154:ILE:HD13	1.90	0.53
1:A:282:SER:HA	1:A:285:LYS:CD	2.38	0.53
1:A:654:ILE:C	1:A:655:ASN:HD22	2.17	0.53
1:A:700:TYR:CD1	1:A:727:GLN:C	2.86	0.53
1:B:387:ASN:HB3	1:B:477:MET:SD	2.49	0.53
1:B:671:ARG:HG3	1:B:671:ARG:HH11	1.74	0.53
1:C:180:ASP:CG	1:C:181:ILE:N	2.59	0.53
1:C:387:ASN:HB3	1:C:477:MET:SD	2.49	0.53
1:C:501:LEU:HD23	2:Q:112:LEU:HD21	1.91	0.53
1:C:709:ASN:O	1:C:717:LYS:HE3	2.09	0.53
1:D:110:ASP:O	1:D:111:LEU:C	2.51	0.53
1:D:297:LYS:HE3	1:D:601:GLU:OE1	2.09	0.53
1:D:792:VAL:HG12	1:D:796:ILE:HD11	1.89	0.53
1:E:279:ILE:HG22	1:E:283:LEU:CD1	2.38	0.53
1:E:450:ASN:ND2	1:E:452:GLU:CD	2.67	0.53
1:E:504:ILE:O	1:E:507:GLN:HB3	2.08	0.53
1:E:699:GLY:O	1:E:703:ASP:OD2	2.26	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:357:TRP:CH2	1:F:439:ASN:ND2	2.77	0.53
1:F:405:LEU:HD12	1:F:405:LEU:N	2.24	0.53
2:R:6:GLU:HG3	2:R:7:GLU:H	1.72	0.53
2:T:97:ASN:C	2:T:97:ASN:HD22	2.15	0.53
1:A:371:SER:O	1:A:372:LYS:C	2.51	0.53
1:A:478:ALA:CB	1:A:486:LYS:O	2.56	0.53
1:A:671:ARG:HH11	1:A:671:ARG:HG3	1.74	0.53
1:A:792:VAL:HG12	1:A:796:ILE:HD11	1.90	0.53
1:B:297:LYS:HE3	1:B:601:GLU:OE1	2.09	0.53
1:B:413:LEU:HB2	1:B:419:ILE:CD1	2.39	0.53
1:B:501:LEU:HD23	2:P:112:LEU:HD21	1.91	0.53
1:B:713:SER:O	1:B:714:GLN:C	2.52	0.53
1:B:742:ALA:HB1	1:B:744:GLU:OE1	2.08	0.53
1:C:512:GLU:O	1:C:516:VAL:HG23	2.08	0.53
1:D:597:ASN:ND2	1:D:601:GLU:CB	2.72	0.53
1:E:99:GLU:C	1:E:101:GLY:N	2.60	0.53
1:E:326:ILE:C	1:E:327:LEU:HD12	2.33	0.53
1:E:371:SER:O	1:E:372:LYS:C	2.52	0.53
1:E:666:ASN:O	1:E:670:ILE:HG13	2.08	0.53
1:E:792:VAL:HG12	1:E:796:ILE:HD11	1.91	0.53
2:O:16:PHE:CE1	2:O:27:ILE:HG12	2.44	0.53
2:O:52:ILE:HG23	2:O:53:ASN:N	2.23	0.53
2:Q:16:PHE:HE1	2:Q:27:ILE:HG12	1.74	0.53
1:A:197:LYS:HB3	1:A:197:LYS:HZ2	1.73	0.53
1:A:357:TRP:CH2	1:A:439:ASN:ND2	2.77	0.53
1:A:712:PHE:HB3	1:A:716:LYS:HG2	1.91	0.53
1:A:742:ALA:HB1	1:A:744:GLU:OE1	2.07	0.53
1:C:413:LEU:HB2	1:C:419:ILE:CD1	2.39	0.53
1:C:567:THR:CG2	1:C:568:GLY:H	2.22	0.53
1:D:85:LEU:O	1:D:88:LYS:HB3	2.09	0.53
1:D:579:THR:C	1:D:581:GLN:H	2.16	0.53
1:D:625:LEU:HD12	1:D:626:TYR:N	2.24	0.53
1:E:254:ARG:HD2	1:E:254:ARG:N	2.22	0.53
1:E:413:LEU:HB2	1:E:419:ILE:CD1	2.39	0.53
1:E:549:LEU:HG	1:E:550:SER:O	2.08	0.53
1:F:387:ASN:HB3	1:F:477:MET:SD	2.48	0.53
1:F:443:GLU:O	1:F:455:TYR:HA	2.09	0.53
2:P:16:PHE:CE1	2:P:27:ILE:HG12	2.44	0.53
2:P:137:ASN:OD1	2:P:139:GLU:HB2	2.09	0.53
2:R:138:TYR:CE1	2:R:142:VAL:HG22	2.44	0.53
2:T:121:VAL:C	2:T:123:GLN:N	2.66	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:699:GLY:O	1:A:703:ASP:OD2	2.27	0.53
1:B:395:GLU:C	1:B:397:GLU:H	2.17	0.53
1:B:667:LEU:HB3	2:P:14:GLU:OE2	2.08	0.53
1:B:671:ARG:HH12	1:B:677:GLY:HA3	1.73	0.53
1:D:173:ILE:HD12	1:D:243:LEU:HG	1.91	0.53
1:D:218:LEU:HD21	1:D:225:ILE:CD1	2.39	0.53
1:D:333:LYS:C	1:D:335:ALA:H	2.17	0.53
1:D:713:SER:O	1:D:714:GLN:C	2.52	0.53
1:E:201:ASP:O	1:E:210:PHE:CE2	2.61	0.53
1:E:297:LYS:HZ3	1:E:601:GLU:HB3	1.74	0.53
1:F:165:GLN:O	1:F:167:LYS:N	2.40	0.53
1:F:364:ILE:O	1:F:477:MET:HG2	2.09	0.53
1:F:710:HIS:C	1:F:712:PHE:N	2.59	0.53
2:Q:16:PHE:CE1	2:Q:27:ILE:HG12	2.44	0.53
2:Q:16:PHE:HA	2:Q:35:VAL:HG11	1.91	0.53
2:Q:18:LEU:HD13	2:Q:19:PHE:CE1	2.44	0.53
2:Q:70:THR:C	2:Q:72:MET:H	2.17	0.53
2:T:70:THR:C	2:T:72:MET:H	2.17	0.53
1:A:115:LYS:HG3	1:A:153:ILE:CG2	2.39	0.53
1:A:162:ASN:O	1:A:163:SER:C	2.52	0.53
1:B:181:ILE:HD12	1:B:238:GLN:OE1	2.09	0.53
1:B:549:LEU:HG	1:B:550:SER:O	2.09	0.53
1:B:656:THR:O	1:B:755:ARG:NH1	2.41	0.53
1:B:700:TYR:HD1	1:B:727:GLN:C	2.17	0.53
1:C:90:PRO:HG2	1:C:93:VAL:CB	2.39	0.53
1:C:197:LYS:NZ	1:C:197:LYS:CB	2.72	0.53
1:C:217:LYS:HZ1	1:C:233:ASN:HB3	1.74	0.53
1:C:435:LEU:H	1:C:445:ARG:HA	1.74	0.53
1:D:216:GLU:HG3	1:D:217:LYS:H	1.73	0.53
1:D:700:TYR:CD1	1:D:727:GLN:C	2.87	0.53
1:E:132:GLY:O	1:E:133:GLU:HB2	2.07	0.53
1:E:134:LYS:C	1:E:136:PRO:HD3	2.34	0.53
1:E:540:ARG:HD3	1:E:627:TYR:CZ	2.44	0.53
1:F:93:VAL:HG23	1:F:179:LEU:HD13	1.90	0.53
1:F:230:ILE:HG13	1:F:237:PHE:CE2	2.43	0.53
2:T:10:ALA:O	2:T:14:GLU:HB2	2.08	0.53
1:A:173:ILE:HD12	1:A:243:LEU:HG	1.91	0.52
1:A:625:LEU:HD12	1:A:626:TYR:N	2.24	0.52
1:B:115:LYS:HD3	1:B:153:ILE:HD13	1.91	0.52
1:C:359:PRO:HG2	1:C:444:PHE:CD2	2.44	0.52
1:C:450:ASN:ND2	1:C:452:GLU:CD	2.66	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:597:ASN:ND2	1:C:601:GLU:CB	2.72	0.52
1:C:671:ARG:HH11	1:C:671:ARG:HG3	1.73	0.52
1:D:364:ILE:O	1:D:477:MET:HG2	2.09	0.52
1:D:371:SER:O	1:D:372:LYS:C	2.52	0.52
1:D:387:ASN:HB3	1:D:477:MET:SD	2.49	0.52
1:D:609:GLU:O	1:D:612:GLY:N	2.43	0.52
1:E:173:ILE:HD12	1:E:243:LEU:HG	1.91	0.52
1:E:659:THR:H	1:E:662:GLU:HB3	1.73	0.52
1:F:115:LYS:HB2	1:F:118:GLN:HG2	1.88	0.52
1:F:162:ASN:O	1:F:163:SER:C	2.52	0.52
2:O:16:PHE:HE1	2:O:27:ILE:HG12	1.74	0.52
2:P:16:PHE:HA	2:P:35:VAL:HG11	1.91	0.52
2:S:52:ILE:HG23	2:S:53:ASN:N	2.22	0.52
1:A:639:ASN:H	1:A:639:ASN:HD22	1.54	0.52
1:B:162:ASN:O	1:B:163:SER:C	2.52	0.52
1:B:197:LYS:NZ	1:B:197:LYS:CB	2.72	0.52
1:C:197:LYS:HE2	1:C:263:ASP:HB3	1.92	0.52
1:C:302:LEU:HB2	1:C:602:PHE:CD1	2.37	0.52
1:C:555:GLN:HG3	1:C:556:MET:N	2.24	0.52
1:D:105:TYR:O	1:D:154:ILE:N	2.34	0.52
1:D:549:LEU:HG	1:D:550:SER:O	2.08	0.52
1:E:142:VAL:HG22	1:E:154:ILE:HD12	1.90	0.52
1:E:142:VAL:HA	1:E:154:ILE:HD13	1.90	0.52
1:E:408:LEU:HD12	1:E:408:LEU:N	2.01	0.52
1:F:285:LYS:C	1:F:287:GLY:H	2.17	0.52
1:F:609:GLU:O	1:F:612:GLY:N	2.42	0.52
1:F:671:ARG:HH11	1:F:671:ARG:HG3	1.74	0.52
2:S:138:TYR:CE1	2:S:142:VAL:HG22	2.44	0.52
1:A:494:LEU:HB3	1:A:579:THR:HG22	1.90	0.52
1:A:555:GLN:HG3	1:A:556:MET:N	2.24	0.52
1:C:218:LEU:HD21	1:C:225:ILE:CD1	2.40	0.52
1:C:297:LYS:HE3	1:C:601:GLU:OE1	2.10	0.52
1:C:700:TYR:HD1	1:C:727:GLN:C	2.17	0.52
1:D:142:VAL:HA	1:D:154:ILE:HD13	1.91	0.52
1:D:671:ARG:HD2	2:R:14:GLU:HG2	1.91	0.52
1:E:333:LYS:C	1:E:335:ALA:H	2.16	0.52
1:F:279:ILE:HG22	1:F:283:LEU:CD1	2.39	0.52
1:F:413:LEU:HB2	1:F:419:ILE:CD1	2.40	0.52
1:F:700:TYR:HD1	1:F:727:GLN:C	2.17	0.52
2:Q:106:ARG:HB2	2:Q:121:VAL:HG21	1.91	0.52
2:R:16:PHE:CE1	2:R:27:ILE:HD13	2.44	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:R:70:THR:C	2:R:72:MET:H	2.17	0.52
2:R:114:GLU:OE2	2:R:114:GLU:HA	2.09	0.52
2:S:16:PHE:HA	2:S:35:VAL:HG11	1.91	0.52
2:S:64:ASP:OD2	2:S:67:GLU:HG3	2.09	0.52
2:S:70:THR:C	2:S:72:MET:H	2.18	0.52
1:A:142:VAL:HG22	1:A:154:ILE:HD12	1.90	0.52
1:A:254:ARG:HD2	1:A:254:ARG:N	2.24	0.52
1:A:298:GLY:C	1:A:300:LYS:N	2.66	0.52
1:A:504:ILE:O	1:A:507:GLN:HB3	2.10	0.52
1:A:700:TYR:HD1	1:A:727:GLN:C	2.17	0.52
1:B:85:LEU:O	1:B:88:LYS:HB3	2.10	0.52
1:B:93:VAL:HG23	1:B:179:LEU:HD13	1.90	0.52
1:B:684:ASP:C	1:B:686:ASP:H	2.18	0.52
1:B:788:ASP:O	1:B:789:ASN:C	2.53	0.52
1:C:561:ASN:HA	1:C:564:VAL:HG22	1.92	0.52
1:E:197:LYS:NZ	1:E:197:LYS:CB	2.72	0.52
1:E:395:GLU:C	1:E:397:GLU:H	2.18	0.52
1:E:478:ALA:CB	1:E:486:LYS:O	2.57	0.52
1:E:583:ASN:ND2	1:E:587:PRO:HA	2.24	0.52
1:F:76:LEU:O	1:F:80:GLN:N	2.41	0.52
1:F:85:LEU:O	1:F:88:LYS:HB3	2.10	0.52
1:F:148:GLU:HG3	1:F:149:THR:HG22	1.91	0.52
2:P:120:GLU:HA	2:P:123:GLN:HB2	1.91	0.52
2:Q:64:ASP:OD2	2:Q:67:GLU:HG3	2.09	0.52
2:S:41:GLN:O	2:S:42:ASN:HB3	2.10	0.52
1:A:118:GLN:OE1	1:A:118:GLN:HA	2.09	0.52
1:A:173:ILE:C	1:A:175:LYS:N	2.67	0.52
1:A:656:THR:O	1:A:755:ARG:NH1	2.41	0.52
1:A:730:ASN:C	1:A:732:ILE:N	2.67	0.52
1:C:654:ILE:C	1:C:655:ASN:HD22	2.17	0.52
1:C:730:ASN:C	1:C:732:ILE:N	2.66	0.52
1:D:279:ILE:HG22	1:D:283:LEU:CD1	2.40	0.52
1:D:567:THR:CG2	1:D:568:GLY:H	2.22	0.52
1:E:285:LYS:C	1:E:287:GLY:H	2.17	0.52
1:F:90:PRO:HG2	1:F:93:VAL:CB	2.40	0.52
1:F:140:ARG:HA	1:F:140:ARG:HE	1.75	0.52
1:F:297:LYS:HE3	1:F:601:GLU:OE1	2.10	0.52
1:F:333:LYS:C	1:F:335:ALA:H	2.17	0.52
1:F:678:VAL:HG13	1:F:745:TYR:HD2	1.73	0.52
1:F:730:ASN:C	1:F:732:ILE:N	2.66	0.52
2:R:9:ILE:CD1	2:R:69:LEU:HD11	2.39	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:R:41:GLN:O	2:R:42:ASN:HB3	2.10	0.52
2:S:137:ASN:OD1	2:S:139:GLU:HB2	2.09	0.52
1:A:202:ASP:HB2	1:A:208:LEU:CD2	2.39	0.52
1:A:659:THR:H	1:A:662:GLU:HB3	1.73	0.52
1:B:189:ASP:C	1:B:191:GLU:H	2.15	0.52
1:B:298:GLY:C	1:B:300:LYS:N	2.67	0.52
1:C:678:VAL:HG13	1:C:745:TYR:HD2	1.75	0.52
1:C:684:ASP:C	1:C:686:ASP:H	2.18	0.52
1:E:90:PRO:HG2	1:E:93:VAL:CB	2.38	0.52
1:E:161:ILE:CG2	1:E:168:GLU:HB2	2.40	0.52
1:E:357:TRP:CH2	1:E:439:ASN:ND2	2.77	0.52
1:E:435:LEU:H	1:E:445:ARG:HA	1.74	0.52
1:E:500:SER:HG	1:E:502:THR:HG1	1.57	0.52
1:E:633:ASN:O	1:E:642:TYR:HE1	1.92	0.52
1:E:697:ILE:HD13	1:E:732:ILE:CD1	2.38	0.52
1:F:218:LEU:HD21	1:F:225:ILE:CD1	2.40	0.52
2:O:138:TYR:CE1	2:O:142:VAL:HG22	2.44	0.52
2:Q:58:ASP:C	2:Q:60:ASN:N	2.65	0.52
2:R:16:PHE:HA	2:R:35:VAL:HG11	1.92	0.52
2:T:17:SER:O	2:T:18:LEU:C	2.53	0.52
2:T:94:LYS:HE3	2:T:107:HIS:CD2	2.45	0.52
1:A:93:VAL:HG23	1:A:179:LEU:HD13	1.89	0.52
1:A:140:ARG:HA	1:A:140:ARG:HE	1.75	0.52
1:A:501:LEU:HD23	2:O:112:LEU:HD21	1.92	0.52
1:B:142:VAL:HA	1:B:154:ILE:HD13	1.91	0.52
1:C:443:GLU:O	1:C:455:TYR:HA	2.09	0.52
1:C:609:GLU:O	1:C:612:GLY:N	2.43	0.52
1:C:792:VAL:HG12	1:C:796:ILE:HD11	1.90	0.52
1:D:115:LYS:HG3	1:D:153:ILE:CG2	2.39	0.52
1:D:443:GLU:O	1:D:455:TYR:HA	2.10	0.52
1:D:504:ILE:O	1:D:507:GLN:HB3	2.09	0.52
1:D:555:GLN:HG3	1:D:556:MET:N	2.25	0.52
1:F:90:PRO:HD3	1:F:249:PHE:CE2	2.44	0.52
1:F:118:GLN:OE1	1:F:118:GLN:HA	2.10	0.52
1:F:298:GLY:C	1:F:300:LYS:N	2.66	0.52
1:F:371:SER:O	1:F:372:LYS:C	2.52	0.52
1:F:555:GLN:HG3	1:F:556:MET:N	2.24	0.52
2:Q:41:GLN:O	2:Q:42:ASN:HB3	2.09	0.52
2:R:64:ASP:OD2	2:R:67:GLU:HG3	2.10	0.52
2:T:41:GLN:O	2:T:42:ASN:HB3	2.09	0.52
1:B:173:ILE:HD12	1:B:243:LEU:HG	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:175:LYS:O	1:B:178:SER:N	2.42	0.52
1:B:254:ARG:HD2	1:B:254:ARG:N	2.23	0.52
1:B:435:LEU:H	1:B:445:ARG:HA	1.75	0.52
1:B:504:ILE:O	1:B:507:GLN:HB3	2.10	0.52
1:C:76:LEU:O	1:C:80:GLN:N	2.42	0.52
1:C:285:LYS:C	1:C:287:GLY:H	2.18	0.52
1:C:357:TRP:CH2	1:C:439:ASN:ND2	2.78	0.52
1:E:609:GLU:O	1:E:612:GLY:N	2.43	0.52
1:E:666:ASN:HB2	1:E:748:TYR:OH	2.10	0.52
1:F:197:LYS:HE2	1:F:263:ASP:HB3	1.92	0.52
1:F:501:LEU:HD23	2:T:112:LEU:HD21	1.91	0.52
1:F:567:THR:CG2	1:F:568:GLY:H	2.22	0.52
2:O:36:MET:HE3	2:O:43:PRO:CG	2.32	0.52
2:O:41:GLN:O	2:O:42:ASN:HB3	2.10	0.52
2:O:94:LYS:HE3	2:O:107:HIS:CD2	2.44	0.52
2:S:106:ARG:HB2	2:S:121:VAL:HG21	1.92	0.52
1:A:148:GLU:HG3	1:A:149:THR:HG22	1.91	0.52
1:A:181:ILE:HD12	1:A:238:GLN:OE1	2.09	0.52
1:A:413:LEU:HB2	1:A:419:ILE:CD1	2.39	0.52
1:B:202:ASP:HB2	1:B:208:LEU:CD2	2.40	0.52
1:B:333:LYS:C	1:B:335:ALA:H	2.17	0.52
1:B:567:THR:CG2	1:B:568:GLY:H	2.22	0.52
1:B:659:THR:H	1:B:662:GLU:HB3	1.74	0.52
1:B:730:ASN:C	1:B:732:ILE:N	2.67	0.52
1:C:202:ASP:HB2	1:C:208:LEU:CD2	2.39	0.52
1:C:615:ILE:CD1	1:C:645:TRP:HH2	2.23	0.52
1:D:435:LEU:H	1:D:445:ARG:HA	1.75	0.52
1:D:742:ALA:HB1	1:D:744:GLU:OE1	2.09	0.52
1:E:162:ASN:O	1:E:163:SER:C	2.53	0.52
1:E:597:ASN:ND2	1:E:601:GLU:CB	2.73	0.52
1:E:667:LEU:HB3	2:S:14:GLU:OE2	2.09	0.52
1:E:700:TYR:HD1	1:E:727:GLN:C	2.17	0.52
1:F:583:ASN:ND2	1:F:587:PRO:HA	2.25	0.52
1:F:792:VAL:HG12	1:F:796:ILE:HD11	1.91	0.52
2:P:36:MET:HE3	2:P:43:PRO:CG	2.31	0.52
2:P:94:LYS:HE3	2:P:107:HIS:CD2	2.45	0.52
2:P:138:TYR:CE1	2:P:142:VAL:HG22	2.45	0.52
2:S:97:ASN:C	2:S:99:TYR:H	2.17	0.52
1:A:397:GLU:O	1:A:479:LYS:HA	2.10	0.52
1:A:730:ASN:O	1:A:732:ILE:N	2.43	0.52
1:B:148:GLU:HG3	1:B:149:THR:HG22	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:709:ASN:O	1:B:717:LYS:HE3	2.09	0.52
1:B:715:GLU:HA	1:B:718:ARG:HH12	1.70	0.52
1:C:142:VAL:HA	1:C:154:ILE:HD13	1.91	0.52
1:C:173:ILE:HD12	1:C:243:LEU:HG	1.92	0.52
1:D:162:ASN:O	1:D:163:SER:C	2.53	0.52
1:D:709:ASN:O	1:D:717:LYS:HE3	2.10	0.52
1:E:298:GLY:C	1:E:300:LYS:N	2.66	0.52
1:F:175:LYS:O	1:F:178:SER:N	2.43	0.52
1:F:395:GLU:C	1:F:397:GLU:H	2.18	0.52
1:F:435:LEU:H	1:F:445:ARG:HA	1.75	0.52
1:F:561:ASN:HA	1:F:564:VAL:HG22	1.91	0.52
1:F:597:ASN:ND2	1:F:601:GLU:CB	2.73	0.52
1:F:730:ASN:O	1:F:732:ILE:N	2.42	0.52
2:S:16:PHE:CE1	2:S:27:ILE:CD1	2.93	0.52
1:B:140:ARG:HA	1:B:140:ARG:HE	1.76	0.51
1:B:357:TRP:CH2	1:B:439:ASN:ND2	2.77	0.51
1:C:118:GLN:OE1	1:C:118:GLN:HA	2.10	0.51
1:C:500:SER:HG	1:C:502:THR:HG1	1.57	0.51
1:D:561:ASN:HA	1:D:564:VAL:HG22	1.92	0.51
1:E:140:ARG:HA	1:E:140:ARG:HE	1.75	0.51
1:E:148:GLU:HG3	1:E:149:THR:HG22	1.91	0.51
1:E:175:LYS:O	1:E:178:SER:N	2.42	0.51
1:E:189:ASP:C	1:E:191:GLU:H	2.15	0.51
1:E:297:LYS:HE3	1:E:601:GLU:OE1	2.10	0.51
1:E:597:ASN:CB	1:E:598:PRO:HD2	2.37	0.51
1:E:709:ASN:O	1:E:717:LYS:HE3	2.09	0.51
1:F:671:ARG:HD2	2:T:14:GLU:HG2	1.92	0.51
2:P:6:GLU:HG3	2:P:7:GLU:H	1.73	0.51
2:P:121:VAL:C	2:P:123:GLN:N	2.67	0.51
2:Q:94:LYS:HE3	2:Q:107:HIS:CD2	2.45	0.51
2:Q:97:ASN:C	2:Q:99:TYR:H	2.18	0.51
2:Q:117:THR:OG1	2:Q:119:GLU:HG2	2.10	0.51
2:T:97:ASN:C	2:T:99:TYR:H	2.18	0.51
1:A:395:GLU:C	1:A:397:GLU:H	2.18	0.51
1:B:90:PRO:HG2	1:B:93:VAL:CB	2.40	0.51
1:B:134:LYS:C	1:B:136:PRO:HD3	2.35	0.51
1:B:161:ILE:CG2	1:B:168:GLU:HB2	2.40	0.51
1:C:162:ASN:O	1:C:163:SER:C	2.53	0.51
1:D:379:ALA:O	1:D:383:GLY:N	2.43	0.51
1:D:413:LEU:HB2	1:D:419:ILE:CD1	2.40	0.51
1:D:730:ASN:C	1:D:732:ILE:N	2.67	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:443:GLU:O	1:E:455:TYR:HA	2.10	0.51
1:F:625:LEU:HD12	1:F:625:LEU:C	2.36	0.51
2:T:97:ASN:ND2	2:T:97:ASN:H	2.07	0.51
1:A:609:GLU:O	1:A:612:GLY:N	2.44	0.51
1:B:77:ASP:O	1:B:81:GLN:CB	2.56	0.51
1:B:561:ASN:HA	1:B:564:VAL:HG22	1.91	0.51
1:B:615:ILE:CD1	1:B:645:TRP:HH2	2.22	0.51
1:C:148:GLU:HG3	1:C:149:THR:HG22	1.91	0.51
1:E:214:PHE:CD1	1:E:218:LEU:HD23	2.44	0.51
1:F:581:GLN:HE22	1:F:629:ASN:H	1.53	0.51
2:S:121:VAL:C	2:S:123:GLN:N	2.67	0.51
1:A:90:PRO:HG2	1:A:93:VAL:CB	2.40	0.51
1:A:297:LYS:HE3	1:A:601:GLU:OE1	2.10	0.51
1:B:597:ASN:ND2	1:B:601:GLU:CB	2.74	0.51
1:C:372:LYS:HG3	1:C:373:LYS:HG2	1.93	0.51
1:D:89:ILE:HG22	1:D:93:VAL:CG1	2.30	0.51
1:D:148:GLU:HG3	1:D:149:THR:HG22	1.91	0.51
1:E:565:LYS:C	1:E:567:THR:H	2.19	0.51
1:E:654:ILE:C	1:E:655:ASN:HD22	2.18	0.51
1:F:173:ILE:HD12	1:F:243:LEU:HG	1.92	0.51
1:F:201:ASP:O	1:F:210:PHE:CE2	2.62	0.51
1:F:690:LYS:O	1:F:691:LYS:C	2.52	0.51
2:P:16:PHE:CE1	2:P:27:ILE:CD1	2.94	0.51
2:Q:11:GLU:C	2:Q:13:LYS:N	2.69	0.51
2:R:18:LEU:HD13	2:R:19:PHE:CE1	2.45	0.51
2:R:36:MET:HE3	2:R:43:PRO:CG	2.32	0.51
2:S:68:PHE:C	2:S:70:THR:N	2.67	0.51
1:A:561:ASN:HA	1:A:564:VAL:HG22	1.92	0.51
1:A:678:VAL:HG13	1:A:745:TYR:HD2	1.73	0.51
1:B:694:VAL:HG22	2:P:18:LEU:HD21	1.90	0.51
1:C:238:GLN:O	1:C:240:ALA:N	2.44	0.51
1:C:666:ASN:HB2	1:C:748:TYR:OH	2.11	0.51
1:C:671:ARG:HD2	2:Q:14:GLU:HG2	1.91	0.51
1:D:297:LYS:HZ3	1:D:601:GLU:HB3	1.75	0.51
1:E:93:VAL:HG23	1:E:179:LEU:HD13	1.92	0.51
1:E:684:ASP:C	1:E:686:ASP:H	2.19	0.51
1:F:202:ASP:HB2	1:F:208:LEU:CD2	2.40	0.51
1:F:684:ASP:C	1:F:686:ASP:H	2.19	0.51
2:O:76:MET:C	2:O:78:ASP:H	2.19	0.51
2:P:133:ASP:OD2	2:P:135:GLN:HG3	2.10	0.51
2:Q:51:MET:HB2	2:Q:71:MET:HE2	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:R:58:ASP:C	2:R:60:ASN:N	2.65	0.51
2:T:9:ILE:CD1	2:T:69:LEU:HD11	2.39	0.51
1:A:77:ASP:O	1:A:81:GLN:CB	2.58	0.51
1:A:530:THR:HG21	2:O:145:MET:HE3	1.92	0.51
1:A:567:THR:CG2	1:A:568:GLY:H	2.23	0.51
1:A:684:ASP:C	1:A:686:ASP:H	2.18	0.51
1:B:397:GLU:O	1:B:479:LYS:HA	2.10	0.51
1:B:654:ILE:C	1:B:655:ASN:HD22	2.19	0.51
1:C:713:SER:O	1:C:714:GLN:C	2.51	0.51
1:D:142:VAL:HG22	1:D:154:ILE:HD12	1.92	0.51
1:D:217:LYS:HZ1	1:D:233:ASN:HB3	1.76	0.51
1:D:285:LYS:C	1:D:287:GLY:H	2.18	0.51
1:D:395:GLU:C	1:D:397:GLU:H	2.17	0.51
1:D:397:GLU:O	1:D:479:LYS:HA	2.10	0.51
1:E:730:ASN:O	1:E:732:ILE:N	2.44	0.51
1:E:736:LEU:HD21	1:E:750:GLN:CD	2.35	0.51
1:E:788:ASP:O	1:E:789:ASN:C	2.54	0.51
2:Q:120:GLU:HA	2:Q:123:GLN:HB2	1.93	0.51
2:Q:138:TYR:CE1	2:Q:142:VAL:HG22	2.45	0.51
2:S:13:LYS:HA	2:S:16:PHE:HB3	1.93	0.51
2:S:51:MET:HB2	2:S:71:MET:HE2	1.92	0.51
1:B:118:GLN:OE1	1:B:118:GLN:HA	2.09	0.51
1:B:635:ILE:H	1:B:635:ILE:HD12	1.75	0.51
1:C:214:PHE:CD1	1:C:218:LEU:HD23	2.46	0.51
1:C:478:ALA:CB	1:C:486:LYS:O	2.57	0.51
1:D:140:ARG:HA	1:D:140:ARG:HE	1.75	0.51
1:D:788:ASP:O	1:D:789:ASN:C	2.54	0.51
1:E:77:ASP:O	1:E:81:GLN:CB	2.56	0.51
1:E:202:ASP:HB2	1:E:208:LEU:CD2	2.41	0.51
2:O:117:THR:OG1	2:O:119:GLU:HG2	2.11	0.51
2:P:97:ASN:ND2	2:P:97:ASN:H	2.08	0.51
2:R:51:MET:HB2	2:R:71:MET:HE2	1.92	0.51
2:R:133:ASP:OD2	2:R:135:GLN:HG3	2.11	0.51
2:T:18:LEU:HD13	2:T:19:PHE:CE1	2.46	0.51
1:A:218:LEU:HD21	1:A:225:ILE:CD1	2.40	0.51
1:A:435:LEU:H	1:A:445:ARG:HA	1.75	0.51
1:A:579:THR:C	1:A:581:GLN:N	2.68	0.51
1:A:648:PRO:O	1:A:651:LYS:N	2.44	0.51
1:A:736:LEU:HD21	1:A:750:GLN:CD	2.36	0.51
1:B:628:PHE:CE2	2:P:90:ARG:CZ	2.94	0.51
1:D:197:LYS:NZ	1:D:197:LYS:CB	2.72	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:609:GLU:O	1:D:613:ARG:N	2.33	0.51
1:D:690:LYS:O	1:D:691:LYS:C	2.54	0.51
1:F:79:ILE:C	1:F:81:GLN:N	2.61	0.51
1:F:142:VAL:HA	1:F:154:ILE:HD13	1.92	0.51
1:F:483:GLY:C	1:F:484:VAL:HG22	2.36	0.51
1:F:540:ARG:HD3	1:F:627:TYR:CZ	2.46	0.51
2:P:9:ILE:CD1	2:P:69:LEU:HD11	2.39	0.51
2:P:64:ASP:OD2	2:P:67:GLU:HG3	2.11	0.51
1:A:372:LYS:HG3	1:A:373:LYS:HG2	1.93	0.51
1:A:583:ASN:ND2	1:A:587:PRO:HA	2.26	0.51
1:A:709:ASN:O	1:A:717:LYS:HE3	2.11	0.51
1:A:764:LEU:C	1:A:766:HIS:N	2.69	0.51
1:C:397:GLU:O	1:C:479:LYS:HA	2.10	0.51
1:C:632:TYR:O	1:C:633:ASN:HB2	2.11	0.51
1:C:764:LEU:C	1:C:766:HIS:N	2.69	0.51
1:D:202:ASP:HB2	1:D:208:LEU:CD2	2.40	0.51
1:D:478:ALA:CB	1:D:486:LYS:O	2.57	0.51
1:D:501:LEU:HD23	2:R:112:LEU:HD21	1.92	0.51
1:D:583:ASN:ND2	1:D:587:PRO:HA	2.26	0.51
1:E:199:LEU:O	1:E:201:ASP:N	2.33	0.51
1:E:579:THR:C	1:E:581:GLN:N	2.69	0.51
1:E:635:ILE:HD12	1:E:635:ILE:H	1.76	0.51
1:F:83:GLN:O	1:F:84:ASP:C	2.54	0.51
1:F:397:GLU:O	1:F:479:LYS:HA	2.11	0.51
2:O:17:SER:O	2:O:18:LEU:C	2.52	0.51
2:O:114:GLU:OE2	2:O:114:GLU:HA	2.11	0.51
2:P:18:LEU:HB3	2:P:19:PHE:CD1	2.46	0.51
2:P:97:ASN:C	2:P:99:TYR:H	2.19	0.51
2:Q:17:SER:O	2:Q:20:ASP:N	2.40	0.51
2:R:97:ASN:ND2	2:R:97:ASN:H	2.08	0.51
2:S:94:LYS:HE3	2:S:107:HIS:CD2	2.46	0.51
2:S:114:GLU:HA	2:S:114:GLU:OE2	2.10	0.51
2:S:120:GLU:HA	2:S:123:GLN:HB2	1.92	0.51
2:T:13:LYS:O	2:T:15:ALA:N	2.44	0.51
2:T:106:ARG:HB2	2:T:121:VAL:HG21	1.93	0.51
1:A:285:LYS:C	1:A:287:GLY:H	2.17	0.51
1:A:581:GLN:HE22	1:A:629:ASN:H	1.53	0.51
1:B:483:GLY:C	1:B:484:VAL:HG22	2.36	0.51
1:B:530:THR:HG21	2:P:145:MET:HE3	1.93	0.51
1:C:282:SER:HA	1:C:285:LYS:CD	2.39	0.51
1:C:525:LYS:O	1:C:529:VAL:HG23	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:115:LYS:HB2	1:E:118:GLN:HG2	1.87	0.51
1:E:142:VAL:HG22	1:E:154:ILE:CG2	2.41	0.51
1:E:730:ASN:C	1:E:732:ILE:N	2.67	0.51
1:F:515:LYS:NZ	1:F:515:LYS:HB3	2.26	0.51
1:F:565:LYS:C	1:F:567:THR:H	2.19	0.51
2:O:16:PHE:HA	2:O:35:VAL:HG11	1.93	0.51
2:R:11:GLU:C	2:R:13:LYS:N	2.69	0.51
2:R:106:ARG:HB2	2:R:121:VAL:HG21	1.93	0.51
2:R:117:THR:OG1	2:R:119:GLU:HG2	2.11	0.51
2:R:120:GLU:HA	2:R:123:GLN:HB2	1.93	0.51
2:R:121:VAL:C	2:R:123:GLN:N	2.67	0.51
2:S:18:LEU:HB3	2:S:19:PHE:CD1	2.46	0.51
2:T:51:MET:HB2	2:T:71:MET:HE2	1.93	0.51
1:A:85:LEU:O	1:A:88:LYS:HB3	2.11	0.50
1:A:130:SER:HB2	1:A:170:TYR:HE2	1.76	0.50
1:A:241:PHE:CE1	1:A:268:MET:HE1	2.45	0.50
1:A:379:ALA:O	1:A:383:GLY:N	2.43	0.50
1:B:225:ILE:HG12	1:B:229:PHE:CE2	2.46	0.50
1:B:372:LYS:HG3	1:B:373:LYS:HG2	1.93	0.50
1:B:443:GLU:O	1:B:455:TYR:HA	2.11	0.50
1:B:525:LYS:O	1:B:529:VAL:HG23	2.11	0.50
1:B:555:GLN:HG3	1:B:556:MET:N	2.25	0.50
1:D:170:TYR:O	1:D:174:GLY:N	2.44	0.50
1:D:173:ILE:C	1:D:175:LYS:N	2.66	0.50
1:D:372:LYS:HG3	1:D:373:LYS:HG2	1.94	0.50
1:D:700:TYR:HD1	1:D:727:GLN:C	2.18	0.50
1:E:118:GLN:OE1	1:E:118:GLN:HA	2.10	0.50
1:E:389:LYS:O	1:E:390:SER:C	2.54	0.50
1:E:764:LEU:C	1:E:766:HIS:N	2.68	0.50
1:F:99:GLU:C	1:F:101:GLY:N	2.61	0.50
1:F:112:VAL:O	1:F:114:HIS:N	2.38	0.50
1:F:282:SER:HA	1:F:285:LYS:CD	2.37	0.50
1:F:372:LYS:HG3	1:F:373:LYS:HG2	1.93	0.50
1:F:709:ASN:HB2	2:T:130:ILE:HG23	1.93	0.50
2:O:106:ARG:HB2	2:O:121:VAL:HG21	1.92	0.50
2:P:41:GLN:O	2:P:42:ASN:HB3	2.10	0.50
2:R:97:ASN:C	2:R:99:TYR:H	2.18	0.50
2:T:16:PHE:CE1	2:T:27:ILE:CD1	2.94	0.50
2:T:114:GLU:HA	2:T:114:GLU:OE2	2.10	0.50
1:A:161:ILE:CG2	1:A:168:GLU:HB2	2.41	0.50
1:A:788:ASP:O	1:A:789:ASN:C	2.53	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:327:LEU:HD12	1:B:327:LEU:N	2.25	0.50
1:B:515:LYS:NZ	1:B:515:LYS:HB3	2.26	0.50
1:B:609:GLU:O	1:B:612:GLY:N	2.44	0.50
1:C:83:GLN:O	1:C:84:ASP:C	2.54	0.50
1:C:170:TYR:O	1:C:174:GLY:N	2.45	0.50
1:D:666:ASN:HB2	1:D:748:TYR:OH	2.11	0.50
1:E:238:GLN:O	1:E:240:ALA:N	2.43	0.50
1:E:244:ALA:HB1	1:E:268:MET:HE3	1.93	0.50
1:E:397:GLU:O	1:E:479:LYS:HA	2.10	0.50
1:E:515:LYS:NZ	1:E:515:LYS:HB3	2.26	0.50
1:F:156:ILE:N	1:F:156:ILE:HD12	2.26	0.50
1:F:302:LEU:HB2	1:F:602:PHE:CD1	2.37	0.50
2:P:70:THR:C	2:P:72:MET:H	2.17	0.50
2:P:76:MET:C	2:P:78:ASP:H	2.19	0.50
2:S:17:SER:O	2:S:18:LEU:C	2.54	0.50
1:A:443:GLU:O	1:A:455:TYR:HA	2.11	0.50
1:B:105:TYR:HB2	1:B:153:ILE:HA	1.94	0.50
1:B:285:LYS:C	1:B:287:GLY:H	2.18	0.50
1:B:730:ASN:O	1:B:732:ILE:N	2.45	0.50
1:C:164:GLU:C	1:C:166:SER:H	2.19	0.50
1:C:216:GLU:HG3	1:C:217:LYS:N	2.27	0.50
1:C:310:GLU:O	1:C:314:ALA:HB2	2.12	0.50
1:C:395:GLU:C	1:C:397:GLU:H	2.17	0.50
1:C:633:ASN:O	1:C:642:TYR:HE1	1.93	0.50
1:D:565:LYS:C	1:D:567:THR:H	2.18	0.50
1:D:654:ILE:C	1:D:655:ASN:HD22	2.18	0.50
1:E:216:GLU:HG3	1:E:217:LYS:N	2.27	0.50
1:E:684:ASP:C	1:E:686:ASP:N	2.70	0.50
1:E:709:ASN:HB2	2:S:130:ILE:HG23	1.93	0.50
1:F:648:PRO:O	1:F:651:LYS:N	2.45	0.50
2:P:117:THR:OG1	2:P:119:GLU:HG2	2.12	0.50
2:Q:17:SER:O	2:Q:18:LEU:C	2.54	0.50
2:S:66:PRO:O	2:S:68:PHE:N	2.45	0.50
2:T:6:GLU:HG3	2:T:7:GLU:H	1.72	0.50
1:A:105:TYR:HB2	1:A:153:ILE:HA	1.94	0.50
1:A:318:ILE:CG2	1:A:322:LEU:HD12	2.42	0.50
1:B:216:GLU:HG3	1:B:217:LYS:N	2.27	0.50
1:B:295:VAL:O	1:B:296:LEU:C	2.55	0.50
1:B:318:ILE:CG2	1:B:322:LEU:HD12	2.41	0.50
1:C:85:LEU:O	1:C:88:LYS:HB3	2.11	0.50
1:C:583:ASN:ND2	1:C:587:PRO:HA	2.27	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:788:ASP:O	1:C:789:ASN:C	2.53	0.50
1:D:678:VAL:HG13	1:D:745:TYR:HD2	1.75	0.50
1:E:225:ILE:HG12	1:E:229:PHE:CE2	2.46	0.50
1:F:164:GLU:C	1:F:166:SER:H	2.19	0.50
1:F:214:PHE:CD1	1:F:218:LEU:HD23	2.46	0.50
1:F:389:LYS:O	1:F:390:SER:C	2.54	0.50
1:F:699:GLY:O	1:F:703:ASP:OD2	2.30	0.50
2:O:6:GLU:O	2:O:9:ILE:N	2.41	0.50
2:O:18:LEU:HD13	2:O:19:PHE:CE1	2.46	0.50
2:O:120:GLU:HA	2:O:123:GLN:HB2	1.92	0.50
2:Q:133:ASP:OD2	2:Q:135:GLN:HG3	2.12	0.50
1:A:156:ILE:N	1:A:156:ILE:HD12	2.27	0.50
1:A:201:ASP:O	1:A:210:PHE:HE2	1.94	0.50
1:B:173:ILE:C	1:B:175:LYS:N	2.67	0.50
1:B:579:THR:C	1:B:581:GLN:N	2.70	0.50
1:B:666:ASN:HB2	1:B:748:TYR:OH	2.11	0.50
1:C:140:ARG:HA	1:C:140:ARG:HE	1.76	0.50
1:D:97:TYR:HD2	1:D:102:GLY:HA3	1.77	0.50
1:D:238:GLN:O	1:D:240:ALA:N	2.44	0.50
1:D:298:GLY:C	1:D:300:LYS:N	2.66	0.50
1:D:736:LEU:HD21	1:D:750:GLN:CD	2.37	0.50
1:E:90:PRO:HG2	1:E:93:VAL:CG1	2.41	0.50
1:E:372:LYS:HG3	1:E:373:LYS:HG2	1.93	0.50
1:E:501:LEU:HD23	2:S:112:LEU:HD21	1.93	0.50
1:E:555:GLN:HG3	1:E:556:MET:N	2.25	0.50
1:E:694:VAL:HG22	2:S:18:LEU:HD21	1.91	0.50
1:F:509:PRO:HD2	1:F:536:TYR:CE2	2.47	0.50
1:F:654:ILE:C	1:F:655:ASN:HD22	2.19	0.50
1:F:666:ASN:HB2	1:F:748:TYR:OH	2.10	0.50
2:O:16:PHE:CE1	2:O:27:ILE:CD1	2.95	0.50
2:O:68:PHE:C	2:O:70:THR:N	2.69	0.50
2:P:8:GLN:O	2:P:12:PHE:CD2	2.65	0.50
2:P:18:LEU:HD13	2:P:19:PHE:CE1	2.46	0.50
2:P:51:MET:HB2	2:P:71:MET:HE2	1.94	0.50
2:P:106:ARG:HB2	2:P:121:VAL:HG21	1.93	0.50
2:Q:114:GLU:HA	2:Q:114:GLU:OE2	2.11	0.50
2:S:9:ILE:O	2:S:10:ALA:C	2.53	0.50
2:S:9:ILE:CD1	2:S:69:LEU:HD11	2.39	0.50
2:S:27:ILE:HB	2:S:31:GLU:HB3	1.94	0.50
2:T:27:ILE:HB	2:T:31:GLU:HB3	1.94	0.50
1:A:225:ILE:HG12	1:A:229:PHE:CE2	2.46	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:238:GLN:O	1:A:240:ALA:N	2.45	0.50
1:A:363:TYR:O	1:A:365:PRO:HD3	2.12	0.50
1:A:540:ARG:HD3	1:A:627:TYR:CZ	2.46	0.50
1:A:555:GLN:CG	1:A:556:MET:N	2.75	0.50
1:B:142:VAL:HG22	1:B:154:ILE:CG2	2.41	0.50
1:B:363:TYR:O	1:B:365:PRO:HD3	2.12	0.50
1:B:764:LEU:C	1:B:766:HIS:N	2.66	0.50
1:C:244:ALA:HB1	1:C:268:MET:HE3	1.93	0.50
1:C:515:LYS:NZ	1:C:515:LYS:HB3	2.26	0.50
1:C:540:ARG:HD3	1:C:627:TYR:CZ	2.46	0.50
1:C:565:LYS:C	1:C:567:THR:H	2.19	0.50
1:D:310:GLU:O	1:D:314:ALA:HB2	2.12	0.50
1:E:112:VAL:O	1:E:114:HIS:N	2.40	0.50
1:E:164:GLU:C	1:E:166:SER:H	2.19	0.50
1:E:188:LEU:CD2	1:E:188:LEU:N	2.47	0.50
1:E:217:LYS:HZ1	1:E:233:ASN:HB3	1.76	0.50
1:E:628:PHE:CD1	1:E:645:TRP:CD1	2.99	0.50
1:F:176:GLY:C	1:F:178:SER:N	2.70	0.50
1:F:500:SER:HG	1:F:502:THR:HG1	1.60	0.50
1:F:632:TYR:O	1:F:633:ASN:HB2	2.12	0.50
2:O:9:ILE:CD1	2:O:69:LEU:HD11	2.38	0.50
2:P:13:LYS:HA	2:P:16:PHE:HB3	1.94	0.50
2:Q:13:LYS:HA	2:Q:16:PHE:HB3	1.94	0.50
2:R:9:ILE:O	2:R:10:ALA:C	2.55	0.50
2:R:25:GLY:HA3	2:R:65:PHE:CZ	2.47	0.50
2:R:27:ILE:HB	2:R:31:GLU:HB3	1.94	0.50
2:S:117:THR:OG1	2:S:119:GLU:HG2	2.11	0.50
1:B:90:PRO:HG2	1:B:93:VAL:CG1	2.42	0.50
1:B:238:GLN:O	1:B:240:ALA:N	2.45	0.50
1:B:736:LEU:HD21	1:B:750:GLN:CD	2.37	0.50
1:C:201:ASP:O	1:C:210:PHE:HE2	1.95	0.50
1:C:275:GLY:HA2	1:C:278:LYS:CD	2.42	0.50
1:C:389:LYS:O	1:C:390:SER:C	2.53	0.50
1:C:736:LEU:HD21	1:C:750:GLN:CD	2.36	0.50
1:D:684:ASP:C	1:D:686:ASP:H	2.18	0.50
1:D:730:ASN:O	1:D:732:ILE:N	2.45	0.50
1:E:327:LEU:HD12	1:E:327:LEU:N	2.26	0.50
1:E:656:THR:O	1:E:755:ARG:NH1	2.43	0.50
1:E:659:THR:O	1:E:660:SER:C	2.55	0.50
1:F:135:VAL:N	1:F:136:PRO:CD	2.75	0.50
1:F:217:LYS:HZ1	1:F:233:ASN:HB3	1.76	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:628:PHE:HD1	1:F:632:TYR:HD1	1.60	0.50
1:F:788:ASP:O	1:F:789:ASN:C	2.53	0.50
2:P:49:GLN:HB3	2:P:53:ASN:HB2	1.93	0.50
2:Q:68:PHE:C	2:Q:70:THR:N	2.69	0.50
2:R:49:GLN:HB3	2:R:53:ASN:HB2	1.93	0.50
1:A:310:GLU:O	1:A:314:ALA:HB2	2.11	0.50
1:A:345:THR:HG21	1:A:491:ASP:HA	1.94	0.50
1:A:409:ARG:O	1:A:412:GLU:HB3	2.12	0.50
1:B:156:ILE:HD12	1:B:156:ILE:N	2.27	0.50
1:B:270:LYS:HA	1:B:273:LYS:CD	2.41	0.50
1:B:500:SER:HG	1:B:502:THR:HG1	1.58	0.50
1:C:387:ASN:O	1:C:390:SER:HB2	2.12	0.50
1:C:416:ASN:HB3	1:C:418:ILE:HG12	1.94	0.50
1:D:118:GLN:HA	1:D:118:GLN:OE1	2.11	0.50
1:D:189:ASP:C	1:D:191:GLU:N	2.70	0.50
1:D:327:LEU:HD12	1:D:327:LEU:N	2.26	0.50
1:D:363:TYR:O	1:D:365:PRO:HD3	2.12	0.50
1:D:509:PRO:HD2	1:D:536:TYR:CE2	2.46	0.50
1:E:115:LYS:HD3	1:E:153:ILE:HD13	1.93	0.50
1:E:218:LEU:HD21	1:E:225:ILE:CD1	2.42	0.50
1:E:379:ALA:O	1:E:383:GLY:N	2.43	0.50
1:E:648:PRO:O	1:E:651:LYS:N	2.45	0.50
1:F:238:GLN:O	1:F:240:ALA:N	2.45	0.50
1:F:525:LYS:O	1:F:529:VAL:HG23	2.11	0.50
1:F:555:GLN:CG	1:F:556:MET:N	2.75	0.50
1:F:579:THR:C	1:F:581:GLN:N	2.70	0.50
2:O:8:GLN:O	2:O:12:PHE:CD2	2.65	0.50
2:O:13:LYS:HA	2:O:16:PHE:HB3	1.94	0.50
2:O:25:GLY:HA3	2:O:65:PHE:CZ	2.47	0.50
2:O:97:ASN:C	2:O:99:TYR:H	2.19	0.50
2:P:114:GLU:HA	2:P:114:GLU:OE2	2.11	0.50
2:Q:27:ILE:HB	2:Q:31:GLU:HB3	1.94	0.50
2:R:13:LYS:HA	2:R:16:PHE:HB3	1.94	0.50
2:R:16:PHE:CE1	2:R:27:ILE:CD1	2.95	0.50
2:R:66:PRO:O	2:R:68:PHE:N	2.45	0.50
2:R:76:MET:C	2:R:78:ASP:H	2.20	0.50
1:A:164:GLU:C	1:A:166:SER:H	2.19	0.50
1:B:241:PHE:CE1	1:B:268:MET:HE1	2.47	0.50
1:B:244:ALA:HB1	1:B:268:MET:HE3	1.94	0.50
1:B:597:ASN:CB	1:B:598:PRO:HD2	2.36	0.50
1:D:164:GLU:C	1:D:166:SER:H	2.19	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:216:GLU:HG3	1:D:217:LYS:N	2.27	0.50
1:D:659:THR:H	1:D:662:GLU:HB3	1.75	0.50
1:E:561:ASN:HA	1:E:564:VAL:HG22	1.92	0.50
1:F:105:TYR:HB2	1:F:153:ILE:HA	1.93	0.50
1:F:628:PHE:CD1	1:F:645:TRP:CD1	3.00	0.50
2:O:11:GLU:C	2:O:13:LYS:N	2.69	0.50
2:S:88:ALA:O	2:S:91:VAL:HB	2.11	0.50
2:T:13:LYS:HA	2:T:16:PHE:HB3	1.94	0.50
2:T:16:PHE:HA	2:T:35:VAL:HG11	1.94	0.50
1:A:115:LYS:O	1:A:117:LEU:N	2.45	0.49
1:A:483:GLY:C	1:A:484:VAL:HG22	2.37	0.49
1:A:716:LYS:O	1:A:717:LYS:C	2.55	0.49
1:B:218:LEU:HD21	1:B:225:ILE:CD1	2.41	0.49
1:B:389:LYS:O	1:B:390:SER:C	2.55	0.49
1:C:130:SER:HB2	1:C:170:TYR:HE2	1.77	0.49
1:C:173:ILE:HG13	1:C:242:SER:CB	2.34	0.49
1:C:483:GLY:C	1:C:484:VAL:HG22	2.36	0.49
1:D:105:TYR:HB2	1:D:153:ILE:HA	1.94	0.49
1:D:153:ILE:C	1:D:154:ILE:HG12	2.37	0.49
1:D:295:VAL:O	1:D:296:LEU:C	2.55	0.49
1:E:105:TYR:HB2	1:E:153:ILE:HA	1.94	0.49
1:E:115:LYS:O	1:E:117:LEU:N	2.45	0.49
1:E:270:LYS:HA	1:E:273:LYS:CD	2.41	0.49
1:E:318:ILE:CG2	1:E:322:LEU:HD12	2.41	0.49
1:E:481:VAL:O	1:E:482:GLU:HB2	2.12	0.49
1:F:764:LEU:C	1:F:766:HIS:N	2.70	0.49
2:O:49:GLN:O	2:O:53:ASN:N	2.40	0.49
2:P:11:GLU:C	2:P:13:LYS:N	2.69	0.49
2:Q:49:GLN:HB3	2:Q:53:ASN:HB2	1.93	0.49
2:Q:76:MET:C	2:Q:78:ASP:H	2.19	0.49
2:R:17:SER:O	2:R:18:LEU:C	2.54	0.49
1:B:357:TRP:CZ2	1:B:370:LEU:HD23	2.47	0.49
1:B:379:ALA:O	1:B:383:GLY:N	2.44	0.49
1:B:684:ASP:C	1:B:686:ASP:N	2.70	0.49
1:C:105:TYR:HB2	1:C:153:ILE:HA	1.94	0.49
1:C:173:ILE:C	1:C:175:LYS:N	2.67	0.49
1:D:318:ILE:CG2	1:D:322:LEU:HD12	2.42	0.49
1:D:710:HIS:C	1:D:712:PHE:N	2.59	0.49
1:E:310:GLU:O	1:E:314:ALA:HB2	2.12	0.49
1:E:344:ALA:HA	1:E:569:TYR:OH	2.12	0.49
1:E:357:TRP:CZ2	1:E:370:LEU:HD23	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:363:TYR:O	1:E:365:PRO:HD3	2.13	0.49
1:E:483:GLY:C	1:E:484:VAL:HG22	2.38	0.49
1:E:517:VAL:HG23	1:E:518:ASN:HD22	1.77	0.49
1:E:567:THR:CG2	1:E:568:GLY:H	2.24	0.49
1:F:318:ILE:CG2	1:F:322:LEU:HD12	2.42	0.49
2:O:18:LEU:HB3	2:O:19:PHE:CD1	2.48	0.49
2:O:97:ASN:ND2	2:O:97:ASN:H	2.08	0.49
2:O:133:ASP:OD2	2:O:135:GLN:HG3	2.12	0.49
2:Q:36:MET:HE3	2:Q:43:PRO:CG	2.31	0.49
2:S:6:GLU:O	2:S:9:ILE:N	2.42	0.49
2:S:25:GLY:HA3	2:S:65:PHE:CZ	2.47	0.49
2:T:133:ASP:OD2	2:T:135:GLN:HG3	2.12	0.49
1:A:175:LYS:O	1:A:178:SER:N	2.43	0.49
1:A:244:ALA:HB1	1:A:268:MET:HE3	1.95	0.49
1:A:666:ASN:HB2	1:A:748:TYR:OH	2.12	0.49
1:B:481:VAL:O	1:B:482:GLU:HB2	2.12	0.49
1:B:540:ARG:HD3	1:B:627:TYR:CZ	2.47	0.49
1:B:632:TYR:O	1:B:633:ASN:HB2	2.12	0.49
1:B:776:LEU:HD23	1:B:776:LEU:C	2.37	0.49
1:C:89:ILE:HG22	1:C:90:PRO:HD2	1.93	0.49
1:C:175:LYS:O	1:C:178:SER:N	2.44	0.49
1:C:176:GLY:C	1:C:178:SER:N	2.70	0.49
1:C:625:LEU:HD12	1:C:625:LEU:C	2.37	0.49
1:C:628:PHE:HD1	1:C:632:TYR:HD1	1.60	0.49
1:D:90:PRO:HG2	1:D:93:VAL:CB	2.41	0.49
1:D:225:ILE:HG12	1:D:229:PHE:CE2	2.47	0.49
1:D:351:HIS:HB2	1:D:386:GLU:CG	2.43	0.49
1:D:540:ARG:HD3	1:D:627:TYR:CZ	2.47	0.49
1:F:161:ILE:CG2	1:F:168:GLU:HB2	2.41	0.49
1:F:170:TYR:O	1:F:174:GLY:N	2.46	0.49
1:F:225:ILE:HG12	1:F:229:PHE:CE2	2.46	0.49
1:F:310:GLU:O	1:F:314:ALA:HB2	2.12	0.49
1:F:344:ALA:HA	1:F:569:TYR:OH	2.13	0.49
1:F:653:LYS:O	1:F:655:ASN:N	2.45	0.49
1:F:659:THR:O	1:F:660:SER:C	2.55	0.49
2:O:66:PRO:O	2:O:68:PHE:N	2.45	0.49
2:O:93:ASP:OD1	2:O:97:ASN:ND2	2.46	0.49
2:P:9:ILE:O	2:P:10:ALA:C	2.54	0.49
2:Q:6:GLU:HG3	2:Q:7:GLU:H	1.74	0.49
2:Q:93:ASP:OD1	2:Q:97:ASN:ND2	2.45	0.49
2:S:18:LEU:HD13	2:S:19:PHE:CE1	2.48	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:T:6:GLU:O	2:T:9:ILE:N	2.41	0.49
1:A:153:ILE:C	1:A:154:ILE:HG12	2.36	0.49
1:A:189:ASP:C	1:A:191:GLU:N	2.70	0.49
1:A:217:LYS:HZ1	1:A:233:ASN:HB3	1.77	0.49
1:A:327:LEU:HD12	1:A:327:LEU:N	2.26	0.49
1:B:164:GLU:C	1:B:166:SER:H	2.19	0.49
1:B:275:GLY:HA2	1:B:278:LYS:CD	2.42	0.49
1:B:523:LEU:HD11	2:P:144:MET:HG3	1.95	0.49
1:B:697:ILE:HG23	1:B:732:ILE:HD11	1.94	0.49
1:C:142:VAL:HG22	1:C:154:ILE:CG2	2.41	0.49
1:C:153:ILE:C	1:C:154:ILE:HG12	2.38	0.49
1:C:225:ILE:HG12	1:C:229:PHE:CE2	2.46	0.49
1:C:409:ARG:O	1:C:412:GLU:HB3	2.13	0.49
1:C:653:LYS:O	1:C:655:ASN:N	2.46	0.49
1:C:776:LEU:HD23	1:C:776:LEU:C	2.38	0.49
1:D:244:ALA:HB1	1:D:268:MET:HE3	1.95	0.49
1:D:481:VAL:O	1:D:482:GLU:HB2	2.12	0.49
1:E:295:VAL:O	1:E:296:LEU:C	2.56	0.49
1:E:716:LYS:O	1:E:717:LYS:C	2.55	0.49
1:F:363:TYR:O	1:F:365:PRO:HD3	2.12	0.49
2:O:9:ILE:O	2:O:10:ALA:C	2.55	0.49
2:O:121:VAL:C	2:O:123:GLN:N	2.67	0.49
2:P:68:PHE:C	2:P:70:THR:N	2.68	0.49
2:Q:18:LEU:HB3	2:Q:19:PHE:CD1	2.47	0.49
2:R:8:GLN:O	2:R:12:PHE:CD2	2.65	0.49
2:S:49:GLN:HB3	2:S:53:ASN:HB2	1.93	0.49
2:T:88:ALA:O	2:T:91:VAL:HB	2.12	0.49
2:T:117:THR:OG1	2:T:119:GLU:HG2	2.11	0.49
1:A:217:LYS:HZ1	1:A:236:GLU:HB2	1.77	0.49
1:A:295:VAL:O	1:A:296:LEU:C	2.55	0.49
1:A:565:LYS:C	1:A:567:THR:H	2.19	0.49
1:A:748:TYR:O	1:A:751:TYR:N	2.46	0.49
1:B:115:LYS:O	1:B:117:LEU:N	2.45	0.49
1:B:286:GLU:O	1:B:290:LYS:HB2	2.13	0.49
1:B:709:ASN:HB2	2:P:130:ILE:HG23	1.95	0.49
1:B:747:ASN:O	1:B:750:GLN:HB2	2.12	0.49
1:C:90:PRO:HG2	1:C:93:VAL:CG1	2.41	0.49
1:C:199:LEU:C	1:C:201:ASP:N	2.70	0.49
1:C:296:LEU:HD23	1:C:296:LEU:N	2.28	0.49
1:C:363:TYR:O	1:C:365:PRO:HD3	2.11	0.49
1:C:555:GLN:CG	1:C:556:MET:N	2.75	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:241:PHE:CE1	1:D:268:MET:HE1	2.47	0.49
1:D:275:GLY:HA2	1:D:278:LYS:CD	2.43	0.49
1:D:409:ARG:O	1:D:412:GLU:HB3	2.13	0.49
1:D:441:VAL:HG11	1:D:462:ILE:O	2.13	0.49
1:D:662:GLU:OE2	1:D:755:ARG:NH2	2.41	0.49
1:E:176:GLY:C	1:E:178:SER:N	2.70	0.49
1:F:216:GLU:HG3	1:F:217:LYS:N	2.27	0.49
1:F:345:THR:HG21	1:F:491:ASP:HA	1.94	0.49
1:F:376:GLN:O	1:F:378:LEU:N	2.46	0.49
2:Q:25:GLY:HA3	2:Q:65:PHE:CZ	2.47	0.49
2:R:17:SER:O	2:R:20:ASP:N	2.42	0.49
2:T:58:ASP:C	2:T:60:ASN:N	2.65	0.49
1:A:170:TYR:O	1:A:174:GLY:N	2.45	0.49
1:A:210:PHE:C	1:A:210:PHE:CD1	2.91	0.49
1:A:509:PRO:HD2	1:A:536:TYR:CE2	2.46	0.49
1:A:515:LYS:NZ	1:A:515:LYS:HB3	2.27	0.49
1:B:136:PRO:O	1:B:137:PHE:C	2.56	0.49
1:B:217:LYS:HZ1	1:B:233:ASN:HB3	1.76	0.49
1:B:478:ALA:CB	1:B:486:LYS:O	2.57	0.49
1:B:497:LEU:CD1	1:B:556:MET:HG2	2.40	0.49
1:B:711:ILE:C	1:B:712:PHE:HD2	2.20	0.49
1:C:123:GLU:O	1:C:146:LYS:NZ	2.45	0.49
1:C:379:ALA:O	1:C:383:GLY:N	2.44	0.49
1:D:109:ILE:HD13	1:D:157:LYS:HZ3	1.78	0.49
1:D:345:THR:HG21	1:D:491:ASP:HA	1.95	0.49
1:D:448:ASP:O	1:D:449:GLU:C	2.56	0.49
1:D:530:THR:HG21	2:R:145:MET:HE3	1.94	0.49
1:D:581:GLN:HE22	1:D:629:ASN:H	1.52	0.49
1:D:627:TYR:CD1	1:D:627:TYR:C	2.91	0.49
1:D:632:TYR:O	1:D:633:ASN:HB2	2.12	0.49
1:D:764:LEU:C	1:D:766:HIS:N	2.70	0.49
1:E:170:TYR:O	1:E:174:GLY:N	2.46	0.49
1:E:282:SER:HA	1:E:285:LYS:CD	2.38	0.49
1:F:244:ALA:HB1	1:F:268:MET:HE3	1.94	0.49
2:O:17:SER:O	2:O:20:ASP:N	2.40	0.49
2:O:49:GLN:HB3	2:O:53:ASN:HB2	1.94	0.49
2:O:70:THR:C	2:O:72:MET:H	2.19	0.49
2:P:66:PRO:O	2:P:68:PHE:N	2.46	0.49
2:S:133:ASP:OD2	2:S:135:GLN:HG3	2.12	0.49
2:T:25:GLY:HA3	2:T:65:PHE:CZ	2.47	0.49
2:T:76:MET:C	2:T:78:ASP:H	2.19	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:184:LYS:HE2	1:A:193:LEU:CD1	2.31	0.49
1:A:286:GLU:O	1:A:290:LYS:HB2	2.13	0.49
1:A:344:ALA:HA	1:A:569:TYR:OH	2.13	0.49
1:A:387:ASN:O	1:A:390:SER:HB2	2.12	0.49
1:A:552:TRP:O	1:A:553:GLN:C	2.55	0.49
1:A:690:LYS:O	1:A:691:LYS:C	2.55	0.49
1:B:109:ILE:HD13	1:B:157:LYS:HZ3	1.76	0.49
1:B:310:GLU:O	1:B:314:ALA:HB2	2.12	0.49
1:B:565:LYS:C	1:B:567:THR:H	2.20	0.49
1:B:716:LYS:O	1:B:717:LYS:C	2.55	0.49
1:B:764:LEU:O	1:B:766:HIS:N	2.45	0.49
1:C:210:PHE:C	1:C:210:PHE:CD1	2.91	0.49
1:C:318:ILE:CG2	1:C:322:LEU:HD12	2.42	0.49
1:C:376:GLN:O	1:C:378:LEU:N	2.46	0.49
1:C:523:LEU:HD11	2:Q:144:MET:HG3	1.94	0.49
1:C:716:LYS:O	1:C:717:LYS:C	2.55	0.49
1:D:123:GLU:O	1:D:146:LYS:NZ	2.46	0.49
1:D:210:PHE:C	1:D:210:PHE:CD1	2.91	0.49
1:D:357:TRP:CZ2	1:D:370:LEU:HD23	2.47	0.49
1:D:387:ASN:O	1:D:390:SER:HB2	2.13	0.49
1:D:555:GLN:CG	1:D:556:MET:N	2.75	0.49
1:D:776:LEU:HD23	1:D:776:LEU:C	2.38	0.49
1:E:345:THR:HG21	1:E:491:ASP:HA	1.94	0.49
1:E:376:GLN:O	1:E:378:LEU:N	2.45	0.49
1:E:552:TRP:O	1:E:553:GLN:C	2.56	0.49
1:E:719:LYS:O	1:E:721:SER:N	2.46	0.49
1:E:747:ASN:O	1:E:750:GLN:HB2	2.13	0.49
1:F:295:VAL:O	1:F:296:LEU:C	2.55	0.49
1:F:351:HIS:HB2	1:F:386:GLU:CG	2.43	0.49
1:F:530:THR:HG21	2:T:145:MET:HE3	1.95	0.49
1:F:633:ASN:O	1:F:642:TYR:HE1	1.95	0.49
1:F:719:LYS:O	1:F:721:SER:N	2.46	0.49
2:O:51:MET:HB2	2:O:71:MET:HE2	1.94	0.49
2:Q:16:PHE:CE1	2:Q:27:ILE:CD1	2.94	0.49
2:S:97:ASN:ND2	2:S:97:ASN:H	2.10	0.49
2:T:9:ILE:O	2:T:10:ALA:C	2.54	0.49
2:T:17:SER:O	2:T:19:PHE:N	2.45	0.49
2:T:120:GLU:HA	2:T:123:GLN:HB2	1.93	0.49
1:A:135:VAL:N	1:A:136:PRO:CD	2.75	0.49
1:A:176:GLY:C	1:A:178:SER:N	2.71	0.49
1:A:199:LEU:C	1:A:201:ASP:N	2.70	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:322:LEU:O	1:A:324:THR:HG22	2.12	0.49
1:A:351:HIS:HB2	1:A:386:GLU:CG	2.43	0.49
1:A:632:TYR:O	1:A:633:ASN:HB2	2.12	0.49
1:B:176:GLY:C	1:B:178:SER:N	2.70	0.49
1:B:225:ILE:HA	1:B:229:PHE:CE2	2.47	0.49
1:B:408:LEU:HD12	1:B:408:LEU:N	2.02	0.49
1:B:509:PRO:HD2	1:B:536:TYR:CE2	2.47	0.49
1:B:517:VAL:HG23	1:B:518:ASN:HD22	1.77	0.49
1:B:555:GLN:CG	1:B:556:MET:N	2.76	0.49
1:B:581:GLN:HE22	1:B:629:ASN:H	1.52	0.49
1:C:220:LEU:HD11	1:C:223:LYS:HB2	1.95	0.49
1:C:241:PHE:CE1	1:C:268:MET:HE1	2.47	0.49
1:C:351:HIS:HB2	1:C:386:GLU:CG	2.43	0.49
1:D:136:PRO:O	1:D:137:PHE:C	2.55	0.49
1:D:142:VAL:HG22	1:D:154:ILE:CG2	2.42	0.49
1:D:161:ILE:CG2	1:D:168:GLU:HB2	2.42	0.49
1:D:214:PHE:CD1	1:D:218:LEU:HD23	2.48	0.49
1:D:709:ASN:HB2	2:R:130:ILE:HG23	1.94	0.49
1:E:89:ILE:HG22	1:E:90:PRO:HD2	1.94	0.49
1:E:210:PHE:C	1:E:210:PHE:CD1	2.91	0.49
1:E:530:THR:HG21	2:S:145:MET:HE3	1.95	0.49
1:E:555:GLN:CG	1:E:556:MET:N	2.76	0.49
1:F:275:GLY:HA2	1:F:278:LYS:CD	2.43	0.49
1:F:416:ASN:HB3	1:F:418:ILE:HG12	1.95	0.49
2:P:17:SER:O	2:P:18:LEU:C	2.54	0.49
2:Q:70:THR:C	2:Q:72:MET:N	2.71	0.49
2:T:18:LEU:HB3	2:T:19:PHE:CD1	2.47	0.49
1:A:297:LYS:HZ1	1:A:601:GLU:HB3	1.75	0.49
1:A:441:VAL:HG11	1:A:462:ILE:O	2.13	0.49
1:A:610:MET:O	1:A:614:PHE:N	2.40	0.49
1:A:755:ARG:O	1:A:756:ILE:C	2.56	0.49
1:B:73:ASN:C	1:B:74:GLU:O	2.56	0.49
1:B:344:ALA:HA	1:B:569:TYR:OH	2.13	0.49
1:B:441:VAL:HG11	1:B:462:ILE:O	2.12	0.49
1:B:690:LYS:O	1:B:691:LYS:C	2.54	0.49
1:C:156:ILE:N	1:C:156:ILE:HD12	2.28	0.49
1:C:244:ALA:CB	1:C:268:MET:HE3	2.43	0.49
1:C:270:LYS:HA	1:C:273:LYS:CD	2.41	0.49
1:C:295:VAL:O	1:C:296:LEU:C	2.56	0.49
1:C:344:ALA:HA	1:C:569:TYR:OH	2.13	0.49
1:C:448:ASP:O	1:C:449:GLU:C	2.56	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:659:THR:O	1:C:660:SER:C	2.56	0.49
1:C:755:ARG:O	1:C:756:ILE:C	2.56	0.49
1:D:189:ASP:HB3	1:D:190:PRO:CD	2.43	0.49
1:D:483:GLY:C	1:D:484:VAL:HG22	2.37	0.49
1:E:156:ILE:HD12	1:E:156:ILE:N	2.28	0.49
1:E:191:GLU:O	1:E:193:LEU:N	2.46	0.49
1:E:776:LEU:HD23	1:E:776:LEU:C	2.38	0.49
1:F:123:GLU:O	1:F:146:LYS:NZ	2.45	0.49
1:F:479:LYS:HG2	1:F:488:LEU:CD2	2.37	0.49
1:F:628:PHE:CE2	2:T:90:ARG:CZ	2.95	0.49
2:R:18:LEU:HB3	2:R:19:PHE:CD1	2.48	0.49
2:T:66:PRO:O	2:T:68:PHE:N	2.45	0.49
2:T:68:PHE:C	2:T:70:THR:N	2.68	0.49
2:T:138:TYR:CE1	2:T:142:VAL:CG2	2.96	0.49
1:B:162:ASN:HA	1:B:165:GLN:HG3	1.95	0.49
1:B:164:GLU:C	1:B:166:SER:N	2.71	0.49
1:C:115:LYS:O	1:C:117:LEU:N	2.46	0.49
1:C:189:ASP:C	1:C:191:GLU:N	2.71	0.49
1:D:217:LYS:HZ1	1:D:236:GLU:HB2	1.77	0.49
1:D:635:ILE:H	1:D:635:ILE:HD12	1.77	0.49
1:D:719:LYS:O	1:D:721:SER:N	2.45	0.49
1:E:136:PRO:O	1:E:137:PHE:C	2.56	0.49
1:E:416:ASN:HB3	1:E:418:ILE:HG12	1.94	0.49
1:F:115:LYS:HD3	1:F:153:ILE:HD13	1.93	0.49
1:F:115:LYS:O	1:F:117:LEU:N	2.46	0.49
1:F:217:LYS:CB	1:F:236:GLU:CD	2.86	0.49
1:F:297:LYS:CB	1:F:297:LYS:HZ2	2.26	0.49
1:F:414:LYS:C	1:F:416:ASN:H	2.21	0.49
1:F:448:ASP:O	1:F:449:GLU:C	2.56	0.49
1:F:730:ASN:C	1:F:732:ILE:H	2.20	0.49
2:O:27:ILE:HB	2:O:31:GLU:HB3	1.94	0.49
2:P:70:THR:C	2:P:72:MET:N	2.71	0.49
2:R:6:GLU:O	2:R:9:ILE:N	2.42	0.49
2:S:8:GLN:O	2:S:12:PHE:CD2	2.66	0.49
1:A:244:ALA:CB	1:A:268:MET:HE3	2.43	0.48
1:A:628:PHE:CE2	2:O:90:ARG:CZ	2.96	0.48
1:A:684:ASP:C	1:A:686:ASP:N	2.70	0.48
1:B:210:PHE:C	1:B:210:PHE:CD1	2.91	0.48
1:B:387:ASN:O	1:B:390:SER:HB2	2.12	0.48
1:B:658:PRO:HG2	1:B:701:LEU:HD22	1.95	0.48
1:C:115:LYS:HD3	1:C:153:ILE:HD13	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:162:ASN:HA	1:C:165:GLN:HG3	1.95	0.48
1:C:530:THR:HG21	2:Q:145:MET:HE3	1.95	0.48
1:C:684:ASP:C	1:C:686:ASP:N	2.70	0.48
1:C:765:THR:HA	1:C:769:SER:HB2	1.95	0.48
1:D:162:ASN:HA	1:D:165:GLN:HG3	1.95	0.48
1:D:716:LYS:O	1:D:717:LYS:C	2.56	0.48
1:E:162:ASN:HA	1:E:165:GLN:HG3	1.95	0.48
1:E:186:LYS:CD	1:E:234:LEU:HD12	2.43	0.48
1:E:462:ILE:CG1	1:E:463:THR:N	2.74	0.48
1:E:697:ILE:HG23	1:E:732:ILE:HD11	1.95	0.48
1:F:89:ILE:HG22	1:F:90:PRO:HD2	1.94	0.48
1:F:90:PRO:HG2	1:F:93:VAL:CG1	2.43	0.48
1:F:357:TRP:CZ2	1:F:370:LEU:HD23	2.47	0.48
1:F:379:ALA:O	1:F:383:GLY:N	2.44	0.48
1:F:409:ARG:O	1:F:412:GLU:HB3	2.13	0.48
1:F:412:GLU:C	1:F:414:LYS:H	2.22	0.48
1:F:711:ILE:C	1:F:712:PHE:HD2	2.21	0.48
2:O:88:ALA:O	2:O:91:VAL:HB	2.13	0.48
1:A:90:PRO:HG2	1:A:93:VAL:CG1	2.43	0.48
1:A:297:LYS:HB3	1:A:297:LYS:HZ3	1.78	0.48
1:A:389:LYS:O	1:A:390:SER:C	2.54	0.48
1:A:633:ASN:O	1:A:642:TYR:HE1	1.96	0.48
1:B:405:LEU:HD13	1:B:453:VAL:CG2	2.43	0.48
1:B:765:THR:HA	1:B:769:SER:HB2	1.96	0.48
1:C:97:TYR:HD2	1:C:102:GLY:HA3	1.79	0.48
1:C:135:VAL:N	1:C:136:PRO:CD	2.76	0.48
1:C:136:PRO:O	1:C:137:PHE:C	2.56	0.48
1:C:191:GLU:O	1:C:193:LEU:N	2.46	0.48
1:C:357:TRP:CZ2	1:C:370:LEU:HD23	2.47	0.48
1:C:627:TYR:CD1	1:C:627:TYR:C	2.91	0.48
1:C:680:LYS:HG2	1:C:681:ASP:N	2.28	0.48
1:D:83:GLN:O	1:D:84:ASP:C	2.55	0.48
1:D:164:GLU:C	1:D:166:SER:N	2.71	0.48
1:D:201:ASP:O	1:D:210:PHE:HE2	1.96	0.48
1:D:389:LYS:O	1:D:390:SER:C	2.53	0.48
1:D:416:ASN:HB3	1:D:418:ILE:HG12	1.95	0.48
1:D:691:LYS:O	1:D:692:GLU:C	2.56	0.48
1:E:100:LEU:O	1:E:150:PRO:HD2	2.13	0.48
1:E:210:PHE:C	1:E:212:GLN:N	2.71	0.48
1:E:302:LEU:HB2	1:E:602:PHE:CD1	2.37	0.48
1:E:509:PRO:HD2	1:E:536:TYR:CE2	2.48	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:711:ILE:C	1:E:712:PHE:HD2	2.21	0.48
1:F:353:LYS:HG2	1:F:368:GLN:HE22	1.77	0.48
1:F:387:ASN:O	1:F:390:SER:HB2	2.12	0.48
1:F:556:MET:O	1:F:560:LEU:HG	2.13	0.48
1:F:736:LEU:HD21	1:F:750:GLN:CD	2.37	0.48
2:R:13:LYS:HE2	2:R:65:PHE:HB3	1.95	0.48
2:R:138:TYR:O	2:R:141:PHE:HB3	2.13	0.48
1:A:353:LYS:HG2	1:A:368:GLN:HE22	1.78	0.48
1:A:448:ASP:O	1:A:449:GLU:C	2.56	0.48
1:A:747:ASN:O	1:A:750:GLN:HB2	2.14	0.48
1:B:220:LEU:HD11	1:B:223:LYS:HB2	1.95	0.48
1:C:217:LYS:CB	1:C:236:GLU:CD	2.86	0.48
1:C:387:ASN:N	1:C:387:ASN:ND2	2.50	0.48
1:C:509:PRO:HD2	1:C:536:TYR:CE2	2.48	0.48
1:C:552:TRP:O	1:C:553:GLN:C	2.55	0.48
1:C:667:LEU:O	1:C:669:SER:N	2.47	0.48
1:C:719:LYS:O	1:C:721:SER:N	2.46	0.48
1:D:115:LYS:O	1:D:117:LEU:N	2.46	0.48
1:D:135:VAL:N	1:D:136:PRO:CD	2.76	0.48
1:D:579:THR:C	1:D:581:GLN:N	2.70	0.48
1:D:628:PHE:HD1	1:D:632:TYR:HD1	1.60	0.48
1:D:659:THR:O	1:D:660:SER:C	2.56	0.48
1:D:755:ARG:O	1:D:756:ILE:C	2.57	0.48
1:E:197:LYS:HB3	1:E:197:LYS:HZ2	1.76	0.48
1:E:217:LYS:HZ1	1:E:236:GLU:HB2	1.78	0.48
1:E:244:ALA:CB	1:E:268:MET:HE3	2.43	0.48
1:E:275:GLY:HA2	1:E:278:LYS:CD	2.43	0.48
1:E:414:LYS:C	1:E:416:ASN:H	2.21	0.48
1:E:764:LEU:O	1:E:766:HIS:N	2.47	0.48
1:F:162:ASN:HA	1:F:165:GLN:HG3	1.95	0.48
1:F:189:ASP:C	1:F:191:GLU:N	2.71	0.48
1:F:288:VAL:O	1:F:292:ARG:HG2	2.14	0.48
1:F:481:VAL:O	1:F:482:GLU:HB2	2.12	0.48
1:F:517:VAL:HG23	1:F:518:ASN:HD22	1.78	0.48
1:F:627:TYR:CD1	1:F:627:TYR:C	2.90	0.48
1:F:635:ILE:H	1:F:635:ILE:HD12	1.78	0.48
2:T:70:THR:C	2:T:72:MET:N	2.71	0.48
1:A:164:GLU:C	1:A:166:SER:N	2.72	0.48
1:A:216:GLU:HG3	1:A:217:LYS:N	2.27	0.48
1:A:275:GLY:HA2	1:A:278:LYS:CD	2.43	0.48
1:A:525:LYS:O	1:A:529:VAL:HG23	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:214:PHE:CD1	1:B:218:LEU:HD23	2.46	0.48
1:B:217:LYS:HZ1	1:B:236:GLU:HB2	1.79	0.48
1:B:648:PRO:O	1:B:651:LYS:N	2.45	0.48
1:C:95:GLU:O	1:C:99:GLU:HB2	2.14	0.48
1:C:186:LYS:CD	1:C:234:LEU:HD12	2.43	0.48
1:C:288:VAL:O	1:C:292:ARG:HG2	2.13	0.48
1:C:517:VAL:HG23	1:C:518:ASN:HD22	1.77	0.48
1:D:73:ASN:C	1:D:74:GLU:O	2.56	0.48
1:D:130:SER:HB2	1:D:170:TYR:HE2	1.76	0.48
1:D:156:ILE:HD12	1:D:156:ILE:N	2.29	0.48
1:D:176:GLY:C	1:D:178:SER:N	2.70	0.48
1:D:220:LEU:HD11	1:D:223:LYS:HB2	1.95	0.48
1:D:249:PHE:O	1:D:250:ALA:C	2.57	0.48
1:D:517:VAL:HG23	1:D:518:ASN:HD22	1.77	0.48
1:D:525:LYS:O	1:D:529:VAL:HG23	2.13	0.48
1:D:653:LYS:O	1:D:655:ASN:N	2.46	0.48
1:D:684:ASP:C	1:D:686:ASP:N	2.69	0.48
1:D:765:THR:HA	1:D:769:SER:HB2	1.95	0.48
1:E:153:ILE:C	1:E:154:ILE:HG12	2.37	0.48
1:F:270:LYS:HA	1:F:273:LYS:CD	2.42	0.48
1:F:353:LYS:HG2	1:F:368:GLN:NE2	2.29	0.48
1:F:552:TRP:O	1:F:553:GLN:C	2.54	0.48
1:F:691:LYS:NZ	2:T:20:ASP:O	2.43	0.48
1:F:746:LYS:HD2	1:F:747:ASN:ND2	2.29	0.48
1:F:776:LEU:HD23	1:F:776:LEU:C	2.37	0.48
2:O:13:LYS:HE2	2:O:65:PHE:HB3	1.95	0.48
2:P:27:ILE:HB	2:P:31:GLU:HB3	1.95	0.48
2:P:88:ALA:O	2:P:91:VAL:HB	2.14	0.48
2:Q:9:ILE:CD1	2:Q:69:LEU:HD11	2.39	0.48
2:Q:38:SER:C	2:Q:40:GLY:H	2.20	0.48
2:Q:138:TYR:O	2:Q:141:PHE:HB3	2.14	0.48
2:R:138:TYR:CE1	2:R:142:VAL:CG2	2.96	0.48
2:S:76:MET:C	2:S:78:ASP:H	2.20	0.48
1:A:136:PRO:O	1:A:137:PHE:C	2.56	0.48
1:A:142:VAL:HG22	1:A:154:ILE:CG2	2.42	0.48
1:A:375:GLY:HA2	1:A:464:VAL:HG11	1.96	0.48
1:A:635:ILE:H	1:A:635:ILE:HD12	1.77	0.48
1:A:765:THR:HA	1:A:769:SER:HB2	1.96	0.48
1:B:189:ASP:C	1:B:191:GLU:N	2.71	0.48
1:B:345:THR:HG21	1:B:491:ASP:HA	1.94	0.48
1:B:448:ASP:O	1:B:449:GLU:C	2.56	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:77:ASP:O	1:C:81:GLN:CB	2.59	0.48
1:C:161:ILE:CG2	1:C:168:GLU:HB2	2.42	0.48
1:D:112:VAL:O	1:D:114:HIS:N	2.39	0.48
1:D:191:GLU:O	1:D:193:LEU:N	2.46	0.48
1:D:648:PRO:O	1:D:651:LYS:N	2.46	0.48
1:D:794:GLN:HE21	1:D:794:GLN:HB3	1.53	0.48
1:E:115:LYS:HB3	1:E:118:GLN:HG2	1.95	0.48
1:E:441:VAL:HG11	1:E:462:ILE:O	2.14	0.48
1:F:241:PHE:CE1	1:F:268:MET:HE1	2.48	0.48
1:F:755:ARG:O	1:F:756:ILE:C	2.56	0.48
1:F:765:THR:HA	1:F:769:SER:HB2	1.96	0.48
2:O:27:ILE:HD12	2:O:32:LEU:HA	1.95	0.48
2:P:16:PHE:CD1	2:P:27:ILE:HD13	2.49	0.48
2:P:25:GLY:HA3	2:P:65:PHE:CZ	2.48	0.48
2:P:49:GLN:O	2:P:53:ASN:N	2.41	0.48
2:S:11:GLU:C	2:S:13:LYS:N	2.69	0.48
2:T:101:SER:H	2:T:104:GLU:CG	2.27	0.48
1:A:131:ARG:HG3	1:A:243:LEU:HD22	1.96	0.48
1:A:162:ASN:HA	1:A:165:GLN:HG3	1.96	0.48
1:A:191:GLU:O	1:A:193:LEU:N	2.46	0.48
1:A:376:GLN:O	1:A:378:LEU:N	2.46	0.48
1:A:445:ARG:HG2	1:A:471:TRP:CZ3	2.48	0.48
1:A:462:ILE:HD11	1:A:466:GLY:HA2	1.95	0.48
1:A:746:LYS:HD2	1:A:747:ASN:ND2	2.29	0.48
1:B:95:GLU:O	1:B:99:GLU:HB2	2.14	0.48
1:B:234:LEU:HD23	1:B:235:THR:N	2.27	0.48
1:B:351:HIS:HB2	1:B:386:GLU:CG	2.43	0.48
1:B:351:HIS:HB2	1:B:386:GLU:HG3	1.96	0.48
1:B:409:ARG:O	1:B:412:GLU:HB3	2.14	0.48
1:C:164:GLU:C	1:C:166:SER:N	2.71	0.48
1:C:414:LYS:C	1:C:416:ASN:H	2.21	0.48
1:C:462:ILE:HD11	1:C:466:GLY:HA2	1.95	0.48
1:C:690:LYS:O	1:C:691:LYS:C	2.53	0.48
1:D:376:GLN:O	1:D:378:LEU:N	2.46	0.48
1:D:552:TRP:O	1:D:553:GLN:C	2.55	0.48
1:D:658:PRO:HG3	1:D:752:LEU:HD22	1.94	0.48
1:E:217:LYS:CB	1:E:236:GLU:CD	2.87	0.48
1:E:241:PHE:CE1	1:E:268:MET:HE1	2.48	0.48
1:E:351:HIS:HB2	1:E:386:GLU:CG	2.43	0.48
1:E:375:GLY:HA2	1:E:464:VAL:HG11	1.96	0.48
1:E:755:ARG:O	1:E:756:ILE:C	2.56	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:77:ASP:O	1:F:81:GLN:CB	2.59	0.48
1:F:100:LEU:O	1:F:150:PRO:HD2	2.13	0.48
1:F:130:SER:HB2	1:F:170:TYR:HE2	1.78	0.48
1:F:441:VAL:HG11	1:F:462:ILE:O	2.14	0.48
1:F:462:ILE:HD11	1:F:466:GLY:HA2	1.95	0.48
1:F:478:ALA:CB	1:F:486:LYS:O	2.58	0.48
2:O:138:TYR:CE1	2:O:142:VAL:CG2	2.96	0.48
2:R:70:THR:C	2:R:72:MET:N	2.71	0.48
2:R:106:ARG:NH2	2:R:118:ASP:OD2	2.47	0.48
1:A:214:PHE:CD1	1:A:218:LEU:HD23	2.48	0.48
1:A:416:ASN:HB3	1:A:418:ILE:HG12	1.95	0.48
1:A:481:VAL:O	1:A:482:GLU:HB2	2.13	0.48
1:B:414:LYS:C	1:B:416:ASN:H	2.21	0.48
1:B:556:MET:O	1:B:560:LEU:HG	2.14	0.48
1:C:345:THR:HG21	1:C:491:ASP:HA	1.95	0.48
1:C:441:VAL:HG11	1:C:462:ILE:O	2.13	0.48
1:C:730:ASN:C	1:C:732:ILE:H	2.21	0.48
1:D:282:SER:HA	1:D:285:LYS:CD	2.39	0.48
1:D:414:LYS:C	1:D:416:ASN:H	2.21	0.48
1:D:747:ASN:O	1:D:750:GLN:HB2	2.14	0.48
1:E:73:ASN:C	1:E:74:GLU:O	2.56	0.48
1:E:225:ILE:HA	1:E:229:PHE:CE2	2.47	0.48
1:E:387:ASN:O	1:E:390:SER:HB2	2.13	0.48
1:E:409:ARG:O	1:E:412:GLU:HB3	2.13	0.48
1:E:556:MET:O	1:E:560:LEU:HG	2.13	0.48
1:E:625:LEU:HD12	1:E:625:LEU:C	2.39	0.48
1:E:632:TYR:O	1:E:633:ASN:HB2	2.13	0.48
1:F:210:PHE:C	1:F:210:PHE:CD1	2.91	0.48
1:F:220:LEU:HD11	1:F:223:LYS:HB2	1.96	0.48
1:F:747:ASN:O	1:F:750:GLN:HB2	2.14	0.48
2:Q:9:ILE:O	2:Q:10:ALA:C	2.54	0.48
1:A:220:LEU:HD11	1:A:223:LYS:HB2	1.96	0.48
1:A:296:LEU:HD23	1:A:296:LEU:N	2.28	0.48
1:B:153:ILE:C	1:B:154:ILE:HG12	2.38	0.48
1:C:186:LYS:HE3	1:C:234:LEU:CD1	2.44	0.48
1:C:217:LYS:NZ	1:C:236:GLU:HB2	2.29	0.48
1:C:234:LEU:HD23	1:C:235:THR:N	2.27	0.48
1:C:286:GLU:O	1:C:290:LYS:HB2	2.13	0.48
1:C:434:LEU:HD22	1:C:445:ARG:HB3	1.96	0.48
1:D:180:ASP:O	1:D:183:SER:N	2.36	0.48
1:D:218:LEU:HG	1:D:218:LEU:O	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:344:ALA:HA	1:D:569:TYR:OH	2.14	0.48
1:E:159:TYR:HD1	1:E:159:TYR:N	2.10	0.48
1:E:448:ASP:O	1:E:449:GLU:C	2.56	0.48
1:E:765:THR:HA	1:E:769:SER:HB2	1.96	0.48
2:P:126:ARG:HG3	2:P:126:ARG:NH2	2.28	0.48
2:R:3:GLN:N	2:R:77:LYS:HD3	2.29	0.48
2:S:27:ILE:HD12	2:S:32:LEU:HA	1.96	0.48
2:T:8:GLN:O	2:T:12:PHE:CD2	2.66	0.48
2:T:27:ILE:HD12	2:T:32:LEU:HA	1.96	0.48
1:A:112:VAL:O	1:A:114:HIS:N	2.39	0.48
1:A:217:LYS:CB	1:A:236:GLU:CD	2.87	0.48
1:A:357:TRP:CZ2	1:A:370:LEU:HD23	2.48	0.48
1:A:639:ASN:ND2	1:A:639:ASN:N	2.62	0.48
1:B:170:TYR:O	1:B:174:GLY:N	2.46	0.48
1:B:376:GLN:O	1:B:378:LEU:N	2.46	0.48
1:B:746:LYS:HD2	1:B:747:ASN:ND2	2.29	0.48
1:C:658:PRO:HG3	1:C:752:LEU:HD22	1.94	0.48
1:C:709:ASN:HB2	2:Q:130:ILE:HG23	1.95	0.48
1:C:747:ASN:O	1:C:750:GLN:HB2	2.14	0.48
1:D:217:LYS:NZ	1:D:236:GLU:HB2	2.29	0.48
1:D:391:ILE:O	1:D:392:THR:C	2.57	0.48
1:D:412:GLU:C	1:D:414:LYS:H	2.22	0.48
1:D:434:LEU:HD22	1:D:445:ARG:HB3	1.95	0.48
1:D:633:ASN:O	1:D:642:TYR:HE1	1.96	0.48
1:D:746:LYS:HD2	1:D:747:ASN:ND2	2.29	0.48
1:E:83:GLN:O	1:E:84:ASP:C	2.56	0.48
1:E:690:LYS:O	1:E:691:LYS:C	2.55	0.48
1:E:730:ASN:C	1:E:732:ILE:H	2.22	0.48
1:E:797:ILE:O	1:E:797:ILE:HG13	2.14	0.48
1:F:217:LYS:NZ	1:F:236:GLU:HB2	2.29	0.48
1:F:329:ARG:HB3	1:F:330:PRO:CD	2.44	0.48
1:F:497:LEU:CD1	1:F:556:MET:HG2	2.40	0.48
2:P:27:ILE:HD12	2:P:32:LEU:HA	1.96	0.48
2:S:16:PHE:CD1	2:S:27:ILE:HD13	2.49	0.48
1:A:462:ILE:CG1	1:A:463:THR:N	2.75	0.48
1:A:517:VAL:HG23	1:A:518:ASN:HD22	1.78	0.48
1:A:523:LEU:HD11	2:O:144:MET:HG3	1.95	0.48
1:A:687:GLU:O	1:A:690:LYS:N	2.47	0.48
1:A:711:ILE:C	1:A:712:PHE:HD2	2.21	0.48
1:A:730:ASN:C	1:A:732:ILE:H	2.22	0.48
1:B:416:ASN:HB3	1:B:418:ILE:HG12	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:678:VAL:HG13	1:B:745:TYR:HD2	1.75	0.48
1:B:687:GLU:O	1:B:690:LYS:N	2.47	0.48
1:B:691:LYS:O	1:B:692:GLU:C	2.56	0.48
1:C:218:LEU:HG	1:C:218:LEU:O	2.14	0.48
1:D:170:TYR:HA	1:D:173:ILE:HG22	1.95	0.48
1:D:217:LYS:CB	1:D:236:GLU:CD	2.87	0.48
1:D:296:LEU:HD23	1:D:296:LEU:N	2.29	0.48
1:D:353:LYS:HG2	1:D:368:GLN:HE22	1.78	0.48
1:D:559:ARG:HG3	1:D:559:ARG:HH11	1.79	0.48
1:E:189:ASP:C	1:E:191:GLU:N	2.71	0.48
1:F:88:LYS:NZ	1:F:172:GLU:CD	2.72	0.48
1:F:136:PRO:O	1:F:137:PHE:C	2.56	0.48
1:F:159:TYR:HD1	1:F:159:TYR:N	2.10	0.48
2:O:16:PHE:CD1	2:O:27:ILE:HD13	2.49	0.48
2:O:66:PRO:C	2:O:68:PHE:N	2.71	0.48
2:O:138:TYR:O	2:O:141:PHE:HB3	2.14	0.48
2:Q:30:LYS:HD3	2:Q:30:LYS:N	2.12	0.48
2:Q:70:THR:O	2:Q:72:MET:N	2.47	0.48
2:T:49:GLN:HB3	2:T:53:ASN:HB2	1.94	0.48
1:A:327:LEU:HG	1:A:595:ILE:HG12	1.96	0.47
1:A:671:ARG:NH1	1:A:671:ARG:HG3	2.29	0.47
1:B:89:ILE:HG22	1:B:90:PRO:HD2	1.95	0.47
1:B:218:LEU:O	1:B:218:LEU:HG	2.14	0.47
1:B:680:LYS:HG2	1:B:681:ASP:N	2.29	0.47
1:B:797:ILE:HG13	1:B:797:ILE:O	2.14	0.47
1:C:130:SER:O	1:C:131:ARG:C	2.57	0.47
1:D:100:LEU:O	1:D:150:PRO:HD2	2.14	0.47
1:D:479:LYS:HG2	1:D:488:LEU:CD2	2.37	0.47
1:D:748:TYR:O	1:D:751:TYR:N	2.47	0.47
1:E:424:LYS:HB3	1:E:424:LYS:NZ	2.27	0.47
1:E:628:PHE:CE2	2:S:90:ARG:CZ	2.97	0.47
1:E:628:PHE:HD1	1:E:632:TYR:HD1	1.61	0.47
1:F:523:LEU:HD11	2:T:144:MET:HG3	1.96	0.47
1:F:671:ARG:NH1	1:F:671:ARG:HG3	2.29	0.47
1:F:748:TYR:O	1:F:751:TYR:N	2.46	0.47
2:R:68:PHE:C	2:R:70:THR:N	2.68	0.47
2:R:70:THR:O	2:R:72:MET:N	2.47	0.47
2:S:70:THR:O	2:S:72:MET:N	2.47	0.47
1:A:83:GLN:O	1:A:84:ASP:C	2.56	0.47
1:A:89:ILE:HG22	1:A:93:VAL:CG1	2.32	0.47
1:A:170:TYR:HA	1:A:173:ILE:HG22	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:414:LYS:C	1:A:416:ASN:H	2.21	0.47
1:A:625:LEU:HD12	1:A:625:LEU:C	2.38	0.47
1:A:627:TYR:CD1	1:A:627:TYR:C	2.92	0.47
1:A:709:ASN:HB2	2:O:130:ILE:HG23	1.96	0.47
1:B:83:GLN:O	1:B:84:ASP:C	2.55	0.47
1:B:375:GLY:HA2	1:B:464:VAL:HG11	1.96	0.47
1:B:462:ILE:HD11	1:B:466:GLY:HA2	1.95	0.47
1:B:627:TYR:C	1:B:627:TYR:CD1	2.92	0.47
1:B:653:LYS:O	1:B:655:ASN:N	2.46	0.47
1:C:249:PHE:O	1:C:250:ALA:C	2.58	0.47
1:C:479:LYS:HG2	1:C:488:LEU:CD2	2.36	0.47
1:D:270:LYS:HA	1:D:273:LYS:CD	2.41	0.47
1:D:351:HIS:HB2	1:D:386:GLU:HG3	1.95	0.47
1:D:523:LEU:HD11	2:R:144:MET:HG3	1.95	0.47
1:E:164:GLU:C	1:E:166:SER:N	2.72	0.47
1:E:329:ARG:HB3	1:E:330:PRO:CD	2.44	0.47
1:E:711:ILE:O	1:E:712:PHE:HD2	1.97	0.47
1:F:95:GLU:O	1:F:99:GLU:HB2	2.14	0.47
1:F:186:LYS:CD	1:F:234:LEU:HD12	2.44	0.47
1:F:244:ALA:CB	1:F:268:MET:HE3	2.44	0.47
1:F:286:GLU:O	1:F:290:LYS:HB2	2.13	0.47
1:F:629:ASN:HB3	1:F:632:TYR:CZ	2.50	0.47
2:P:106:ARG:NH2	2:P:118:ASP:OD2	2.47	0.47
2:Q:8:GLN:O	2:Q:12:PHE:CD2	2.66	0.47
2:Q:13:LYS:HE2	2:Q:65:PHE:HB3	1.95	0.47
2:Q:51:MET:CB	2:Q:71:MET:HE2	2.44	0.47
2:R:19:PHE:CD1	2:R:19:PHE:N	2.83	0.47
2:S:138:TYR:CE1	2:S:142:VAL:CG2	2.96	0.47
1:A:186:LYS:HE3	1:A:234:LEU:CD1	2.44	0.47
1:A:297:LYS:HZ3	1:A:601:GLU:HB3	1.79	0.47
1:A:412:GLU:C	1:A:414:LYS:H	2.21	0.47
1:A:658:PRO:HG3	1:A:752:LEU:HD22	1.96	0.47
1:B:135:VAL:HG22	1:B:135:VAL:O	2.14	0.47
1:B:730:ASN:C	1:B:732:ILE:H	2.23	0.47
1:C:135:VAL:O	1:C:135:VAL:HG22	2.14	0.47
1:C:329:ARG:HB3	1:C:330:PRO:CD	2.44	0.47
1:C:375:GLY:HA2	1:C:464:VAL:HG11	1.96	0.47
1:D:175:LYS:HB2	1:D:175:LYS:HZ2	1.78	0.47
1:D:244:ALA:CB	1:D:268:MET:HE3	2.44	0.47
1:D:711:ILE:C	1:D:712:PHE:HD2	2.22	0.47
1:E:201:ASP:O	1:E:210:PHE:HE2	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:581:GLN:HE22	1:E:629:ASN:H	1.53	0.47
1:E:671:ARG:NH1	1:E:671:ARG:HG3	2.29	0.47
1:F:135:VAL:O	1:F:135:VAL:HG22	2.14	0.47
1:F:218:LEU:HG	1:F:218:LEU:O	2.14	0.47
1:F:225:ILE:HA	1:F:229:PHE:CE2	2.47	0.47
1:F:322:LEU:O	1:F:324:THR:HG22	2.14	0.47
2:S:66:PRO:C	2:S:68:PHE:N	2.71	0.47
1:A:90:PRO:C	1:A:92:ASP:N	2.72	0.47
1:A:180:ASP:O	1:A:183:SER:N	2.36	0.47
1:A:189:ASP:HB3	1:A:190:PRO:CD	2.44	0.47
1:A:434:LEU:HD22	1:A:445:ARG:HB3	1.96	0.47
1:A:628:PHE:HD1	1:A:632:TYR:HD1	1.61	0.47
1:A:691:LYS:O	1:A:692:GLU:C	2.57	0.47
1:C:100:LEU:O	1:C:150:PRO:HD2	2.13	0.47
1:C:372:LYS:O	1:C:374:HIS:N	2.47	0.47
1:C:746:LYS:HD2	1:C:747:ASN:ND2	2.29	0.47
1:D:199:LEU:C	1:D:201:ASP:N	2.70	0.47
1:D:375:GLY:HA2	1:D:464:VAL:HG11	1.96	0.47
1:D:697:ILE:HG23	1:D:732:ILE:HD11	1.95	0.47
1:E:457:THR:CG2	1:E:469:PHE:H	2.27	0.47
1:E:627:TYR:C	1:E:627:TYR:CD1	2.92	0.47
1:F:170:TYR:HA	1:F:173:ILE:HG22	1.96	0.47
1:F:375:GLY:HA2	1:F:464:VAL:HG11	1.96	0.47
1:F:797:ILE:O	1:F:797:ILE:HG13	2.15	0.47
2:O:38:SER:C	2:O:40:GLY:H	2.22	0.47
2:Q:88:ALA:O	2:Q:91:VAL:HB	2.14	0.47
2:R:81:SER:O	2:R:82:GLU:C	2.58	0.47
2:S:70:THR:C	2:S:72:MET:N	2.71	0.47
1:A:100:LEU:O	1:A:150:PRO:HD2	2.14	0.47
1:B:189:ASP:HB3	1:B:190:PRO:CD	2.45	0.47
1:B:191:GLU:O	1:B:193:LEU:N	2.46	0.47
1:B:217:LYS:CB	1:B:236:GLU:CD	2.87	0.47
1:B:296:LEU:HD23	1:B:296:LEU:N	2.27	0.47
1:B:353:LYS:HG2	1:B:368:GLN:HE22	1.79	0.47
1:B:755:ARG:O	1:B:756:ILE:C	2.57	0.47
1:C:115:LYS:HE3	1:C:116:GLU:HG2	1.95	0.47
1:C:170:TYR:HA	1:C:173:ILE:HG22	1.95	0.47
1:C:556:MET:O	1:C:560:LEU:HG	2.14	0.47
1:D:90:PRO:HG2	1:D:93:VAL:CG1	2.44	0.47
1:D:90:PRO:C	1:D:92:ASP:H	2.23	0.47
1:D:171:TYR:O	1:D:175:LYS:NZ	2.48	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:286:GLU:O	1:D:290:LYS:HB2	2.14	0.47
1:D:405:LEU:HD13	1:D:453:VAL:CG2	2.44	0.47
1:D:515:LYS:HB3	1:D:515:LYS:NZ	2.29	0.47
1:D:667:LEU:O	1:D:669:SER:N	2.48	0.47
1:D:687:GLU:O	1:D:690:LYS:N	2.47	0.47
1:E:173:ILE:C	1:E:175:LYS:N	2.67	0.47
1:E:296:LEU:HD23	1:E:296:LEU:N	2.29	0.47
1:E:658:PRO:HG3	1:E:752:LEU:HD22	1.97	0.47
1:F:112:VAL:C	1:F:114:HIS:N	2.73	0.47
1:F:201:ASP:O	1:F:210:PHE:HE2	1.97	0.47
1:F:296:LEU:HD23	1:F:296:LEU:N	2.29	0.47
1:F:658:PRO:HG3	1:F:752:LEU:HD22	1.96	0.47
1:F:667:LEU:O	1:F:669:SER:N	2.48	0.47
2:P:66:PRO:C	2:P:68:PHE:N	2.72	0.47
2:Q:13:LYS:O	2:Q:15:ALA:N	2.48	0.47
2:R:13:LYS:O	2:R:15:ALA:N	2.48	0.47
2:S:51:MET:CB	2:S:71:MET:HE2	2.45	0.47
1:A:90:PRO:C	1:A:92:ASP:H	2.23	0.47
1:A:318:ILE:H	1:A:318:ILE:CD1	2.23	0.47
1:A:329:ARG:HB3	1:A:330:PRO:CD	2.44	0.47
1:A:351:HIS:HB2	1:A:386:GLU:HG3	1.97	0.47
1:A:628:PHE:CD1	1:A:645:TRP:CD1	3.02	0.47
1:B:73:ASN:O	1:B:74:GLU:C	2.56	0.47
1:B:170:TYR:HA	1:B:173:ILE:HG22	1.95	0.47
1:B:186:LYS:CD	1:B:234:LEU:HD12	2.45	0.47
1:B:201:ASP:O	1:B:210:PHE:HE2	1.96	0.47
1:B:559:ARG:HH11	1:B:559:ARG:HG3	1.80	0.47
1:B:633:ASN:O	1:B:642:TYR:HE1	1.97	0.47
1:C:210:PHE:C	1:C:210:PHE:HD1	2.23	0.47
1:C:687:GLU:O	1:C:690:LYS:N	2.47	0.47
1:C:711:ILE:C	1:C:712:PHE:HD2	2.22	0.47
1:D:77:ASP:O	1:D:81:GLN:CB	2.58	0.47
1:D:671:ARG:NH1	1:D:671:ARG:HG3	2.29	0.47
1:E:90:PRO:C	1:E:92:ASP:N	2.72	0.47
1:E:180:ASP:O	1:E:183:SER:N	2.35	0.47
1:E:186:LYS:HE3	1:E:234:LEU:CD1	2.44	0.47
1:E:299:GLU:HA	1:E:302:LEU:HB3	1.97	0.47
1:E:353:LYS:HG2	1:E:368:GLN:HE22	1.78	0.47
1:E:667:LEU:O	1:E:669:SER:N	2.48	0.47
1:F:73:ASN:C	1:F:74:GLU:O	2.56	0.47
1:F:189:ASP:HB3	1:F:190:PRO:CD	2.45	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:299:GLU:HA	1:F:302:LEU:HB3	1.96	0.47
2:O:117:THR:O	2:O:119:GLU:N	2.48	0.47
2:Q:66:PRO:O	2:Q:68:PHE:N	2.47	0.47
2:T:3:GLN:N	2:T:77:LYS:HD3	2.30	0.47
2:T:16:PHE:CD1	2:T:27:ILE:HD13	2.49	0.47
2:T:70:THR:O	2:T:72:MET:N	2.47	0.47
2:T:93:ASP:OD1	2:T:97:ASN:ND2	2.47	0.47
2:T:117:THR:O	2:T:119:GLU:N	2.48	0.47
1:A:88:LYS:NZ	1:A:172:GLU:CD	2.71	0.47
1:A:234:LEU:HD23	1:A:235:THR:N	2.28	0.47
1:A:299:GLU:HA	1:A:302:LEU:HB3	1.97	0.47
1:A:308:VAL:CG2	1:A:336:THR:O	2.63	0.47
1:A:431:LYS:O	1:A:432:TYR:HD2	1.98	0.47
1:A:443:GLU:CD	1:A:458:LYS:HG2	2.40	0.47
1:B:90:PRO:C	1:B:92:ASP:N	2.73	0.47
1:B:299:GLU:HA	1:B:302:LEU:HB3	1.97	0.47
1:B:329:ARG:HB3	1:B:330:PRO:CD	2.45	0.47
1:B:365:PRO:O	1:B:366:PHE:C	2.58	0.47
1:B:431:LYS:O	1:B:432:TYR:HD2	1.97	0.47
1:B:443:GLU:CD	1:B:458:LYS:HG2	2.40	0.47
1:B:445:ARG:HG2	1:B:471:TRP:CZ3	2.49	0.47
1:B:628:PHE:HD1	1:B:632:TYR:HD1	1.61	0.47
1:B:776:LEU:O	1:B:780:LEU:CD2	2.63	0.47
1:B:794:GLN:HE21	1:B:794:GLN:HB3	1.53	0.47
1:C:175:LYS:HB2	1:C:175:LYS:HZ2	1.79	0.47
1:C:351:HIS:HB2	1:C:386:GLU:HG3	1.96	0.47
1:C:405:LEU:N	1:C:405:LEU:CD1	2.78	0.47
1:C:412:GLU:C	1:C:414:LYS:H	2.22	0.47
1:C:443:GLU:CD	1:C:458:LYS:HG2	2.40	0.47
1:C:457:THR:CG2	1:C:469:PHE:H	2.27	0.47
1:C:554:LYS:O	1:C:557:LEU:N	2.48	0.47
1:C:628:PHE:CE2	2:Q:90:ARG:CZ	2.97	0.47
1:D:115:LYS:HE3	1:D:116:GLU:HG2	1.95	0.47
1:D:337:ASN:C	1:D:339:ILE:N	2.72	0.47
1:D:353:LYS:HG2	1:D:368:GLN:NE2	2.30	0.47
1:D:372:LYS:O	1:D:374:HIS:N	2.47	0.47
1:D:444:PHE:CD2	1:D:455:TYR:HB3	2.50	0.47
1:D:680:LYS:HG2	1:D:681:ASP:N	2.30	0.47
1:E:210:PHE:C	1:E:210:PHE:HD1	2.23	0.47
1:E:286:GLU:O	1:E:290:LYS:HB2	2.14	0.47
1:E:499:PRO:CG	1:E:504:ILE:HD11	2.44	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:523:LEU:HD11	2:S:144:MET:HG3	1.96	0.47
1:F:164:GLU:C	1:F:166:SER:N	2.71	0.47
1:F:191:GLU:O	1:F:193:LEU:N	2.47	0.47
1:F:210:PHE:C	1:F:210:PHE:HD1	2.23	0.47
1:F:654:ILE:HG22	1:F:655:ASN:ND2	2.29	0.47
1:F:680:LYS:HG2	1:F:681:ASP:N	2.29	0.47
1:F:691:LYS:O	1:F:692:GLU:C	2.57	0.47
2:O:48:LEU:HA	2:O:51:MET:CE	2.41	0.47
2:O:117:THR:C	2:O:119:GLU:N	2.73	0.47
2:P:6:GLU:O	2:P:9:ILE:N	2.42	0.47
2:P:19:PHE:CD1	2:P:19:PHE:N	2.83	0.47
2:P:70:THR:O	2:P:72:MET:N	2.47	0.47
2:P:121:VAL:C	2:P:123:GLN:H	2.22	0.47
2:P:138:TYR:CE1	2:P:142:VAL:CG2	2.98	0.47
2:Q:3:GLN:N	2:Q:77:LYS:HD3	2.30	0.47
2:Q:16:PHE:CD1	2:Q:27:ILE:HD13	2.49	0.47
2:Q:81:SER:O	2:Q:82:GLU:C	2.56	0.47
2:Q:117:THR:O	2:Q:119:GLU:N	2.48	0.47
2:R:27:ILE:HD12	2:R:32:LEU:HA	1.96	0.47
2:R:49:GLN:O	2:R:53:ASN:N	2.41	0.47
2:R:66:PRO:C	2:R:68:PHE:N	2.71	0.47
2:S:132:GLY:C	2:S:134:GLY:H	2.22	0.47
2:T:13:LYS:HE2	2:T:65:PHE:HB3	1.95	0.47
2:T:89:PHE:HB2	2:T:141:PHE:CE2	2.50	0.47
1:A:171:TYR:O	1:A:175:LYS:NZ	2.48	0.47
1:A:186:LYS:CD	1:A:234:LEU:HD12	2.45	0.47
1:A:405:LEU:HD13	1:A:453:VAL:CG2	2.44	0.47
1:A:559:ARG:HH11	1:A:559:ARG:HG3	1.80	0.47
1:A:719:LYS:O	1:A:721:SER:N	2.48	0.47
1:B:244:ALA:CB	1:B:268:MET:HE3	2.44	0.47
1:B:639:ASN:ND2	1:B:639:ASN:N	2.61	0.47
1:B:719:LYS:O	1:B:721:SER:N	2.48	0.47
1:C:225:ILE:HA	1:C:229:PHE:CE2	2.47	0.47
1:C:308:VAL:CG2	1:C:336:THR:O	2.62	0.47
1:C:353:LYS:HG2	1:C:368:GLN:NE2	2.30	0.47
1:C:365:PRO:O	1:C:366:PHE:C	2.57	0.47
1:C:559:ARG:HH11	1:C:559:ARG:HG3	1.80	0.47
1:C:671:ARG:NH1	1:C:671:ARG:HG3	2.29	0.47
1:C:708:ALA:O	1:C:711:ILE:HG12	2.15	0.47
1:D:191:GLU:O	1:D:192:PHE:C	2.58	0.47
1:D:556:MET:O	1:D:560:LEU:HG	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:628:PHE:CD1	1:D:645:TRP:CD1	3.03	0.47
1:E:130:SER:HB2	1:E:170:TYR:HE2	1.77	0.47
1:E:170:TYR:HA	1:E:173:ILE:HG22	1.97	0.47
1:E:326:ILE:CG2	1:E:328:PHE:CE1	2.97	0.47
1:E:351:HIS:HB2	1:E:386:GLU:HG3	1.96	0.47
1:E:353:LYS:HG2	1:E:368:GLN:NE2	2.30	0.47
1:E:479:LYS:HG2	1:E:488:LEU:CD2	2.36	0.47
1:E:497:LEU:CD1	1:E:556:MET:HG2	2.39	0.47
1:F:153:ILE:C	1:F:154:ILE:HG12	2.38	0.47
1:F:171:TYR:O	1:F:175:LYS:NZ	2.47	0.47
1:F:559:ARG:HH11	1:F:559:ARG:HG3	1.80	0.47
2:O:3:GLN:N	2:O:77:LYS:HD3	2.30	0.47
2:O:70:THR:C	2:O:72:MET:N	2.72	0.47
2:P:58:ASP:C	2:P:60:ASN:N	2.65	0.47
2:P:101:SER:OG	2:P:104:GLU:HG2	2.15	0.47
2:Q:66:PRO:C	2:Q:68:PHE:N	2.72	0.47
2:S:38:SER:C	2:S:40:GLY:H	2.22	0.47
2:S:138:TYR:O	2:S:141:PHE:HB3	2.14	0.47
2:T:138:TYR:O	2:T:141:PHE:HB3	2.15	0.47
1:A:293:ILE:HD11	1:A:617:LYS:HD3	1.97	0.47
1:A:408:LEU:HD12	1:A:408:LEU:N	2.03	0.47
1:A:556:MET:O	1:A:560:LEU:HG	2.14	0.47
1:A:776:LEU:HD23	1:A:776:LEU:C	2.38	0.47
1:A:797:ILE:O	1:A:797:ILE:HG13	2.14	0.47
1:B:100:LEU:O	1:B:150:PRO:HD2	2.15	0.47
1:B:337:ASN:C	1:B:339:ILE:N	2.71	0.47
1:B:457:THR:CG2	1:B:469:PHE:H	2.27	0.47
1:B:629:ASN:HB3	1:B:632:TYR:CZ	2.50	0.47
1:C:89:ILE:HG22	1:C:93:VAL:CG1	2.33	0.47
1:C:299:GLU:HA	1:C:302:LEU:HB3	1.97	0.47
1:C:363:TYR:O	1:C:365:PRO:CD	2.63	0.47
1:C:797:ILE:O	1:C:797:ILE:HG13	2.15	0.47
1:D:131:ARG:HG3	1:D:243:LEU:HD22	1.96	0.47
1:D:175:LYS:O	1:D:176:GLY:C	2.58	0.47
1:D:184:LYS:HE2	1:D:193:LEU:CD1	2.30	0.47
1:D:299:GLU:HA	1:D:302:LEU:HB3	1.97	0.47
1:D:443:GLU:CD	1:D:458:LYS:HG2	2.40	0.47
1:E:135:VAL:HG22	1:E:135:VAL:O	2.14	0.47
1:E:220:LEU:HD11	1:E:223:LYS:HB2	1.97	0.47
1:E:629:ASN:HB3	1:E:632:TYR:CZ	2.50	0.47
1:E:653:LYS:O	1:E:655:ASN:N	2.47	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:680:LYS:HG2	1:E:681:ASP:N	2.29	0.47
1:E:687:GLU:O	1:E:690:LYS:N	2.48	0.47
1:E:746:LYS:HD2	1:E:747:ASN:ND2	2.29	0.47
1:F:142:VAL:HG22	1:F:154:ILE:CG2	2.41	0.47
1:F:318:ILE:H	1:F:318:ILE:CD1	2.23	0.47
1:F:326:ILE:CG2	1:F:328:PHE:CE1	2.98	0.47
1:F:443:GLU:CD	1:F:458:LYS:HG2	2.40	0.47
2:O:81:SER:O	2:O:82:GLU:C	2.58	0.47
2:Q:48:LEU:C	2:Q:51:MET:HE2	2.40	0.47
2:Q:132:GLY:C	2:Q:134:GLY:H	2.23	0.47
1:A:123:GLU:O	1:A:146:LYS:NZ	2.48	0.47
1:A:135:VAL:O	1:A:135:VAL:HG22	2.14	0.47
1:A:201:ASP:CG	1:A:225:ILE:HD12	2.40	0.47
1:A:353:LYS:HG2	1:A:368:GLN:NE2	2.30	0.47
1:A:457:THR:CG2	1:A:469:PHE:H	2.27	0.47
1:A:497:LEU:CD1	1:A:556:MET:HG2	2.40	0.47
1:A:680:LYS:HG2	1:A:681:ASP:N	2.30	0.47
1:B:282:SER:HA	1:B:285:LYS:CD	2.39	0.47
1:B:327:LEU:HG	1:B:595:ILE:HG12	1.97	0.47
1:B:648:PRO:O	1:B:649:ILE:C	2.57	0.47
1:C:191:GLU:O	1:C:192:PHE:C	2.58	0.47
1:C:217:LYS:HB3	1:C:236:GLU:CD	2.40	0.47
1:C:455:TYR:C	1:C:455:TYR:CD2	2.93	0.47
1:C:691:LYS:O	1:C:692:GLU:C	2.57	0.47
1:C:715:GLU:HA	1:C:718:ARG:HH12	1.71	0.47
1:D:135:VAL:O	1:D:135:VAL:HG22	2.14	0.47
1:D:190:PRO:O	1:D:191:GLU:C	2.58	0.47
1:D:288:VAL:O	1:D:292:ARG:HG2	2.14	0.47
1:D:326:ILE:CG2	1:D:328:PHE:CE1	2.98	0.47
1:D:365:PRO:O	1:D:366:PHE:C	2.57	0.47
1:D:457:THR:CG2	1:D:469:PHE:H	2.27	0.47
1:E:95:GLU:O	1:E:99:GLU:HB2	2.15	0.47
1:E:112:VAL:C	1:E:114:HIS:N	2.73	0.47
1:E:234:LEU:HD23	1:E:235:THR:N	2.27	0.47
1:E:400:LYS:HA	1:E:476:VAL:O	2.15	0.47
1:E:405:LEU:HD13	1:E:453:VAL:CG2	2.45	0.47
1:E:431:LYS:O	1:E:432:TYR:HD2	1.97	0.47
1:F:173:ILE:C	1:F:175:LYS:N	2.69	0.47
1:F:178:SER:OG	1:F:179:LEU:N	2.47	0.47
1:F:191:GLU:O	1:F:192:PHE:C	2.58	0.47
1:F:217:LYS:HZ1	1:F:236:GLU:HB2	1.79	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:434:LEU:HD22	1:F:445:ARG:HB3	1.96	0.47
1:F:684:ASP:C	1:F:686:ASP:N	2.70	0.47
2:P:81:SER:O	2:P:82:GLU:C	2.57	0.47
2:R:16:PHE:CD1	2:R:27:ILE:HD13	2.50	0.47
2:R:101:SER:H	2:R:104:GLU:CG	2.28	0.47
2:R:132:GLY:C	2:R:134:GLY:H	2.23	0.47
2:S:19:PHE:CD1	2:S:19:PHE:N	2.83	0.47
2:T:11:GLU:C	2:T:13:LYS:N	2.69	0.47
1:A:95:GLU:O	1:A:99:GLU:HB2	2.14	0.46
1:A:115:LYS:HB3	1:A:118:GLN:HG2	1.97	0.46
1:A:405:LEU:CD1	1:A:405:LEU:N	2.78	0.46
1:A:697:ILE:HG23	1:A:732:ILE:HD11	1.95	0.46
1:B:89:ILE:HG22	1:B:93:VAL:CG1	2.33	0.46
1:B:191:GLU:O	1:B:192:PHE:C	2.58	0.46
1:B:602:PHE:N	1:B:602:PHE:CD2	2.83	0.46
1:B:671:ARG:NH1	1:B:671:ARG:HG3	2.29	0.46
1:C:159:TYR:HD1	1:C:159:TYR:N	2.10	0.46
1:C:190:PRO:HB2	1:C:195:LEU:CD1	2.45	0.46
1:C:628:PHE:CD1	1:C:645:TRP:CD1	3.04	0.46
1:D:322:LEU:O	1:D:324:THR:HG22	2.15	0.46
1:D:462:ILE:HD11	1:D:466:GLY:HA2	1.96	0.46
1:D:628:PHE:CE2	2:R:90:ARG:CZ	2.98	0.46
1:D:694:VAL:CG2	1:D:695:LYS:N	2.78	0.46
1:E:412:GLU:C	1:E:414:LYS:H	2.23	0.46
1:E:445:ARG:HG2	1:E:471:TRP:CZ3	2.50	0.46
1:E:776:LEU:O	1:E:780:LEU:CD2	2.63	0.46
1:F:167:LYS:HG3	1:F:168:GLU:H	1.80	0.46
1:F:249:PHE:O	1:F:250:ALA:C	2.58	0.46
1:F:403:LEU:HG	1:F:405:LEU:HD11	1.97	0.46
1:F:405:LEU:HD13	1:F:453:VAL:CG2	2.44	0.46
1:F:540:ARG:NH1	1:F:627:TYR:CE1	2.83	0.46
2:O:13:LYS:O	2:O:15:ALA:N	2.47	0.46
2:P:132:GLY:C	2:P:134:GLY:H	2.23	0.46
2:Q:19:PHE:CD1	2:Q:19:PHE:N	2.84	0.46
2:R:51:MET:CB	2:R:71:MET:HE2	2.45	0.46
2:S:13:LYS:HE2	2:S:65:PHE:HB3	1.96	0.46
2:S:101:SER:H	2:S:104:GLU:CG	2.28	0.46
2:S:121:VAL:C	2:S:123:GLN:H	2.22	0.46
1:A:456:LYS:HG2	1:A:471:TRP:CD2	2.50	0.46
1:B:142:VAL:HA	1:B:154:ILE:CD1	2.45	0.46
1:B:190:PRO:O	1:B:195:LEU:HD13	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:667:LEU:O	1:B:669:SER:N	2.49	0.46
1:C:189:ASP:HB3	1:C:190:PRO:CD	2.45	0.46
1:C:391:ILE:O	1:C:392:THR:C	2.58	0.46
1:C:579:THR:C	1:C:581:GLN:N	2.69	0.46
1:C:629:ASN:HB3	1:C:632:TYR:CZ	2.50	0.46
1:D:186:LYS:CD	1:D:234:LEU:HD12	2.45	0.46
1:D:210:PHE:C	1:D:210:PHE:HD1	2.23	0.46
1:D:329:ARG:HB3	1:D:330:PRO:CD	2.45	0.46
1:D:797:ILE:HG13	1:D:797:ILE:O	2.14	0.46
1:E:249:PHE:O	1:E:250:ALA:C	2.57	0.46
1:E:322:LEU:O	1:E:324:THR:HG22	2.15	0.46
1:E:323:ASN:ND2	1:E:598:PRO:HB2	2.30	0.46
1:E:697:ILE:C	1:E:699:GLY:N	2.74	0.46
1:F:90:PRO:C	1:F:92:ASP:H	2.23	0.46
2:O:17:SER:O	2:O:19:PHE:N	2.48	0.46
2:O:58:ASP:O	2:O:60:ASN:N	2.45	0.46
2:O:106:ARG:NH2	2:O:118:ASP:OD2	2.48	0.46
2:P:13:LYS:HE2	2:P:65:PHE:HB3	1.96	0.46
2:Q:138:TYR:CE1	2:Q:142:VAL:CG2	2.98	0.46
2:R:30:LYS:HD3	2:R:30:LYS:N	2.12	0.46
2:S:137:ASN:O	2:S:138:TYR:C	2.58	0.46
2:T:36:MET:O	2:T:37:ARG:C	2.59	0.46
2:T:132:GLY:C	2:T:134:GLY:H	2.23	0.46
1:A:197:LYS:CA	1:A:197:LYS:HZ3	2.28	0.46
1:A:210:PHE:C	1:A:210:PHE:HD1	2.23	0.46
1:A:217:LYS:NZ	1:A:236:GLU:HB2	2.30	0.46
1:A:337:ASN:C	1:A:339:ILE:N	2.72	0.46
1:A:376:GLN:CB	1:A:379:ALA:HB3	2.46	0.46
1:B:628:PHE:CD1	1:B:645:TRP:CD1	3.03	0.46
1:B:708:ALA:O	1:B:711:ILE:HG12	2.15	0.46
1:C:322:LEU:O	1:C:324:THR:HG22	2.16	0.46
1:C:353:LYS:HG2	1:C:368:GLN:HE22	1.79	0.46
1:C:445:ARG:HG2	1:C:471:TRP:CZ3	2.50	0.46
1:C:694:VAL:CG2	1:C:695:LYS:N	2.78	0.46
1:D:89:ILE:HG22	1:D:90:PRO:HD2	1.97	0.46
1:D:90:PRO:C	1:D:92:ASP:N	2.72	0.46
1:D:95:GLU:O	1:D:99:GLU:HB2	2.15	0.46
1:D:186:LYS:HE3	1:D:234:LEU:CD1	2.45	0.46
1:D:504:ILE:N	1:D:504:ILE:HD12	2.30	0.46
1:D:730:ASN:C	1:D:732:ILE:H	2.22	0.46
1:D:776:LEU:O	1:D:780:LEU:CD2	2.64	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:142:VAL:HA	1:E:154:ILE:CD1	2.46	0.46
1:F:131:ARG:HG3	1:F:243:LEU:HD22	1.98	0.46
1:F:210:PHE:C	1:F:212:GLN:N	2.73	0.46
1:F:279:ILE:O	1:F:283:LEU:HB2	2.16	0.46
1:F:297:LYS:HB3	1:F:297:LYS:HZ2	1.79	0.46
1:F:457:THR:CG2	1:F:469:PHE:H	2.28	0.46
1:F:716:LYS:O	1:F:717:LYS:C	2.57	0.46
2:O:30:LYS:HD3	2:O:30:LYS:N	2.12	0.46
2:P:101:SER:H	2:P:104:GLU:CG	2.28	0.46
2:Q:6:GLU:O	2:Q:9:ILE:N	2.42	0.46
2:Q:17:SER:O	2:Q:19:PHE:N	2.48	0.46
2:Q:121:VAL:C	2:Q:123:GLN:N	2.67	0.46
2:S:49:GLN:O	2:S:53:ASN:N	2.41	0.46
2:T:102:ALA:CB	2:T:125:ILE:HG13	2.42	0.46
2:T:121:VAL:C	2:T:123:GLN:H	2.21	0.46
1:A:73:ASN:C	1:A:74:GLU:O	2.57	0.46
1:A:131:ARG:HG3	1:A:243:LEU:CD2	2.46	0.46
1:A:218:LEU:HG	1:A:218:LEU:O	2.15	0.46
1:A:444:PHE:CD2	1:A:455:TYR:HB3	2.51	0.46
1:A:711:ILE:O	1:A:712:PHE:HD2	1.99	0.46
1:A:764:LEU:O	1:A:766:HIS:N	2.47	0.46
1:A:776:LEU:O	1:A:780:LEU:CD2	2.64	0.46
1:B:353:LYS:HG2	1:B:368:GLN:NE2	2.30	0.46
1:B:427:ASP:O	1:B:428:ASN:HB2	2.15	0.46
1:B:540:ARG:NH1	1:B:627:TYR:CE1	2.83	0.46
1:B:625:LEU:HD12	1:B:625:LEU:C	2.39	0.46
1:C:167:LYS:HG3	1:C:168:GLU:H	1.80	0.46
1:C:182:ILE:O	1:C:187:SER:HB2	2.14	0.46
1:C:403:LEU:HG	1:C:405:LEU:HD11	1.97	0.46
1:C:456:LYS:HG2	1:C:471:TRP:CD2	2.50	0.46
1:C:635:ILE:H	1:C:635:ILE:HD12	1.80	0.46
1:D:142:VAL:HA	1:D:154:ILE:CD1	2.45	0.46
1:D:201:ASP:CG	1:D:225:ILE:HD12	2.39	0.46
1:D:400:LYS:HA	1:D:476:VAL:O	2.15	0.46
1:D:403:LEU:HG	1:D:405:LEU:HD11	1.97	0.46
1:E:217:LYS:NZ	1:E:236:GLU:HB2	2.30	0.46
1:E:444:PHE:CD2	1:E:455:TYR:HB3	2.50	0.46
1:E:462:ILE:HD11	1:E:466:GLY:HA2	1.96	0.46
1:E:559:ARG:HH11	1:E:559:ARG:HG3	1.80	0.46
1:E:691:LYS:O	1:E:692:GLU:C	2.58	0.46
1:F:186:LYS:HE3	1:F:234:LEU:CD1	2.45	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:217:LYS:HB3	1:F:236:GLU:CD	2.41	0.46
1:F:776:LEU:O	1:F:780:LEU:CD2	2.63	0.46
2:O:64:ASP:OD1	2:O:66:PRO:HD2	2.16	0.46
2:O:94:LYS:HE3	2:O:107:HIS:HD2	1.81	0.46
2:P:93:ASP:OD1	2:P:97:ASN:ND2	2.48	0.46
2:Q:63:ILE:HD11	2:Q:67:GLU:C	2.41	0.46
2:R:64:ASP:OD1	2:R:66:PRO:HD2	2.16	0.46
2:R:117:THR:C	2:R:119:GLU:N	2.73	0.46
2:S:81:SER:O	2:S:82:GLU:C	2.57	0.46
1:A:115:LYS:HE3	1:A:116:GLU:HG2	1.95	0.46
1:A:191:GLU:O	1:A:192:PHE:C	2.58	0.46
1:A:365:PRO:O	1:A:366:PHE:C	2.58	0.46
1:A:372:LYS:O	1:A:374:HIS:N	2.47	0.46
1:B:175:LYS:O	1:B:176:GLY:C	2.59	0.46
1:B:178:SER:OG	1:B:179:LEU:N	2.49	0.46
1:B:186:LYS:HE3	1:B:234:LEU:CD1	2.46	0.46
1:B:210:PHE:C	1:B:210:PHE:HD1	2.23	0.46
1:B:302:LEU:HB2	1:B:602:PHE:CD1	2.37	0.46
1:B:659:THR:O	1:B:660:SER:C	2.55	0.46
1:C:178:SER:OG	1:C:179:LEU:N	2.49	0.46
1:C:444:PHE:CD2	1:C:455:TYR:HB3	2.50	0.46
1:D:170:TYR:HA	1:D:173:ILE:CG2	2.46	0.46
1:D:431:LYS:O	1:D:432:TYR:HD2	1.99	0.46
1:E:73:ASN:O	1:E:74:GLU:C	2.58	0.46
1:E:90:PRO:C	1:E:92:ASP:H	2.23	0.46
1:E:443:GLU:CD	1:E:458:LYS:HG2	2.40	0.46
1:E:639:ASN:ND2	1:E:639:ASN:N	2.61	0.46
1:F:115:LYS:HB3	1:F:118:GLN:HG2	1.97	0.46
1:F:142:VAL:HA	1:F:154:ILE:CD1	2.46	0.46
1:F:190:PRO:O	1:F:191:GLU:C	2.58	0.46
1:F:387:ASN:N	1:F:387:ASN:ND2	2.50	0.46
1:F:400:LYS:HA	1:F:476:VAL:O	2.16	0.46
2:O:48:LEU:C	2:O:51:MET:HE2	2.41	0.46
2:O:101:SER:H	2:O:104:GLU:CG	2.28	0.46
2:O:137:ASN:O	2:O:138:TYR:C	2.59	0.46
2:Q:27:ILE:HD12	2:Q:32:LEU:HA	1.97	0.46
2:R:38:SER:C	2:R:40:GLY:H	2.22	0.46
2:S:13:LYS:O	2:S:15:ALA:N	2.49	0.46
1:A:403:LEU:HG	1:A:405:LEU:HD11	1.96	0.46
1:A:576:ASN:N	1:A:576:ASN:ND2	2.61	0.46
1:A:671:ARG:C	1:A:673:SER:H	2.23	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:115:LYS:HB3	1:B:118:GLN:HG2	1.97	0.46
1:B:288:VAL:O	1:B:292:ARG:HG2	2.15	0.46
1:B:405:LEU:CD1	1:B:405:LEU:N	2.79	0.46
1:C:112:VAL:C	1:C:114:HIS:N	2.73	0.46
1:C:431:LYS:O	1:C:432:TYR:HD2	1.98	0.46
1:D:73:ASN:O	1:D:74:GLU:C	2.58	0.46
1:D:112:VAL:C	1:D:114:HIS:N	2.73	0.46
1:D:311:HIS:HE1	1:D:339:ILE:HG22	1.81	0.46
1:D:376:GLN:CB	1:D:379:ALA:HB3	2.45	0.46
1:D:453:VAL:HG12	1:D:454:GLN:N	2.31	0.46
1:D:661:ALA:HB2	2:R:39:LEU:C	2.41	0.46
1:D:786:GLU:O	1:D:789:ASN:HB3	2.16	0.46
1:E:178:SER:OG	1:E:179:LEU:N	2.49	0.46
1:E:293:ILE:HD11	1:E:617:LYS:HD3	1.97	0.46
1:E:308:VAL:CG2	1:E:336:THR:O	2.64	0.46
1:E:540:ARG:NH1	1:E:627:TYR:CE1	2.83	0.46
1:E:748:TYR:O	1:E:751:TYR:N	2.46	0.46
1:F:376:GLN:CB	1:F:379:ALA:HB3	2.45	0.46
1:F:694:VAL:CG2	1:F:695:LYS:N	2.77	0.46
2:O:19:PHE:CD1	2:O:19:PHE:N	2.83	0.46
2:O:63:ILE:HD11	2:O:68:PHE:HA	1.98	0.46
2:O:70:THR:O	2:O:72:MET:N	2.48	0.46
2:O:126:ARG:HG3	2:O:126:ARG:NH2	2.29	0.46
2:P:117:THR:O	2:P:119:GLU:N	2.48	0.46
2:Q:106:ARG:NH2	2:Q:118:ASP:OD2	2.48	0.46
2:R:88:ALA:O	2:R:91:VAL:HB	2.15	0.46
2:R:126:ARG:HG3	2:R:126:ARG:NH2	2.27	0.46
1:A:97:TYR:HD2	1:A:102:GLY:HA3	1.78	0.46
1:A:105:TYR:O	1:A:154:ILE:N	2.35	0.46
1:A:602:PHE:CD2	1:A:602:PHE:N	2.84	0.46
1:B:90:PRO:C	1:B:92:ASP:H	2.24	0.46
1:B:177:ILE:C	1:B:180:ASP:OD1	2.58	0.46
1:B:456:LYS:HG2	1:B:471:TRP:CD2	2.51	0.46
1:B:493:ASP:OD2	1:B:577:HIS:HE1	1.98	0.46
1:B:694:VAL:CG2	1:B:695:LYS:N	2.79	0.46
1:C:112:VAL:O	1:C:114:HIS:N	2.39	0.46
1:C:279:ILE:O	1:C:283:LEU:HB2	2.16	0.46
1:C:497:LEU:CD1	1:C:556:MET:HG2	2.41	0.46
1:C:499:PRO:CG	1:C:504:ILE:HD11	2.45	0.46
1:C:648:PRO:O	1:C:651:LYS:N	2.47	0.46
1:C:671:ARG:C	1:C:673:SER:H	2.23	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:178:SER:OG	1:D:179:LEU:N	2.49	0.46
1:E:285:LYS:C	1:E:287:GLY:N	2.74	0.46
1:E:297:LYS:HZ1	1:E:601:GLU:HB3	1.80	0.46
1:E:337:ASN:C	1:E:339:ILE:N	2.72	0.46
1:E:456:LYS:HG2	1:E:471:TRP:CD2	2.51	0.46
1:E:658:PRO:HG2	1:E:701:LEU:HD22	1.97	0.46
1:F:137:PHE:O	1:F:139:SER:N	2.48	0.46
1:F:192:PHE:HA	1:F:195:LEU:CB	2.45	0.46
1:F:687:GLU:O	1:F:690:LYS:N	2.48	0.46
2:O:48:LEU:HD12	2:O:51:MET:CE	2.46	0.46
2:O:121:VAL:C	2:O:123:GLN:H	2.23	0.46
2:O:146:THR:O	2:O:147:ALA:C	2.59	0.46
2:Q:63:ILE:HD11	2:Q:68:PHE:HA	1.98	0.46
2:R:117:THR:O	2:R:119:GLU:N	2.48	0.46
1:A:89:ILE:HG22	1:A:90:PRO:HD2	1.97	0.46
1:A:481:VAL:HG21	1:A:486:LYS:HB2	1.98	0.46
1:A:661:ALA:HB2	2:O:38:SER:O	2.16	0.46
1:B:137:PHE:O	1:B:139:SER:N	2.48	0.46
1:B:338:LEU:HD21	1:B:409:ARG:NH1	2.30	0.46
1:B:412:GLU:C	1:B:414:LYS:H	2.22	0.46
1:B:630:ARG:CD	2:P:83:GLU:HG2	2.45	0.46
1:B:786:GLU:O	1:B:789:ASN:HB3	2.16	0.46
1:B:792:VAL:O	1:B:793:PHE:C	2.59	0.46
1:C:137:PHE:O	1:C:139:SER:N	2.49	0.46
1:C:190:PRO:O	1:C:195:LEU:HD13	2.15	0.46
1:C:210:PHE:C	1:C:212:GLN:N	2.73	0.46
1:C:481:VAL:O	1:C:482:GLU:HB2	2.14	0.46
1:C:602:PHE:CD2	1:C:602:PHE:N	2.84	0.46
1:C:724:ARG:HG3	1:C:724:ARG:NH1	2.30	0.46
1:C:748:TYR:O	1:C:751:TYR:N	2.48	0.46
1:C:792:VAL:O	1:C:793:PHE:C	2.59	0.46
1:D:115:LYS:HB3	1:D:118:GLN:HG2	1.97	0.46
1:D:293:ILE:HD11	1:D:617:LYS:HD3	1.97	0.46
1:D:297:LYS:HZ1	1:D:601:GLU:HB3	1.80	0.46
1:D:363:TYR:O	1:D:365:PRO:CD	2.64	0.46
1:D:625:LEU:HD12	1:D:625:LEU:C	2.40	0.46
1:D:691:LYS:NZ	2:R:20:ASP:O	2.47	0.46
1:D:746:LYS:HD2	1:D:747:ASN:HD22	1.81	0.46
1:E:137:PHE:O	1:E:139:SER:N	2.49	0.46
1:E:171:TYR:O	1:E:175:LYS:NZ	2.48	0.46
1:E:190:PRO:O	1:E:191:GLU:C	2.58	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:218:LEU:O	1:E:218:LEU:HG	2.15	0.46
1:E:311:HIS:HE1	1:E:339:ILE:HG22	1.79	0.46
1:E:376:GLN:CB	1:E:379:ALA:HB3	2.45	0.46
1:E:391:ILE:O	1:E:392:THR:C	2.58	0.46
1:E:403:LEU:HG	1:E:405:LEU:HD11	1.97	0.46
1:E:405:LEU:CD1	1:E:405:LEU:N	2.78	0.46
1:F:177:ILE:C	1:F:180:ASP:OD1	2.59	0.46
1:F:285:LYS:C	1:F:287:GLY:N	2.74	0.46
1:F:431:LYS:O	1:F:432:TYR:HD2	1.98	0.46
1:F:444:PHE:CD2	1:F:455:TYR:HB3	2.50	0.46
2:O:51:MET:CB	2:O:71:MET:HE2	2.46	0.46
2:P:58:ASP:O	2:P:60:ASN:N	2.46	0.46
2:P:138:TYR:O	2:P:141:PHE:HB3	2.15	0.46
2:R:63:ILE:HD11	2:R:68:PHE:HA	1.98	0.46
2:S:17:SER:O	2:S:19:PHE:N	2.49	0.46
2:S:63:ILE:HD11	2:S:68:PHE:HA	1.98	0.46
2:T:51:MET:CB	2:T:71:MET:HE2	2.45	0.46
1:A:288:VAL:O	1:A:292:ARG:HG2	2.15	0.46
1:A:363:TYR:O	1:A:365:PRO:CD	2.64	0.46
1:A:391:ILE:O	1:A:392:THR:C	2.59	0.46
1:A:654:ILE:HG22	1:A:655:ASN:ND2	2.30	0.46
1:A:694:VAL:CG2	1:A:695:LYS:N	2.79	0.46
1:B:171:TYR:O	1:B:175:LYS:NZ	2.49	0.46
1:B:249:PHE:O	1:B:250:ALA:C	2.57	0.46
1:B:322:LEU:O	1:B:324:THR:HG22	2.15	0.46
1:B:376:GLN:CB	1:B:379:ALA:HB3	2.46	0.46
1:B:552:TRP:O	1:B:553:GLN:C	2.57	0.46
1:B:658:PRO:HG3	1:B:752:LEU:HD22	1.98	0.46
1:B:746:LYS:HD2	1:B:747:ASN:HD22	1.80	0.46
1:C:293:ILE:HD11	1:C:617:LYS:HD3	1.97	0.46
1:C:610:MET:O	1:C:614:PHE:N	2.40	0.46
1:C:697:ILE:HG23	1:C:732:ILE:HD11	1.96	0.46
1:D:338:LEU:HD21	1:D:409:ARG:NH1	2.31	0.46
1:D:405:LEU:CD1	1:D:405:LEU:N	2.79	0.46
1:D:445:ARG:HG2	1:D:471:TRP:CZ3	2.50	0.46
1:D:670:ILE:O	1:D:671:ARG:C	2.58	0.46
1:E:97:TYR:C	1:E:99:GLU:N	2.73	0.46
1:E:175:LYS:O	1:E:176:GLY:C	2.59	0.46
1:E:199:LEU:C	1:E:201:ASP:N	2.71	0.46
1:E:708:ALA:O	1:E:711:ILE:HG12	2.15	0.46
1:F:130:SER:O	1:F:131:ARG:C	2.58	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:311:HIS:HE1	1:F:339:ILE:HG22	1.81	0.46
1:F:351:HIS:HB2	1:F:386:GLU:HG3	1.96	0.46
1:F:372:LYS:O	1:F:374:HIS:N	2.48	0.46
1:F:767:GLN:CG	1:F:768:LYS:HG2	2.36	0.46
2:P:51:MET:CB	2:P:71:MET:HE2	2.46	0.46
2:R:121:VAL:C	2:R:123:GLN:H	2.23	0.46
2:S:48:LEU:HD12	2:S:51:MET:CE	2.45	0.46
1:A:165:GLN:HE22	1:A:252:ASP:HB3	1.80	0.46
1:A:217:LYS:HB3	1:A:236:GLU:CD	2.41	0.46
1:A:492:TYR:HB2	1:A:575:VAL:HG22	1.98	0.46
1:A:629:ASN:HB3	1:A:632:TYR:CZ	2.51	0.46
1:B:391:ILE:O	1:B:392:THR:C	2.58	0.46
1:B:434:LEU:HD22	1:B:445:ARG:HB3	1.98	0.46
1:B:492:TYR:HB2	1:B:575:VAL:HG22	1.98	0.46
1:C:129:ASN:N	1:C:129:ASN:ND2	2.64	0.46
1:C:142:VAL:HA	1:C:154:ILE:CD1	2.46	0.46
1:C:311:HIS:HE1	1:C:339:ILE:HG22	1.81	0.46
1:C:318:ILE:H	1:C:318:ILE:CD1	2.23	0.46
1:D:167:LYS:HG3	1:D:168:GLU:H	1.80	0.46
1:D:308:VAL:CG2	1:D:336:THR:O	2.64	0.46
1:D:648:PRO:O	1:D:649:ILE:C	2.59	0.46
1:D:654:ILE:HG22	1:D:655:ASN:ND2	2.31	0.46
1:D:694:VAL:HG23	1:D:695:LYS:N	2.31	0.46
1:D:711:ILE:O	1:D:712:PHE:HD2	1.99	0.46
1:E:135:VAL:N	1:E:136:PRO:CD	2.78	0.46
1:E:190:PRO:O	1:E:195:LEU:HD13	2.16	0.46
1:E:197:LYS:HZ3	1:E:197:LYS:CA	2.29	0.46
1:E:318:ILE:H	1:E:318:ILE:CD1	2.23	0.46
1:F:175:LYS:O	1:F:176:GLY:C	2.59	0.46
2:O:140:GLU:HG2	2:O:140:GLU:H	1.55	0.46
2:P:3:GLN:N	2:P:77:LYS:HD3	2.30	0.46
2:P:38:SER:C	2:P:40:GLY:H	2.22	0.46
2:P:48:LEU:HD12	2:P:51:MET:CE	2.46	0.46
2:P:94:LYS:HE3	2:P:107:HIS:HD2	1.82	0.46
2:Q:84:GLU:N	2:Q:84:GLU:OE2	2.49	0.46
2:S:48:LEU:C	2:S:51:MET:HE2	2.41	0.46
2:S:48:LEU:O	2:S:52:ILE:HG22	2.16	0.46
2:S:106:ARG:NH2	2:S:118:ASP:OD2	2.49	0.46
2:T:66:PRO:C	2:T:68:PHE:N	2.72	0.46
1:A:192:PHE:HD1	1:A:192:PHE:H	1.64	0.45
1:A:794:GLN:HE21	1:A:794:GLN:HB3	1.54	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:182:ILE:O	1:B:187:SER:HB2	2.17	0.45
1:B:444:PHE:CD2	1:B:455:TYR:HB3	2.51	0.45
1:C:171:TYR:O	1:C:175:LYS:NZ	2.49	0.45
1:C:209:LEU:HD23	1:C:260:TYR:HA	1.98	0.45
1:C:400:LYS:HA	1:C:476:VAL:O	2.16	0.45
1:C:776:LEU:O	1:C:780:LEU:CD2	2.64	0.45
1:D:131:ARG:HG3	1:D:243:LEU:CD2	2.46	0.45
1:D:137:PHE:O	1:D:139:SER:N	2.48	0.45
1:D:292:ARG:NE	1:D:617:LYS:HE3	2.31	0.45
1:D:718:ARG:HH12	1:D:767:GLN:HE21	1.64	0.45
1:D:792:VAL:CG1	1:D:796:ILE:HD11	2.47	0.45
1:E:455:TYR:C	1:E:455:TYR:CD2	2.94	0.45
1:F:292:ARG:NE	1:F:617:LYS:HE3	2.31	0.45
1:F:338:LEU:HD21	1:F:409:ARG:NH1	2.31	0.45
1:F:602:PHE:CD2	1:F:602:PHE:N	2.84	0.45
2:O:48:LEU:O	2:O:52:ILE:HG22	2.16	0.45
2:P:48:LEU:O	2:P:52:ILE:HG22	2.16	0.45
2:P:63:ILE:HD11	2:P:68:PHE:HA	1.98	0.45
2:P:137:ASN:O	2:P:138:TYR:C	2.59	0.45
2:R:63:ILE:HD11	2:R:67:GLU:C	2.41	0.45
2:S:3:GLN:N	2:S:77:LYS:HD3	2.30	0.45
2:S:63:ILE:HD11	2:S:68:PHE:CA	2.46	0.45
2:S:64:ASP:OD1	2:S:66:PRO:HD2	2.16	0.45
2:T:76:MET:C	2:T:78:ASP:N	2.74	0.45
2:T:81:SER:O	2:T:82:GLU:C	2.58	0.45
1:A:142:VAL:HA	1:A:154:ILE:CD1	2.46	0.45
1:A:182:ILE:O	1:A:187:SER:HB2	2.16	0.45
1:A:190:PRO:O	1:A:191:GLU:C	2.59	0.45
1:A:400:LYS:HA	1:A:476:VAL:O	2.16	0.45
1:A:499:PRO:CG	1:A:504:ILE:HD11	2.46	0.45
1:A:500:SER:HG	1:A:502:THR:HG1	1.58	0.45
1:A:517:VAL:CB	1:A:525:LYS:HZ1	2.20	0.45
1:A:667:LEU:O	1:A:669:SER:N	2.49	0.45
1:B:170:TYR:HA	1:B:173:ILE:CG2	2.46	0.45
1:B:184:LYS:HE2	1:B:193:LEU:CD1	2.31	0.45
1:B:217:LYS:NZ	1:B:236:GLU:HB2	2.30	0.45
1:B:293:ILE:HD11	1:B:617:LYS:HD3	1.98	0.45
1:B:400:LYS:HA	1:B:476:VAL:O	2.17	0.45
1:B:615:ILE:HG13	1:B:615:ILE:H	1.60	0.45
1:B:711:ILE:O	1:B:712:PHE:HD2	1.98	0.45
1:C:88:LYS:NZ	1:C:172:GLU:CD	2.74	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:131:ARG:HG3	1:C:243:LEU:CD2	2.46	0.45
1:C:554:LYS:O	1:C:555:GLN:C	2.59	0.45
1:D:192:PHE:HA	1:D:195:LEU:CB	2.46	0.45
1:D:225:ILE:HA	1:D:229:PHE:CE2	2.47	0.45
1:D:234:LEU:HD23	1:D:235:THR:N	2.27	0.45
1:D:288:VAL:HG12	1:D:292:ARG:HH22	1.81	0.45
1:D:671:ARG:C	1:D:673:SER:H	2.23	0.45
1:D:697:ILE:C	1:D:699:GLY:N	2.73	0.45
1:E:167:LYS:HG3	1:E:168:GLU:H	1.80	0.45
1:E:190:PRO:HB2	1:E:195:LEU:CD1	2.46	0.45
1:E:236:GLU:C	1:E:238:GLN:H	2.24	0.45
1:E:648:PRO:O	1:E:649:ILE:C	2.59	0.45
1:E:656:THR:HG22	1:E:657:ILE:O	2.16	0.45
1:F:234:LEU:HD23	1:F:235:THR:N	2.27	0.45
1:F:255:THR:HG22	1:F:258:GLU:OE2	2.16	0.45
1:F:363:TYR:O	1:F:365:PRO:CD	2.65	0.45
1:F:395:GLU:C	1:F:395:GLU:CD	2.85	0.45
1:F:455:TYR:C	1:F:455:TYR:CD2	2.94	0.45
1:F:481:VAL:HG21	1:F:486:LYS:HB2	1.97	0.45
1:F:697:ILE:C	1:F:699:GLY:N	2.74	0.45
2:Q:101:SER:H	2:Q:104:GLU:CG	2.28	0.45
2:T:63:ILE:HD11	2:T:68:PHE:HA	1.98	0.45
1:A:311:HIS:HE1	1:A:339:ILE:HG22	1.80	0.45
1:A:639:ASN:HD22	1:A:639:ASN:N	2.14	0.45
1:A:657:ILE:CD1	1:A:701:LEU:HD23	2.47	0.45
1:B:130:SER:HB2	1:B:170:TYR:HE2	1.77	0.45
1:B:311:HIS:HE1	1:B:339:ILE:HG22	1.80	0.45
1:B:363:TYR:O	1:B:365:PRO:CD	2.64	0.45
1:B:540:ARG:HD2	1:B:582:ASP:OD1	2.17	0.45
1:C:658:PRO:HG2	1:C:701:LEU:HD22	1.98	0.45
1:C:747:ASN:N	1:C:747:ASN:HD22	2.14	0.45
1:C:764:LEU:O	1:C:766:HIS:N	2.49	0.45
1:E:129:ASN:N	1:E:129:ASN:ND2	2.64	0.45
1:E:192:PHE:HA	1:E:195:LEU:CB	2.46	0.45
1:E:372:LYS:O	1:E:374:HIS:N	2.47	0.45
1:E:671:ARG:C	1:E:673:SER:H	2.23	0.45
1:E:724:ARG:HG3	1:E:724:ARG:NH1	2.31	0.45
1:F:97:TYR:C	1:F:99:GLU:N	2.72	0.45
1:F:100:LEU:CD1	1:F:182:ILE:HG21	2.47	0.45
1:F:170:TYR:HA	1:F:173:ILE:CG2	2.46	0.45
1:F:182:ILE:O	1:F:187:SER:HB2	2.15	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:337:ASN:C	1:F:339:ILE:N	2.71	0.45
1:F:391:ILE:O	1:F:392:THR:C	2.58	0.45
1:F:445:ARG:HG2	1:F:471:TRP:CZ3	2.51	0.45
1:F:630:ARG:CD	2:T:83:GLU:HG2	2.46	0.45
1:F:671:ARG:C	1:F:673:SER:H	2.23	0.45
1:F:708:ALA:O	1:F:711:ILE:HG12	2.16	0.45
2:O:132:GLY:C	2:O:134:GLY:H	2.24	0.45
2:P:63:ILE:HD11	2:P:67:GLU:C	2.41	0.45
2:R:48:LEU:O	2:R:52:ILE:HG22	2.16	0.45
2:R:84:GLU:N	2:R:84:GLU:OE2	2.50	0.45
2:R:94:LYS:HB3	2:R:94:LYS:HZ2	1.81	0.45
2:S:102:ALA:CB	2:S:125:ILE:HG13	2.42	0.45
2:T:38:SER:C	2:T:40:GLY:H	2.23	0.45
2:T:48:LEU:C	2:T:51:MET:HE2	2.41	0.45
2:T:48:LEU:O	2:T:52:ILE:HG22	2.16	0.45
1:A:279:ILE:O	1:A:283:LEU:HB2	2.17	0.45
1:A:453:VAL:HG12	1:A:454:GLN:N	2.31	0.45
1:A:653:LYS:O	1:A:655:ASN:N	2.49	0.45
1:A:659:THR:O	1:A:660:SER:C	2.56	0.45
1:B:190:PRO:O	1:B:191:GLU:C	2.58	0.45
1:B:438:ASN:HD22	1:B:438:ASN:HA	1.60	0.45
1:B:453:VAL:HG12	1:B:454:GLN:N	2.31	0.45
1:B:455:TYR:C	1:B:455:TYR:CD2	2.93	0.45
1:C:90:PRO:C	1:C:92:ASP:H	2.24	0.45
1:C:288:VAL:HG12	1:C:292:ARG:HH22	1.81	0.45
1:C:376:GLN:CB	1:C:379:ALA:HB3	2.46	0.45
1:C:654:ILE:HG22	1:C:655:ASN:ND2	2.31	0.45
1:C:694:VAL:O	1:C:695:LYS:C	2.59	0.45
1:C:711:ILE:O	1:C:712:PHE:HD2	2.00	0.45
1:D:182:ILE:O	1:D:187:SER:HB2	2.16	0.45
1:D:192:PHE:HD1	1:D:192:PHE:H	1.65	0.45
1:D:456:LYS:HG2	1:D:471:TRP:CD2	2.51	0.45
1:D:554:LYS:O	1:D:555:GLN:C	2.59	0.45
1:D:576:ASN:N	1:D:576:ASN:ND2	2.63	0.45
1:D:602:PHE:N	1:D:602:PHE:CD2	2.84	0.45
1:D:792:VAL:O	1:D:793:PHE:C	2.60	0.45
1:E:123:GLU:O	1:E:146:LYS:NZ	2.48	0.45
1:E:177:ILE:C	1:E:180:ASP:OD1	2.59	0.45
1:E:338:LEU:HD21	1:E:409:ARG:NH1	2.31	0.45
1:F:263:ASP:O	1:F:266:GLU:N	2.50	0.45
1:F:456:LYS:HG2	1:F:471:TRP:CD2	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:747:ASN:O	1:F:748:TYR:C	2.60	0.45
1:F:764:LEU:O	1:F:766:HIS:N	2.50	0.45
2:P:17:SER:O	2:P:20:ASP:N	2.40	0.45
2:R:48:LEU:HD12	2:R:51:MET:CE	2.46	0.45
2:R:97:ASN:ND2	2:R:97:ASN:C	2.74	0.45
2:S:58:ASP:O	2:S:60:ASN:N	2.46	0.45
2:T:19:PHE:N	2:T:19:PHE:CD1	2.84	0.45
1:A:175:LYS:O	1:A:176:GLY:C	2.59	0.45
1:A:292:ARG:NE	1:A:617:LYS:HE3	2.32	0.45
1:A:792:VAL:CG1	1:A:796:ILE:HD11	2.47	0.45
1:B:372:LYS:O	1:B:374:HIS:N	2.47	0.45
1:B:409:ARG:CD	1:B:413:LEU:HD21	2.47	0.45
1:B:671:ARG:C	1:B:673:SER:H	2.24	0.45
1:B:724:ARG:HG3	1:B:724:ARG:NH1	2.32	0.45
1:C:220:LEU:HG	1:C:223:LYS:HB3	1.99	0.45
1:C:767:GLN:HG2	1:C:768:LYS:H	1.81	0.45
1:D:694:VAL:O	1:D:695:LYS:C	2.60	0.45
1:D:726:ILE:O	1:D:727:GLN:C	2.60	0.45
1:E:191:GLU:O	1:E:192:PHE:C	2.58	0.45
1:E:363:TYR:O	1:E:365:PRO:CD	2.65	0.45
1:E:409:ARG:CD	1:E:413:LEU:HD21	2.47	0.45
1:E:746:LYS:HD2	1:E:747:ASN:HD22	1.82	0.45
1:F:90:PRO:C	1:F:92:ASP:N	2.72	0.45
1:F:97:TYR:O	1:F:99:GLU:N	2.49	0.45
1:F:131:ARG:HG3	1:F:243:LEU:CD2	2.46	0.45
1:F:499:PRO:CG	1:F:504:ILE:HD11	2.46	0.45
1:F:611:THR:HG22	1:F:615:ILE:CD1	2.45	0.45
1:F:711:ILE:O	1:F:712:PHE:HD2	1.99	0.45
2:P:48:LEU:C	2:P:51:MET:HE2	2.42	0.45
2:P:63:ILE:HD11	2:P:68:PHE:CA	2.47	0.45
2:Q:63:ILE:HD11	2:Q:68:PHE:CA	2.46	0.45
2:Q:121:VAL:C	2:Q:123:GLN:H	2.22	0.45
2:R:63:ILE:HD11	2:R:68:PHE:CA	2.47	0.45
2:S:36:MET:O	2:S:37:ARG:C	2.60	0.45
2:S:117:THR:O	2:S:119:GLU:N	2.49	0.45
2:T:63:ILE:HD11	2:T:67:GLU:C	2.42	0.45
2:T:106:ARG:NH2	2:T:118:ASP:OD2	2.50	0.45
1:A:73:ASN:O	1:A:74:GLU:C	2.58	0.45
1:A:190:PRO:O	1:A:195:LEU:HD13	2.16	0.45
1:A:241:PHE:CD1	1:A:268:MET:HE1	2.52	0.45
1:B:609:GLU:O	1:B:613:ARG:N	2.34	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:792:VAL:CG1	1:B:796:ILE:HD11	2.47	0.45
1:C:192:PHE:HA	1:C:195:LEU:CB	2.46	0.45
1:C:247:TYR:O	1:C:253:HIS:O	2.35	0.45
1:C:746:LYS:HD2	1:C:747:ASN:HD22	1.81	0.45
1:D:190:PRO:O	1:D:195:LEU:HD13	2.16	0.45
1:D:302:LEU:HB2	1:D:602:PHE:CD1	2.37	0.45
1:D:724:ARG:HG3	1:D:724:ARG:NH1	2.30	0.45
1:E:131:ARG:HG3	1:E:243:LEU:HD22	1.98	0.45
1:E:234:LEU:O	1:E:238:GLN:HG3	2.17	0.45
1:E:792:VAL:O	1:E:793:PHE:C	2.59	0.45
1:F:190:PRO:O	1:F:195:LEU:HD13	2.17	0.45
1:F:554:LYS:O	1:F:555:GLN:C	2.59	0.45
1:F:576:ASN:N	1:F:576:ASN:ND2	2.64	0.45
2:Q:65:PHE:CB	2:Q:66:PRO:HD3	2.46	0.45
2:Q:102:ALA:CB	2:Q:125:ILE:HG13	2.42	0.45
2:T:58:ASP:O	2:T:60:ASN:N	2.45	0.45
1:A:175:LYS:O	1:A:177:ILE:N	2.49	0.45
1:A:178:SER:OG	1:A:179:LEU:N	2.49	0.45
1:A:288:VAL:HG12	1:A:292:ARG:HH22	1.82	0.45
1:A:297:LYS:NZ	1:A:297:LYS:CB	2.79	0.45
1:A:325:TYR:CD1	1:A:598:PRO:HD3	2.52	0.45
1:A:395:GLU:C	1:A:395:GLU:CD	2.84	0.45
1:A:401:ILE:HG21	1:A:485:LEU:HB3	1.99	0.45
1:A:455:TYR:C	1:A:455:TYR:CD2	2.93	0.45
1:A:746:LYS:HD2	1:A:747:ASN:HD22	1.81	0.45
1:B:167:LYS:HG3	1:B:168:GLU:H	1.80	0.45
1:B:220:LEU:HG	1:B:223:LYS:HB3	1.99	0.45
1:B:657:ILE:CD1	1:B:701:LEU:HD23	2.47	0.45
1:C:170:TYR:HA	1:C:173:ILE:CG2	2.46	0.45
1:C:355:SER:HA	1:C:372:LYS:H	1.82	0.45
1:C:792:VAL:CG1	1:C:796:ILE:HD11	2.47	0.45
1:D:217:LYS:HB3	1:D:236:GLU:CD	2.42	0.45
1:D:305:SER:C	1:D:307:LEU:N	2.75	0.45
1:D:323:ASN:ND2	1:D:598:PRO:HB2	2.32	0.45
1:E:365:PRO:O	1:E:366:PHE:C	2.57	0.45
1:E:453:VAL:HG12	1:E:454:GLN:N	2.31	0.45
1:F:115:LYS:HE3	1:F:116:GLU:HG2	1.95	0.45
1:F:175:LYS:HB2	1:F:175:LYS:HZ2	1.82	0.45
1:F:201:ASP:CG	1:F:225:ILE:HD12	2.42	0.45
1:F:209:LEU:HD23	1:F:260:TYR:HA	1.98	0.45
1:F:495:PHE:CD1	1:F:495:PHE:O	2.70	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:R:17:SER:O	2:R:19:PHE:N	2.50	0.45
2:R:97:ASN:C	2:R:99:TYR:N	2.75	0.45
2:T:36:MET:HE3	2:T:43:PRO:CG	2.31	0.45
2:T:89:PHE:HB2	2:T:141:PHE:CD2	2.52	0.45
1:A:100:LEU:CD1	1:A:182:ILE:HG21	2.46	0.45
1:A:170:TYR:HA	1:A:173:ILE:CG2	2.47	0.45
1:A:192:PHE:HA	1:A:195:LEU:CB	2.46	0.45
1:A:247:TYR:O	1:A:253:HIS:O	2.35	0.45
1:A:609:GLU:O	1:A:613:ARG:N	2.34	0.45
1:A:708:ALA:O	1:A:711:ILE:HG12	2.16	0.45
1:A:747:ASN:O	1:A:750:GLN:N	2.50	0.45
1:B:192:PHE:HD1	1:B:192:PHE:H	1.65	0.45
1:B:726:ILE:O	1:B:727:GLN:C	2.59	0.45
1:C:297:LYS:NZ	1:C:297:LYS:CB	2.80	0.45
1:C:323:ASN:ND2	1:C:598:PRO:HB2	2.32	0.45
1:C:338:LEU:HD21	1:C:409:ARG:NH1	2.32	0.45
1:C:747:ASN:O	1:C:750:GLN:N	2.50	0.45
1:E:97:TYR:O	1:E:99:GLU:N	2.50	0.45
1:E:97:TYR:HD2	1:E:102:GLY:HA3	1.78	0.45
1:E:217:LYS:HB3	1:E:236:GLU:CD	2.41	0.45
1:E:504:ILE:HD12	1:E:504:ILE:N	2.32	0.45
1:E:629:ASN:CB	1:E:632:TYR:CE1	2.96	0.45
1:F:365:PRO:O	1:F:366:PHE:C	2.57	0.45
1:F:670:ILE:O	1:F:671:ARG:C	2.59	0.45
2:P:64:ASP:OD1	2:P:66:PRO:HD2	2.17	0.45
2:Q:97:ASN:HD22	2:Q:97:ASN:N	2.15	0.45
2:R:37:ARG:HH11	2:R:37:ARG:CG	2.30	0.45
2:S:146:THR:O	2:S:147:ALA:C	2.59	0.45
2:T:144:MET:HE1	2:T:145:MET:CE	2.47	0.45
1:A:71:PHE:CG	1:A:73:ASN:HB2	2.52	0.45
1:A:197:LYS:CB	1:A:197:LYS:HZ3	2.30	0.45
1:A:630:ARG:CD	2:O:83:GLU:HG2	2.47	0.45
1:A:658:PRO:HG2	1:A:701:LEU:HD22	1.99	0.45
1:A:662:GLU:OE2	1:A:755:ARG:NH2	2.42	0.45
1:A:670:ILE:O	1:A:671:ARG:C	2.60	0.45
1:A:691:LYS:NZ	2:O:20:ASP:O	2.47	0.45
1:A:786:GLU:O	1:A:789:ASN:HB3	2.17	0.45
1:B:172:GLU:CB	1:B:246:SER:HA	2.47	0.45
1:B:236:GLU:C	1:B:238:GLN:H	2.24	0.45
1:B:610:MET:O	1:B:614:PHE:N	2.40	0.45
1:C:175:LYS:O	1:C:177:ILE:N	2.50	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:180:ASP:O	1:C:183:SER:N	2.36	0.45
1:C:190:PRO:O	1:C:191:GLU:C	2.59	0.45
1:C:395:GLU:C	1:C:395:GLU:CD	2.85	0.45
1:C:540:ARG:NH1	1:C:627:TYR:CE1	2.85	0.45
1:C:786:GLU:O	1:C:789:ASN:HB3	2.17	0.45
1:D:247:TYR:O	1:D:253:HIS:O	2.35	0.45
1:D:279:ILE:O	1:D:283:LEU:HB2	2.17	0.45
1:D:327:LEU:HG	1:D:595:ILE:HG12	1.97	0.45
1:D:424:LYS:HB3	1:D:424:LYS:NZ	2.29	0.45
1:E:192:PHE:HD1	1:E:192:PHE:H	1.65	0.45
1:E:292:ARG:NE	1:E:617:LYS:HE3	2.32	0.45
1:F:423:LYS:HG2	1:F:424:LYS:N	2.32	0.45
2:O:12:PHE:CE1	2:O:72:MET:CE	2.94	0.45
2:Q:64:ASP:OD1	2:Q:66:PRO:HD2	2.17	0.45
2:R:48:LEU:C	2:R:51:MET:HE2	2.41	0.45
2:S:63:ILE:HD11	2:S:67:GLU:C	2.42	0.45
2:S:76:MET:C	2:S:78:ASP:N	2.75	0.45
2:S:117:THR:C	2:S:119:GLU:N	2.73	0.45
2:T:48:LEU:HD12	2:T:51:MET:CE	2.46	0.45
2:T:117:THR:C	2:T:119:GLU:N	2.73	0.45
1:A:112:VAL:C	1:A:114:HIS:N	2.73	0.45
1:A:167:LYS:HG3	1:A:168:GLU:H	1.81	0.45
1:A:270:LYS:HA	1:A:273:LYS:CD	2.42	0.45
1:A:285:LYS:C	1:A:287:GLY:N	2.74	0.45
1:A:395:GLU:CD	1:A:395:GLU:H	2.25	0.45
1:A:724:ARG:HG3	1:A:724:ARG:NH1	2.30	0.45
1:B:71:PHE:CG	1:B:73:ASN:HB2	2.52	0.45
1:B:279:ILE:O	1:B:283:LEU:HB2	2.17	0.45
1:B:345:THR:CG2	1:B:491:ASP:HA	2.47	0.45
1:B:479:LYS:HG2	1:B:488:LEU:CD2	2.36	0.45
1:B:504:ILE:HD12	1:B:504:ILE:N	2.32	0.45
1:B:694:VAL:HG23	1:B:695:LYS:N	2.32	0.45
1:C:454:GLN:HG2	1:C:473:ASN:HA	1.99	0.45
1:C:697:ILE:C	1:C:699:GLY:N	2.74	0.45
1:D:173:ILE:HG13	1:D:242:SER:CB	2.35	0.45
1:D:438:ASN:HD22	1:D:438:ASN:HA	1.59	0.45
1:E:395:GLU:C	1:E:395:GLU:CD	2.85	0.45
1:E:481:VAL:HG21	1:E:486:LYS:HB2	1.98	0.45
1:E:670:ILE:O	1:E:671:ARG:C	2.60	0.45
1:F:454:GLN:HG2	1:F:473:ASN:HA	1.99	0.45
1:F:540:ARG:HD2	1:F:582:ASP:OD1	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:661:ALA:HB2	2:T:39:LEU:C	2.42	0.45
1:F:697:ILE:HG23	1:F:732:ILE:HD11	1.96	0.45
2:O:97:ASN:ND2	2:O:97:ASN:C	2.75	0.45
2:P:17:SER:O	2:P:19:PHE:N	2.50	0.45
2:R:76:MET:C	2:R:78:ASP:N	2.75	0.45
2:R:137:ASN:O	2:R:138:TYR:C	2.59	0.45
2:T:37:ARG:HH11	2:T:37:ARG:CG	2.30	0.45
2:T:39:LEU:HD23	2:T:39:LEU:HA	1.86	0.45
2:T:146:THR:O	2:T:147:ALA:C	2.59	0.45
1:A:236:GLU:C	1:A:238:GLN:H	2.24	0.44
1:A:360:VAL:O	1:A:361:ALA:C	2.60	0.44
1:A:409:ARG:CZ	1:A:413:LEU:CD2	2.94	0.44
1:A:479:LYS:HG2	1:A:488:LEU:CD2	2.36	0.44
1:A:532:LEU:HD23	1:A:532:LEU:HA	1.80	0.44
1:A:726:ILE:O	1:A:727:GLN:C	2.61	0.44
1:B:83:GLN:O	1:B:86:LEU:N	2.49	0.44
1:B:175:LYS:O	1:B:177:ILE:N	2.50	0.44
1:B:192:PHE:HA	1:B:195:LEU:CB	2.46	0.44
1:B:395:GLU:C	1:B:395:GLU:CD	2.85	0.44
1:C:131:ARG:HG3	1:C:243:LEU:HD22	1.98	0.44
1:C:177:ILE:C	1:C:180:ASP:OD1	2.59	0.44
1:C:581:GLN:HE22	1:C:629:ASN:H	1.56	0.44
1:C:615:ILE:HG13	1:C:615:ILE:H	1.59	0.44
1:C:776:LEU:HD11	1:C:793:PHE:HE1	1.82	0.44
1:D:177:ILE:C	1:D:180:ASP:OD1	2.60	0.44
1:D:395:GLU:C	1:D:395:GLU:CD	2.85	0.44
1:D:423:LYS:HG2	1:D:424:LYS:N	2.32	0.44
1:D:454:GLN:HG2	1:D:473:ASN:HA	1.99	0.44
1:D:694:VAL:C	1:D:696:LYS:N	2.74	0.44
1:D:708:ALA:O	1:D:711:ILE:HG12	2.16	0.44
1:E:175:LYS:O	1:E:177:ILE:N	2.50	0.44
1:E:360:VAL:O	1:E:361:ALA:C	2.59	0.44
1:E:736:LEU:HD21	1:E:750:GLN:HE22	1.82	0.44
1:E:786:GLU:O	1:E:789:ASN:HB3	2.17	0.44
1:F:255:THR:C	1:F:257:LEU:N	2.74	0.44
1:F:293:ILE:HD11	1:F:617:LYS:HD3	1.98	0.44
1:F:325:TYR:CD1	1:F:598:PRO:HD3	2.52	0.44
1:F:554:LYS:O	1:F:557:LEU:N	2.50	0.44
1:F:694:VAL:HG23	1:F:695:LYS:N	2.31	0.44
1:F:747:ASN:N	1:F:747:ASN:HD22	2.16	0.44
2:O:63:ILE:HD11	2:O:68:PHE:CA	2.47	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:Q:37:ARG:HH11	2:Q:37:ARG:CG	2.31	0.44
2:Q:48:LEU:O	2:Q:52:ILE:HG22	2.17	0.44
2:R:48:LEU:HA	2:R:51:MET:CE	2.41	0.44
2:R:144:MET:HE1	2:R:145:MET:HE1	1.99	0.44
2:R:146:THR:O	2:R:147:ALA:C	2.59	0.44
2:T:137:ASN:O	2:T:138:TYR:C	2.59	0.44
1:A:175:LYS:HB2	1:A:175:LYS:HZ2	1.80	0.44
1:A:409:ARG:CD	1:A:413:LEU:HD21	2.46	0.44
1:A:648:PRO:O	1:A:649:ILE:C	2.58	0.44
1:A:747:ASN:O	1:A:748:TYR:C	2.60	0.44
1:B:135:VAL:N	1:B:136:PRO:CD	2.80	0.44
1:B:288:VAL:HG12	1:B:292:ARG:HH22	1.83	0.44
1:B:323:ASN:ND2	1:B:598:PRO:HB2	2.31	0.44
1:B:499:PRO:CG	1:B:504:ILE:HD11	2.45	0.44
1:B:551:ASN:HD22	1:B:551:ASN:HA	1.65	0.44
1:C:83:GLN:O	1:C:86:LEU:N	2.50	0.44
1:C:175:LYS:O	1:C:176:GLY:C	2.60	0.44
1:D:297:LYS:NZ	1:D:297:LYS:CB	2.80	0.44
1:D:333:LYS:O	1:D:335:ALA:N	2.48	0.44
1:D:455:TYR:C	1:D:455:TYR:CD2	2.95	0.44
1:E:288:VAL:O	1:E:292:ARG:HG2	2.17	0.44
1:E:409:ARG:CZ	1:E:413:LEU:CD2	2.94	0.44
1:E:597:ASN:ND2	1:E:601:GLU:H	2.03	0.44
1:F:172:GLU:CB	1:F:246:SER:HA	2.47	0.44
1:F:173:ILE:HG13	1:F:242:SER:CB	2.35	0.44
1:F:405:LEU:CD1	1:F:405:LEU:N	2.80	0.44
1:F:648:PRO:O	1:F:649:ILE:C	2.60	0.44
1:F:718:ARG:HH12	1:F:767:GLN:HE21	1.65	0.44
1:F:719:LYS:O	1:F:720:ILE:C	2.60	0.44
1:F:746:LYS:HD2	1:F:747:ASN:HD22	1.81	0.44
2:P:36:MET:O	2:P:37:ARG:C	2.59	0.44
2:P:86:ARG:O	2:P:86:ARG:CG	2.65	0.44
2:S:89:PHE:HB2	2:S:141:PHE:CE2	2.52	0.44
1:A:137:PHE:O	1:A:139:SER:N	2.50	0.44
1:A:333:LYS:HA	1:A:336:THR:OG1	2.17	0.44
1:A:345:THR:CG2	1:A:491:ASP:HA	2.47	0.44
1:A:504:ILE:N	1:A:504:ILE:HD12	2.31	0.44
1:A:694:VAL:HG23	1:A:695:LYS:N	2.33	0.44
1:B:247:TYR:O	1:B:253:HIS:O	2.35	0.44
1:B:325:TYR:CD1	1:B:598:PRO:HD3	2.52	0.44
1:C:109:ILE:O	1:C:109:ILE:HG22	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:115:LYS:HB3	1:C:118:GLN:HG2	1.97	0.44
1:C:182:ILE:C	1:C:187:SER:HB2	2.42	0.44
1:C:292:ARG:NE	1:C:617:LYS:HE3	2.33	0.44
1:C:326:ILE:CG2	1:C:328:PHE:CE1	3.00	0.44
1:C:540:ARG:HD2	1:C:582:ASP:OD1	2.17	0.44
1:C:700:TYR:CE1	1:C:727:GLN:HB3	2.53	0.44
1:D:220:LEU:HG	1:D:223:LYS:HB3	1.99	0.44
1:D:234:LEU:O	1:D:238:GLN:HG3	2.17	0.44
1:D:540:ARG:HD2	1:D:582:ASP:OD1	2.17	0.44
1:D:629:ASN:HB3	1:D:632:TYR:CZ	2.50	0.44
1:D:630:ARG:CD	2:R:83:GLU:HG2	2.48	0.44
1:D:658:PRO:HG2	1:D:701:LEU:HD22	1.98	0.44
1:E:115:LYS:HE3	1:E:116:GLU:HG2	1.94	0.44
1:E:115:LYS:HE2	1:E:116:GLU:HG2	1.98	0.44
1:E:454:GLN:HG2	1:E:473:ASN:HA	2.00	0.44
1:E:602:PHE:N	1:E:602:PHE:CD2	2.84	0.44
1:F:190:PRO:HB2	1:F:195:LEU:CD1	2.46	0.44
1:F:688:PHE:C	1:F:688:PHE:HD2	2.24	0.44
1:F:726:ILE:O	1:F:727:GLN:C	2.59	0.44
1:F:767:GLN:HG2	1:F:768:LYS:H	1.82	0.44
2:Q:48:LEU:HD12	2:Q:51:MET:CE	2.46	0.44
2:Q:49:GLN:O	2:Q:53:ASN:N	2.41	0.44
2:Q:68:PHE:O	2:Q:70:THR:N	2.51	0.44
2:R:144:MET:HE1	2:R:145:MET:CE	2.46	0.44
2:S:126:ARG:HG3	2:S:126:ARG:NH2	2.30	0.44
2:T:97:ASN:ND2	2:T:97:ASN:C	2.76	0.44
1:A:78:LYS:O	1:A:81:GLN:HB3	2.17	0.44
1:A:97:TYR:CE1	1:A:178:SER:HB2	2.53	0.44
1:A:172:GLU:CB	1:A:246:SER:HA	2.47	0.44
1:A:209:LEU:HD23	1:A:260:TYR:HA	1.99	0.44
1:A:326:ILE:CG2	1:A:328:PHE:CE1	2.98	0.44
1:A:454:GLN:HG2	1:A:473:ASN:HA	2.00	0.44
1:A:747:ASN:HD22	1:A:747:ASN:N	2.16	0.44
1:A:792:VAL:O	1:A:793:PHE:C	2.60	0.44
1:B:199:LEU:C	1:B:201:ASP:N	2.71	0.44
1:B:201:ASP:CG	1:B:225:ILE:HD12	2.42	0.44
1:B:209:LEU:HD23	1:B:260:TYR:HA	1.98	0.44
1:B:308:VAL:CG2	1:B:336:THR:O	2.65	0.44
1:B:326:ILE:CG2	1:B:328:PHE:CE1	2.99	0.44
1:B:401:ILE:HG21	1:B:485:LEU:HB3	2.00	0.44
1:B:403:LEU:HG	1:B:405:LEU:HD11	1.98	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:653:LYS:C	1:B:655:ASN:N	2.75	0.44
1:C:152:LEU:HG	1:C:154:ILE:HD11	2.00	0.44
1:C:255:THR:C	1:C:257:LEU:N	2.75	0.44
1:C:630:ARG:CD	2:Q:83:GLU:HG2	2.47	0.44
1:C:726:ILE:O	1:C:727:GLN:C	2.60	0.44
1:D:307:LEU:HD12	1:D:331:VAL:CG2	2.47	0.44
1:D:587:PRO:C	1:D:588:GLU:HG2	2.43	0.44
1:D:653:LYS:C	1:D:655:ASN:N	2.76	0.44
1:D:666:ASN:HB2	1:D:748:TYR:CZ	2.53	0.44
1:D:767:GLN:HG2	1:D:768:LYS:H	1.82	0.44
1:E:96:ILE:O	1:E:100:LEU:HG	2.17	0.44
1:E:209:LEU:HD23	1:E:260:TYR:HA	1.99	0.44
1:E:297:LYS:NZ	1:E:297:LYS:CB	2.80	0.44
1:E:333:LYS:HA	1:E:336:THR:OG1	2.18	0.44
1:F:71:PHE:CG	1:F:73:ASN:HB2	2.52	0.44
1:F:83:GLN:O	1:F:86:LEU:N	2.51	0.44
1:F:397:GLU:O	1:F:480:ASN:N	2.51	0.44
1:F:792:VAL:O	1:F:793:PHE:C	2.59	0.44
2:O:104:GLU:O	2:O:108:VAL:HG23	2.18	0.44
2:O:144:MET:HE1	2:O:145:MET:CE	2.48	0.44
2:P:65:PHE:CB	2:P:66:PRO:HD3	2.47	0.44
2:P:146:THR:O	2:P:147:ALA:C	2.59	0.44
2:S:117:THR:O	2:S:118:ASP:C	2.60	0.44
2:T:91:VAL:HG12	2:T:92:PHE:N	2.33	0.44
2:T:97:ASN:C	2:T:99:TYR:N	2.76	0.44
1:A:551:ASN:HD22	1:A:551:ASN:HA	1.66	0.44
1:A:587:PRO:C	1:A:588:GLU:HG2	2.42	0.44
1:A:667:LEU:C	1:A:669:SER:N	2.75	0.44
1:B:131:ARG:HG3	1:B:243:LEU:HD22	1.98	0.44
1:B:285:LYS:C	1:B:287:GLY:N	2.75	0.44
1:B:297:LYS:NZ	1:B:297:LYS:CB	2.80	0.44
1:B:333:LYS:HA	1:B:336:THR:OG1	2.17	0.44
1:B:355:SER:HA	1:B:372:LYS:H	1.82	0.44
1:B:735:VAL:HG12	1:B:741:ILE:HD13	2.00	0.44
1:C:96:ILE:O	1:C:100:LEU:HG	2.18	0.44
1:C:109:ILE:C	1:C:111:LEU:N	2.76	0.44
1:C:217:LYS:HZ1	1:C:236:GLU:HB2	1.82	0.44
1:C:327:LEU:HG	1:C:595:ILE:HG12	1.99	0.44
1:C:413:LEU:CB	1:C:419:ILE:HG12	2.43	0.44
1:C:653:LYS:C	1:C:655:ASN:N	2.76	0.44
1:C:661:ALA:HB2	2:Q:39:LEU:C	2.43	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:688:PHE:C	1:C:688:PHE:HD2	2.24	0.44
1:C:694:VAL:HG23	1:C:695:LYS:N	2.32	0.44
1:D:78:LYS:O	1:D:81:GLN:HB3	2.17	0.44
1:D:159:TYR:HD1	1:D:159:TYR:N	2.10	0.44
1:D:165:GLN:HE22	1:D:252:ASP:HB3	1.82	0.44
1:D:318:ILE:H	1:D:318:ILE:CD1	2.23	0.44
1:D:401:ILE:HG21	1:D:485:LEU:HB3	1.99	0.44
1:D:497:LEU:CD1	1:D:556:MET:HG2	2.41	0.44
1:D:554:LYS:O	1:D:557:LEU:N	2.51	0.44
1:E:182:ILE:O	1:E:187:SER:HB2	2.16	0.44
1:E:189:ASP:HB3	1:E:190:PRO:CD	2.47	0.44
1:E:207:ASP:CB	1:E:262:PRO:HG3	2.47	0.44
1:E:327:LEU:HG	1:E:595:ILE:HG12	1.98	0.44
1:E:434:LEU:HD22	1:E:445:ARG:HB3	1.99	0.44
1:E:661:ALA:HB2	2:S:39:LEU:C	2.43	0.44
1:E:741:ILE:H	1:E:741:ILE:HG12	1.60	0.44
1:E:747:ASN:O	1:E:750:GLN:N	2.50	0.44
1:E:792:VAL:CG1	1:E:796:ILE:HD11	2.47	0.44
1:F:120:LEU:CD1	1:F:120:LEU:O	2.66	0.44
1:F:220:LEU:HG	1:F:223:LYS:HB3	1.99	0.44
1:F:305:SER:C	1:F:307:LEU:N	2.76	0.44
1:F:323:ASN:ND2	1:F:598:PRO:HB2	2.32	0.44
1:F:395:GLU:CD	1:F:395:GLU:H	2.25	0.44
1:F:658:PRO:HG2	1:F:701:LEU:HD22	1.99	0.44
1:F:686:ASP:O	1:F:689:ALA:HB3	2.17	0.44
1:F:700:TYR:CE1	1:F:727:GLN:HB3	2.53	0.44
1:F:724:ARG:HG3	1:F:724:ARG:NH1	2.31	0.44
1:F:786:GLU:O	1:F:789:ASN:HB3	2.16	0.44
1:F:792:VAL:CG1	1:F:796:ILE:HD11	2.47	0.44
2:O:63:ILE:HD11	2:O:67:GLU:C	2.42	0.44
2:O:144:MET:HE1	2:O:145:MET:HE1	2.00	0.44
2:Q:144:MET:HE1	2:Q:145:MET:CE	2.47	0.44
2:S:144:MET:HE1	2:S:145:MET:CE	2.47	0.44
2:T:64:ASP:OD1	2:T:66:PRO:HD2	2.17	0.44
2:T:94:LYS:HE3	2:T:107:HIS:HD2	1.81	0.44
1:A:195:LEU:HD12	1:A:230:ILE:HD13	2.00	0.44
1:A:423:LYS:HG2	1:A:424:LYS:N	2.33	0.44
1:B:76:LEU:O	1:B:77:ASP:C	2.60	0.44
1:B:97:TYR:HD2	1:B:102:GLY:HA3	1.80	0.44
1:B:217:LYS:HB3	1:B:236:GLU:CD	2.42	0.44
1:C:201:ASP:CG	1:C:225:ILE:HD12	2.43	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:263:ASP:O	1:C:266:GLU:N	2.51	0.44
1:D:96:ILE:O	1:D:100:LEU:HG	2.18	0.44
1:D:120:LEU:CD1	1:D:120:LEU:O	2.66	0.44
1:D:165:GLN:HG2	1:D:251:PRO:CG	2.47	0.44
1:D:197:LYS:CA	1:D:197:LYS:HZ3	2.31	0.44
1:D:236:GLU:C	1:D:238:GLN:H	2.25	0.44
1:D:241:PHE:CD1	1:D:268:MET:HE1	2.53	0.44
1:E:120:LEU:CD1	1:E:120:LEU:O	2.66	0.44
1:E:395:GLU:CD	1:E:395:GLU:H	2.26	0.44
1:E:401:ILE:HG21	1:E:485:LEU:HB3	2.00	0.44
1:E:654:ILE:HG22	1:E:655:ASN:ND2	2.32	0.44
1:E:719:LYS:O	1:E:720:ILE:C	2.61	0.44
1:F:207:ASP:CB	1:F:262:PRO:HG3	2.48	0.44
1:F:236:GLU:C	1:F:238:GLN:H	2.25	0.44
1:F:308:VAL:CG2	1:F:336:THR:O	2.66	0.44
1:F:345:THR:CG2	1:F:491:ASP:HA	2.47	0.44
1:F:747:ASN:O	1:F:750:GLN:N	2.50	0.44
2:O:117:THR:O	2:O:118:ASP:C	2.60	0.44
2:S:68:PHE:O	2:S:70:THR:N	2.50	0.44
2:T:63:ILE:HD11	2:T:68:PHE:CA	2.47	0.44
2:T:117:THR:O	2:T:118:ASP:C	2.60	0.44
2:T:126:ARG:HG3	2:T:126:ARG:NH2	2.29	0.44
1:A:120:LEU:CD1	1:A:120:LEU:O	2.66	0.44
1:A:177:ILE:C	1:A:180:ASP:OD1	2.60	0.44
1:A:635:ILE:HD12	1:A:635:ILE:N	2.33	0.44
1:B:292:ARG:NE	1:B:617:LYS:HE3	2.33	0.44
1:B:423:LYS:HG2	1:B:424:LYS:N	2.33	0.44
1:B:639:ASN:HD22	1:B:639:ASN:N	2.13	0.44
1:B:700:TYR:CE1	1:B:727:GLN:HB3	2.53	0.44
1:B:748:TYR:O	1:B:751:TYR:N	2.46	0.44
1:B:753:LYS:O	1:B:754:GLU:C	2.61	0.44
1:C:90:PRO:C	1:C:92:ASP:N	2.72	0.44
1:C:384:ASN:ND2	1:C:465:LEU:HD21	2.33	0.44
1:C:587:PRO:C	1:C:588:GLU:HG2	2.43	0.44
1:C:657:ILE:CD1	1:C:701:LEU:HD23	2.48	0.44
1:C:736:LEU:HD21	1:C:750:GLN:HE22	1.83	0.44
1:D:97:TYR:C	1:D:99:GLU:N	2.73	0.44
1:D:109:ILE:C	1:D:111:LEU:N	2.76	0.44
1:D:492:TYR:HB2	1:D:575:VAL:HG22	1.99	0.44
1:D:764:LEU:O	1:D:766:HIS:N	2.50	0.44
1:E:165:GLN:HG2	1:E:251:PRO:CG	2.48	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:170:TYR:HA	1:E:173:ILE:CG2	2.47	0.44
1:E:323:ASN:ND2	1:E:598:PRO:CB	2.81	0.44
1:E:666:ASN:HB2	1:E:748:TYR:CZ	2.52	0.44
1:E:747:ASN:HD22	1:E:747:ASN:N	2.15	0.44
1:F:165:GLN:HG2	1:F:251:PRO:CG	2.48	0.44
1:F:327:LEU:HG	1:F:595:ILE:HG12	2.00	0.44
1:F:504:ILE:HD12	1:F:504:ILE:N	2.32	0.44
1:F:517:VAL:HB	1:F:525:LYS:NZ	2.25	0.44
1:F:587:PRO:C	1:F:588:GLU:HG2	2.43	0.44
1:F:653:LYS:C	1:F:655:ASN:N	2.75	0.44
1:F:776:LEU:HD11	1:F:793:PHE:HE1	1.82	0.44
2:O:76:MET:C	2:O:78:ASP:N	2.75	0.44
2:P:102:ALA:CB	2:P:125:ILE:HG13	2.42	0.44
2:Q:76:MET:C	2:Q:78:ASP:N	2.75	0.44
2:Q:126:ARG:HG3	2:Q:126:ARG:NH2	2.30	0.44
2:Q:146:THR:O	2:Q:147:ALA:C	2.57	0.44
2:R:86:ARG:O	2:R:86:ARG:CG	2.66	0.44
2:R:93:ASP:OD1	2:R:97:ASN:ND2	2.50	0.44
2:R:94:LYS:HB3	2:R:94:LYS:HZ3	1.80	0.44
2:S:65:PHE:CB	2:S:66:PRO:HD3	2.47	0.44
2:T:65:PHE:CB	2:T:66:PRO:HD3	2.47	0.44
1:A:140:ARG:HA	1:A:140:ARG:NE	2.33	0.44
1:A:220:LEU:HG	1:A:223:LYS:HB3	1.99	0.44
1:A:225:ILE:HA	1:A:229:PHE:CE2	2.47	0.44
1:A:249:PHE:O	1:A:250:ALA:C	2.58	0.44
1:A:355:SER:HA	1:A:372:LYS:H	1.83	0.44
1:B:120:LEU:CD1	1:B:120:LEU:O	2.66	0.44
1:B:207:ASP:CB	1:B:262:PRO:HG3	2.48	0.44
1:B:661:ALA:HB2	2:P:39:LEU:C	2.43	0.44
1:B:747:ASN:O	1:B:750:GLN:N	2.50	0.44
1:C:165:GLN:HE22	1:C:252:ASP:HB3	1.81	0.44
1:C:192:PHE:H	1:C:192:PHE:HD1	1.65	0.44
1:C:307:LEU:HD12	1:C:331:VAL:CG2	2.48	0.44
1:C:405:LEU:HD13	1:C:453:VAL:CG2	2.45	0.44
1:C:609:GLU:O	1:C:610:MET:C	2.61	0.44
1:C:662:GLU:OE2	1:C:755:ARG:NH2	2.42	0.44
1:D:130:SER:O	1:D:131:ARG:C	2.61	0.44
1:D:408:LEU:HD12	1:D:408:LEU:N	2.02	0.44
1:D:688:PHE:C	1:D:688:PHE:HD2	2.24	0.44
1:E:71:PHE:CG	1:E:73:ASN:HB2	2.52	0.44
1:E:241:PHE:CD1	1:E:268:MET:HE1	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:279:ILE:O	1:E:283:LEU:HB2	2.17	0.44
1:E:345:THR:CG2	1:E:491:ASP:HA	2.47	0.44
1:E:423:LYS:HG2	1:E:424:LYS:N	2.33	0.44
1:E:576:ASN:N	1:E:576:ASN:ND2	2.65	0.44
1:E:657:ILE:CD1	1:E:701:LEU:HD23	2.47	0.44
1:E:694:VAL:O	1:E:695:LYS:C	2.60	0.44
1:E:719:LYS:C	1:E:721:SER:N	2.76	0.44
1:F:76:LEU:O	1:F:77:ASP:C	2.60	0.44
1:F:109:ILE:C	1:F:111:LEU:N	2.76	0.44
1:F:192:PHE:H	1:F:192:PHE:HD1	1.65	0.44
1:F:199:LEU:C	1:F:201:ASP:N	2.70	0.44
1:F:247:TYR:O	1:F:253:HIS:O	2.35	0.44
1:F:333:LYS:HA	1:F:336:THR:OG1	2.17	0.44
1:F:736:LEU:HD21	1:F:750:GLN:HE22	1.83	0.44
2:P:30:LYS:HD3	2:P:30:LYS:N	2.13	0.44
2:P:97:ASN:O	2:P:99:TYR:HD1	2.01	0.44
2:Q:58:ASP:O	2:Q:60:ASN:N	2.46	0.44
2:Q:144:MET:HE1	2:Q:145:MET:HE1	2.00	0.44
2:R:104:GLU:O	2:R:108:VAL:HG23	2.17	0.44
2:R:117:THR:O	2:R:118:ASP:C	2.60	0.44
2:T:97:ASN:O	2:T:99:TYR:HD1	2.01	0.44
1:A:263:ASP:O	1:A:266:GLU:N	2.50	0.44
1:A:540:ARG:NH1	1:A:627:TYR:CE1	2.85	0.44
1:B:100:LEU:CD1	1:B:182:ILE:HG21	2.48	0.44
1:B:109:ILE:O	1:B:109:ILE:HG22	2.17	0.44
1:B:397:GLU:O	1:B:480:ASN:N	2.50	0.44
1:B:576:ASN:N	1:B:576:ASN:ND2	2.62	0.44
1:B:666:ASN:HB2	1:B:748:TYR:CZ	2.53	0.44
1:C:320:ARG:HG2	1:C:599:GLU:HA	2.00	0.44
1:C:576:ASN:ND2	1:C:576:ASN:N	2.63	0.44
1:C:718:ARG:HH12	1:C:767:GLN:HE21	1.66	0.44
1:D:105:TYR:HB2	1:D:153:ILE:HG12	1.99	0.44
1:E:220:LEU:HG	1:E:223:LYS:HB3	1.99	0.44
1:E:247:TYR:O	1:E:253:HIS:O	2.36	0.44
1:E:288:VAL:HG12	1:E:292:ARG:HH22	1.83	0.44
1:E:384:ASN:ND2	1:E:465:LEU:HD21	2.32	0.44
1:E:397:GLU:O	1:E:480:ASN:N	2.51	0.44
1:E:554:LYS:O	1:E:557:LEU:N	2.51	0.44
1:E:776:LEU:HD11	1:E:793:PHE:HE1	1.82	0.44
1:F:217:LYS:HB2	1:F:236:GLU:CD	2.43	0.44
1:F:241:PHE:CD1	1:F:268:MET:HE1	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:384:ASN:ND2	1:F:465:LEU:HD21	2.33	0.44
2:P:13:LYS:O	2:P:15:ALA:N	2.50	0.44
2:P:68:PHE:O	2:P:70:THR:N	2.51	0.44
2:P:117:THR:O	2:P:118:ASP:C	2.60	0.44
2:Q:89:PHE:HB2	2:Q:141:PHE:CE2	2.53	0.44
2:Q:97:ASN:C	2:Q:99:TYR:N	2.76	0.44
2:R:12:PHE:HZ	2:R:76:MET:HE3	1.83	0.44
2:S:84:GLU:N	2:S:84:GLU:OE2	2.51	0.44
2:T:97:ASN:HD22	2:T:97:ASN:N	2.15	0.44
1:B:83:GLN:O	1:B:85:LEU:N	2.51	0.43
1:B:384:ASN:ND2	1:B:465:LEU:HD21	2.32	0.43
1:B:454:GLN:HG2	1:B:473:ASN:HA	2.00	0.43
1:B:481:VAL:HG21	1:B:486:LYS:HB2	1.99	0.43
1:B:635:ILE:HD12	1:B:635:ILE:N	2.32	0.43
1:B:661:ALA:HB2	2:P:38:SER:O	2.18	0.43
1:C:71:PHE:C	1:C:73:ASN:H	2.26	0.43
1:C:401:ILE:HG21	1:C:485:LEU:HB3	1.99	0.43
1:D:140:ARG:HA	1:D:140:ARG:NE	2.33	0.43
1:D:152:LEU:HG	1:D:154:ILE:HD11	2.00	0.43
1:D:207:ASP:CB	1:D:262:PRO:HG3	2.47	0.43
1:D:285:LYS:C	1:D:287:GLY:N	2.75	0.43
1:D:747:ASN:HD22	1:D:747:ASN:N	2.16	0.43
1:E:263:ASP:O	1:E:266:GLU:N	2.51	0.43
1:E:325:TYR:CD1	1:E:598:PRO:HD3	2.52	0.43
1:E:355:SER:HA	1:E:372:LYS:H	1.83	0.43
1:E:554:LYS:O	1:E:555:GLN:C	2.59	0.43
1:E:726:ILE:O	1:E:727:GLN:C	2.60	0.43
1:F:288:VAL:HG12	1:F:292:ARG:HH22	1.83	0.43
1:F:355:SER:HA	1:F:372:LYS:H	1.83	0.43
1:F:401:ILE:HG21	1:F:485:LEU:HB3	1.99	0.43
1:F:492:TYR:HB2	1:F:575:VAL:HG22	1.99	0.43
1:F:699:GLY:O	1:F:703:ASP:N	2.51	0.43
2:Q:36:MET:O	2:Q:37:ARG:C	2.59	0.43
2:Q:91:VAL:HG12	2:Q:92:PHE:N	2.32	0.43
2:R:68:PHE:O	2:R:70:THR:N	2.51	0.43
2:S:101:SER:OG	2:S:104:GLU:HG2	2.18	0.43
1:A:207:ASP:CB	1:A:262:PRO:HG3	2.48	0.43
1:A:661:ALA:HB2	2:O:39:LEU:C	2.43	0.43
1:A:666:ASN:HB2	1:A:748:TYR:CZ	2.53	0.43
1:A:735:VAL:HG12	1:A:741:ILE:HD13	2.00	0.43
1:B:152:LEU:HG	1:B:154:ILE:HD11	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:165:GLN:HE22	1:B:252:ASP:HB3	1.81	0.43
1:B:240:ALA:C	1:B:242:SER:N	2.76	0.43
1:B:320:ARG:HG2	1:B:599:GLU:HA	2.00	0.43
1:C:76:LEU:O	1:C:77:ASP:C	2.60	0.43
1:C:333:LYS:O	1:C:335:ALA:N	2.47	0.43
1:C:453:VAL:HG12	1:C:454:GLN:N	2.32	0.43
1:C:666:ASN:HB2	1:C:748:TYR:CZ	2.53	0.43
1:C:735:VAL:HG12	1:C:741:ILE:HD13	2.00	0.43
1:D:209:LEU:HD23	1:D:260:TYR:HA	1.99	0.43
1:D:255:THR:C	1:D:257:LEU:N	2.74	0.43
1:D:384:ASN:ND2	1:D:465:LEU:HD21	2.33	0.43
1:D:481:VAL:HG21	1:D:486:LYS:HB2	1.99	0.43
1:E:76:LEU:O	1:E:77:ASP:C	2.60	0.43
1:E:387:ASN:ND2	1:E:387:ASN:H	2.16	0.43
1:E:587:PRO:C	1:E:588:GLU:HG2	2.42	0.43
1:E:633:ASN:O	1:E:642:TYR:CE1	2.70	0.43
1:F:694:VAL:O	1:F:695:LYS:C	2.61	0.43
2:O:137:ASN:H	2:O:140:GLU:HG3	1.83	0.43
2:P:89:PHE:HB2	2:P:141:PHE:CE2	2.53	0.43
2:R:102:ALA:CB	2:R:125:ILE:HG13	2.41	0.43
2:T:49:GLN:O	2:T:53:ASN:N	2.42	0.43
1:B:263:ASP:O	1:B:266:GLU:N	2.51	0.43
1:B:305:SER:C	1:B:307:LEU:N	2.76	0.43
1:B:587:PRO:C	1:B:588:GLU:HG2	2.42	0.43
1:C:83:GLN:O	1:C:85:LEU:N	2.51	0.43
1:C:120:LEU:CD1	1:C:120:LEU:O	2.66	0.43
1:C:234:LEU:O	1:C:238:GLN:HG3	2.18	0.43
1:C:236:GLU:C	1:C:238:GLN:H	2.24	0.43
1:C:255:THR:HG22	1:C:258:GLU:OE2	2.18	0.43
1:C:395:GLU:CD	1:C:395:GLU:H	2.26	0.43
1:C:492:TYR:HB2	1:C:575:VAL:HG22	1.99	0.43
1:C:530:THR:O	1:C:534:ILE:HG13	2.17	0.43
1:C:686:ASP:O	1:C:689:ALA:HB3	2.19	0.43
1:C:691:LYS:NZ	2:Q:20:ASP:O	2.50	0.43
1:D:182:ILE:C	1:D:187:SER:HB2	2.43	0.43
1:D:195:LEU:HD12	1:D:230:ILE:HD13	2.00	0.43
1:D:320:ARG:HG2	1:D:599:GLU:HA	2.00	0.43
1:D:355:SER:HA	1:D:372:LYS:H	1.83	0.43
1:D:530:THR:O	1:D:534:ILE:HG13	2.18	0.43
1:D:700:TYR:CE1	1:D:727:GLN:HB3	2.54	0.43
1:E:105:TYR:O	1:E:154:ILE:N	2.35	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:131:ARG:HG3	1:E:243:LEU:CD2	2.47	0.43
1:E:195:LEU:HD12	1:E:230:ILE:HD13	2.00	0.43
1:E:540:ARG:HD2	1:E:582:ASP:OD1	2.18	0.43
1:E:615:ILE:HD12	1:E:645:TRP:HH2	1.84	0.43
1:E:630:ARG:CD	2:S:83:GLU:HG2	2.48	0.43
1:E:667:LEU:C	1:E:669:SER:N	2.74	0.43
1:E:747:ASN:O	1:E:748:TYR:C	2.61	0.43
1:F:89:ILE:H	1:F:89:ILE:HG13	1.47	0.43
1:F:115:LYS:HE2	1:F:116:GLU:HG2	1.98	0.43
1:F:140:ARG:HA	1:F:140:ARG:NE	2.33	0.43
2:O:39:LEU:HD23	2:O:39:LEU:HA	1.85	0.43
2:O:65:PHE:H	2:O:65:PHE:HD1	1.64	0.43
2:O:86:ARG:O	2:O:86:ARG:CG	2.65	0.43
2:P:76:MET:C	2:P:78:ASP:N	2.74	0.43
2:R:137:ASN:H	2:R:140:GLU:HG3	1.83	0.43
2:S:37:ARG:HH11	2:S:37:ARG:CG	2.31	0.43
2:T:101:SER:OG	2:T:104:GLU:HG2	2.18	0.43
1:A:96:ILE:O	1:A:100:LEU:HG	2.18	0.43
1:A:234:LEU:O	1:A:238:GLN:HG3	2.18	0.43
1:B:182:ILE:C	1:B:187:SER:HB2	2.44	0.43
1:B:318:ILE:H	1:B:318:ILE:CD1	2.23	0.43
1:B:323:ASN:ND2	1:B:598:PRO:CB	2.82	0.43
1:B:353:LYS:CB	1:B:368:GLN:HE22	2.32	0.43
1:C:423:LYS:HG2	1:C:424:LYS:N	2.33	0.43
1:D:76:LEU:O	1:D:77:ASP:C	2.60	0.43
1:D:263:ASP:O	1:D:266:GLU:N	2.51	0.43
1:D:499:PRO:CG	1:D:504:ILE:HD11	2.47	0.43
1:D:540:ARG:NH1	1:D:627:TYR:CE1	2.87	0.43
1:D:747:ASN:O	1:D:748:TYR:C	2.61	0.43
1:E:109:ILE:C	1:E:111:LEU:N	2.76	0.43
1:E:109:ILE:O	1:E:109:ILE:HG22	2.19	0.43
1:E:152:LEU:HG	1:E:154:ILE:HD11	1.99	0.43
1:E:197:LYS:CB	1:E:197:LYS:HZ3	2.30	0.43
1:E:408:LEU:CD1	1:E:408:LEU:N	2.73	0.43
1:E:694:VAL:CG2	1:E:695:LYS:N	2.80	0.43
1:F:73:ASN:O	1:F:74:GLU:C	2.58	0.43
1:F:667:LEU:C	1:F:669:SER:N	2.74	0.43
2:Q:117:THR:C	2:Q:119:GLU:N	2.73	0.43
2:S:94:LYS:HE3	2:S:107:HIS:HD2	1.84	0.43
2:T:16:PHE:CZ	2:T:27:ILE:HG23	2.54	0.43
1:A:165:GLN:HG2	1:A:251:PRO:CG	2.47	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:115:LYS:HE2	1:B:116:GLU:HG2	1.99	0.43
1:B:241:PHE:CD1	1:B:268:MET:HE1	2.53	0.43
1:B:530:THR:HG21	2:P:145:MET:CE	2.49	0.43
1:B:662:GLU:OE2	1:B:755:ARG:NH2	2.42	0.43
1:B:741:ILE:H	1:B:741:ILE:HG12	1.59	0.43
1:B:747:ASN:O	1:B:748:TYR:C	2.61	0.43
1:C:376:GLN:HB3	1:C:379:ALA:CB	2.49	0.43
1:C:462:ILE:CG1	1:C:463:THR:N	2.74	0.43
1:D:297:LYS:HB3	1:D:297:LYS:HZ3	1.83	0.43
1:D:297:LYS:CE	1:D:601:GLU:OE1	2.67	0.43
1:D:687:GLU:HA	1:D:687:GLU:OE2	2.19	0.43
1:D:699:GLY:O	1:D:703:ASP:N	2.51	0.43
1:D:735:VAL:HG12	1:D:741:ILE:HD13	1.99	0.43
1:E:376:GLN:HB3	1:E:379:ALA:CB	2.48	0.43
1:E:492:TYR:HB2	1:E:575:VAL:HG22	2.00	0.43
1:E:735:VAL:HG12	1:E:741:ILE:HD13	2.00	0.43
1:F:97:TYR:CE1	1:F:178:SER:HB2	2.53	0.43
1:F:657:ILE:CD1	1:F:701:LEU:HD23	2.48	0.43
2:O:97:ASN:ND2	2:O:97:ASN:N	2.66	0.43
2:O:97:ASN:C	2:O:99:TYR:N	2.76	0.43
2:P:117:THR:C	2:P:119:GLU:N	2.73	0.43
2:R:89:PHE:HB2	2:R:141:PHE:CE2	2.54	0.43
2:T:138:TYR:CZ	2:T:142:VAL:CG2	3.02	0.43
1:A:302:LEU:HB2	1:A:602:PHE:CD1	2.37	0.43
1:A:718:ARG:HH12	1:A:767:GLN:HE21	1.67	0.43
1:B:165:GLN:HG2	1:B:251:PRO:CG	2.47	0.43
1:B:234:LEU:O	1:B:238:GLN:HG3	2.17	0.43
1:B:395:GLU:CD	1:B:395:GLU:H	2.26	0.43
1:B:687:GLU:HA	1:B:687:GLU:OE2	2.18	0.43
1:B:772:GLU:O	1:B:773:PHE:C	2.61	0.43
1:B:776:LEU:HD11	1:B:793:PHE:HE1	1.82	0.43
1:C:97:TYR:C	1:C:99:GLU:N	2.73	0.43
1:C:325:TYR:CD1	1:C:598:PRO:HD3	2.54	0.43
1:C:504:ILE:HD12	1:C:504:ILE:N	2.33	0.43
1:C:667:LEU:C	1:C:669:SER:N	2.74	0.43
1:D:71:PHE:CG	1:D:73:ASN:HB2	2.52	0.43
1:D:109:ILE:HG22	1:D:109:ILE:O	2.19	0.43
1:D:255:THR:HG22	1:D:258:GLU:OE2	2.18	0.43
1:D:353:LYS:H	1:D:368:GLN:HE22	1.66	0.43
1:D:610:MET:O	1:D:614:PHE:N	2.40	0.43
1:E:184:LYS:HE2	1:E:193:LEU:CD1	2.31	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:186:LYS:HD2	1:E:234:LEU:HD12	2.00	0.43
1:E:635:ILE:HD12	1:E:635:ILE:N	2.33	0.43
1:E:694:VAL:C	1:E:696:LYS:N	2.76	0.43
1:F:96:ILE:O	1:F:100:LEU:HG	2.18	0.43
1:F:152:LEU:HG	1:F:154:ILE:HD11	1.99	0.43
1:F:666:ASN:HB2	1:F:748:TYR:CZ	2.53	0.43
2:O:37:ARG:HH11	2:O:37:ARG:CG	2.31	0.43
2:O:84:GLU:N	2:O:84:GLU:OE2	2.51	0.43
2:R:99:TYR:CD2	2:R:135:GLN:NE2	2.86	0.43
2:S:138:TYR:CZ	2:S:142:VAL:CG2	3.02	0.43
1:A:105:TYR:HB2	1:A:153:ILE:HG12	2.01	0.43
1:A:152:LEU:HG	1:A:154:ILE:HD11	2.00	0.43
1:A:615:ILE:HD12	1:A:645:TRP:HH2	1.83	0.43
1:B:78:LYS:O	1:B:81:GLN:HB3	2.19	0.43
1:B:554:LYS:O	1:B:555:GLN:C	2.58	0.43
1:C:78:LYS:O	1:C:81:GLN:HB3	2.18	0.43
1:C:217:LYS:HB2	1:C:236:GLU:CD	2.44	0.43
1:C:241:PHE:CD1	1:C:268:MET:HE1	2.53	0.43
1:C:345:THR:CG2	1:C:491:ASP:HA	2.48	0.43
1:C:353:LYS:H	1:C:368:GLN:HE22	1.66	0.43
1:C:409:ARG:CD	1:C:413:LEU:HD21	2.47	0.43
1:D:175:LYS:O	1:D:177:ILE:N	2.51	0.43
1:D:345:THR:CG2	1:D:491:ASP:HA	2.48	0.43
1:D:376:GLN:HB3	1:D:379:ALA:CB	2.48	0.43
1:D:635:ILE:HD12	1:D:635:ILE:N	2.34	0.43
1:D:747:ASN:O	1:D:750:GLN:N	2.50	0.43
1:E:105:TYR:HB2	1:E:153:ILE:HG12	2.01	0.43
1:E:741:ILE:O	1:E:742:ALA:C	2.61	0.43
1:F:195:LEU:HD12	1:F:230:ILE:HD13	2.00	0.43
1:F:231:LYS:O	1:F:232:GLU:C	2.62	0.43
1:F:409:ARG:CD	1:F:413:LEU:HD21	2.47	0.43
1:F:741:ILE:O	1:F:742:ALA:C	2.62	0.43
2:R:39:LEU:HD23	2:R:39:LEU:HA	1.86	0.43
2:S:12:PHE:HZ	2:S:76:MET:HE3	1.83	0.43
2:S:97:ASN:C	2:S:99:TYR:N	2.75	0.43
2:T:68:PHE:O	2:T:70:THR:N	2.52	0.43
2:T:97:ASN:ND2	2:T:97:ASN:N	2.66	0.43
2:T:137:ASN:H	2:T:140:GLU:HG3	1.83	0.43
1:A:473:ASN:OD1	1:A:473:ASN:N	2.50	0.43
1:A:700:TYR:CE1	1:A:727:GLN:HB3	2.53	0.43
1:A:776:LEU:HD11	1:A:793:PHE:HE1	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:554:LYS:O	1:B:557:LEU:N	2.51	0.43
1:B:654:ILE:HG22	1:B:655:ASN:ND2	2.33	0.43
1:B:667:LEU:C	1:B:669:SER:N	2.75	0.43
1:C:184:LYS:HE2	1:C:193:LEU:CD1	2.31	0.43
1:C:670:ILE:O	1:C:671:ARG:C	2.61	0.43
1:C:699:GLY:O	1:C:703:ASP:N	2.51	0.43
1:D:97:TYR:O	1:D:99:GLU:N	2.51	0.43
1:D:100:LEU:CD1	1:D:182:ILE:HG21	2.48	0.43
1:D:353:LYS:CB	1:D:368:GLN:HE22	2.32	0.43
1:D:514:ASP:O	1:D:516:VAL:N	2.52	0.43
1:E:165:GLN:HE22	1:E:252:ASP:HB3	1.80	0.43
1:E:231:LYS:O	1:E:232:GLU:C	2.62	0.43
1:E:450:ASN:ND2	1:E:452:GLU:HG3	2.34	0.43
1:E:495:PHE:CD1	1:E:495:PHE:O	2.72	0.43
1:E:686:ASP:O	1:E:689:ALA:HB3	2.18	0.43
1:F:234:LEU:CD2	1:F:235:THR:HG23	2.49	0.43
1:F:234:LEU:O	1:F:238:GLN:HG3	2.18	0.43
2:O:12:PHE:HZ	2:O:76:MET:HE3	1.84	0.43
2:R:7:GLU:C	2:R:9:ILE:H	2.27	0.43
2:R:58:ASP:O	2:R:60:ASN:N	2.45	0.43
1:A:182:ILE:C	1:A:187:SER:HB2	2.44	0.43
1:A:384:ASN:ND2	1:A:465:LEU:HD21	2.33	0.43
1:A:540:ARG:HD2	1:A:582:ASP:OD1	2.19	0.43
1:A:554:LYS:O	1:A:555:GLN:C	2.60	0.43
1:A:699:GLY:O	1:A:703:ASP:N	2.51	0.43
1:B:794:GLN:HE22	1:B:795:LYS:HG3	1.83	0.43
1:C:165:GLN:HG2	1:C:251:PRO:CG	2.47	0.43
1:C:481:VAL:HG21	1:C:486:LYS:HB2	2.00	0.43
1:D:217:LYS:HB2	1:D:236:GLU:CD	2.44	0.43
1:D:294:ASP:O	1:D:610:MET:SD	2.76	0.43
1:D:719:LYS:O	1:D:720:ILE:C	2.61	0.43
1:D:753:LYS:O	1:D:754:GLU:C	2.62	0.43
1:D:776:LEU:HD11	1:D:793:PHE:HE1	1.83	0.43
1:E:140:ARG:HA	1:E:140:ARG:NE	2.33	0.43
1:F:83:GLN:O	1:F:85:LEU:N	2.52	0.43
1:F:109:ILE:HG22	1:F:109:ILE:O	2.18	0.43
1:F:175:LYS:O	1:F:177:ILE:N	2.51	0.43
1:F:353:LYS:H	1:F:368:GLN:HE22	1.67	0.43
1:F:530:THR:O	1:F:534:ILE:HG13	2.18	0.43
1:F:610:MET:O	1:F:614:PHE:N	2.40	0.43
2:P:97:ASN:ND2	2:P:97:ASN:N	2.66	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:P:109:MET:O	2:P:114:GLU:HB3	2.19	0.43
2:T:144:MET:HE1	2:T:145:MET:HE1	2.00	0.43
1:A:294:ASP:O	1:A:610:MET:SD	2.76	0.43
1:B:97:TYR:O	1:B:99:GLU:N	2.51	0.43
1:B:231:LYS:O	1:B:232:GLU:C	2.62	0.43
1:C:140:ARG:HA	1:C:140:ARG:NE	2.33	0.43
1:C:285:LYS:C	1:C:287:GLY:N	2.75	0.43
1:C:333:LYS:HA	1:C:336:THR:OG1	2.18	0.43
1:C:420:LEU:C	1:C:420:LEU:HD13	2.44	0.43
1:C:639:ASN:HD22	1:C:639:ASN:N	2.13	0.43
1:C:694:VAL:C	1:C:696:LYS:N	2.75	0.43
1:C:753:LYS:O	1:C:754:GLU:C	2.62	0.43
1:D:86:LEU:C	1:D:88:LYS:N	2.77	0.43
1:D:313:ASP:O	1:D:316:LYS:HB3	2.19	0.43
1:D:409:ARG:CD	1:D:413:LEU:HD21	2.47	0.43
1:D:473:ASN:OD1	1:D:473:ASN:N	2.50	0.43
1:D:546:LYS:CD	1:D:554:LYS:HE2	2.39	0.43
1:D:657:ILE:CD1	1:D:701:LEU:HD23	2.49	0.43
1:D:686:ASP:O	1:D:689:ALA:HB3	2.18	0.43
1:D:719:LYS:C	1:D:721:SER:N	2.76	0.43
1:E:254:ARG:HB3	1:E:254:ARG:HH11	1.83	0.43
1:E:255:THR:HG22	1:E:258:GLU:OE2	2.18	0.43
1:E:527:LYS:HG3	2:S:145:MET:HA	2.01	0.43
1:E:719:LYS:O	1:E:722:ILE:N	2.52	0.43
1:F:97:TYR:HD2	1:F:102:GLY:HA3	1.77	0.43
2:O:102:ALA:CB	2:O:125:ILE:HG13	2.42	0.43
2:Q:12:PHE:HZ	2:Q:76:MET:HE3	1.84	0.43
2:Q:117:THR:O	2:Q:118:ASP:C	2.60	0.43
2:R:65:PHE:HB2	2:R:66:PRO:CD	2.49	0.43
2:S:48:LEU:HA	2:S:51:MET:CE	2.41	0.43
2:S:86:ARG:O	2:S:86:ARG:CG	2.66	0.43
2:T:86:ARG:O	2:T:86:ARG:CG	2.66	0.43
1:A:97:TYR:O	1:A:99:GLU:N	2.51	0.42
1:A:217:LYS:HB2	1:A:236:GLU:CD	2.43	0.42
1:A:240:ALA:C	1:A:242:SER:N	2.76	0.42
1:A:611:THR:HG22	1:A:615:ILE:CD1	2.47	0.42
1:A:653:LYS:C	1:A:655:ASN:N	2.77	0.42
1:A:719:LYS:O	1:A:720:ILE:C	2.62	0.42
1:B:115:LYS:HE3	1:B:116:GLU:HG2	1.97	0.42
1:B:131:ARG:HG3	1:B:243:LEU:CD2	2.48	0.42
1:B:387:ASN:ND2	1:B:387:ASN:H	2.17	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:697:ILE:C	1:B:699:GLY:N	2.73	0.42
1:B:741:ILE:O	1:B:742:ALA:C	2.62	0.42
1:C:217:LYS:HZ1	1:C:233:ASN:CB	2.32	0.42
1:C:231:LYS:O	1:C:232:GLU:C	2.62	0.42
1:C:648:PRO:O	1:C:649:ILE:C	2.60	0.42
1:D:223:LYS:O	1:D:224:SER:C	2.60	0.42
1:D:295:VAL:HB	1:D:603:ILE:HG23	2.01	0.42
1:D:395:GLU:CD	1:D:395:GLU:H	2.26	0.42
1:D:741:ILE:O	1:D:742:ALA:C	2.62	0.42
1:E:153:ILE:C	1:E:154:ILE:CG1	2.92	0.42
1:E:368:GLN:HG3	1:E:383:GLY:HA3	2.01	0.42
1:E:514:ASP:C	1:E:516:VAL:N	2.77	0.42
1:E:514:ASP:O	1:E:516:VAL:N	2.51	0.42
1:E:688:PHE:C	1:E:688:PHE:HD2	2.25	0.42
1:E:700:TYR:CE1	1:E:727:GLN:HB3	2.54	0.42
1:E:718:ARG:HH12	1:E:767:GLN:HE21	1.67	0.42
1:F:213:LYS:H	1:F:213:LYS:HG3	1.74	0.42
1:F:234:LEU:HD21	1:F:235:THR:HG23	2.01	0.42
1:F:336:THR:O	1:F:339:ILE:HB	2.19	0.42
1:F:514:ASP:C	1:F:516:VAL:N	2.77	0.42
1:F:520:PRO:HG2	1:F:521:ASN:H	1.84	0.42
2:O:36:MET:O	2:O:37:ARG:C	2.60	0.42
2:O:68:PHE:O	2:O:70:THR:N	2.52	0.42
2:O:89:PHE:HB2	2:O:141:PHE:CE2	2.54	0.42
2:O:109:MET:O	2:O:114:GLU:HB3	2.19	0.42
2:Q:101:SER:OG	2:Q:104:GLU:HG2	2.19	0.42
2:R:36:MET:O	2:R:37:ARG:C	2.60	0.42
2:R:97:ASN:ND2	2:R:97:ASN:N	2.66	0.42
2:S:24:ASP:OD1	2:S:25:GLY:N	2.47	0.42
2:T:12:PHE:HZ	2:T:76:MET:HE3	1.84	0.42
1:A:76:LEU:O	1:A:77:ASP:C	2.60	0.42
1:A:190:PRO:HB2	1:A:195:LEU:CD1	2.45	0.42
1:A:462:ILE:CD1	1:A:466:GLY:HA2	2.49	0.42
1:B:105:TYR:O	1:B:154:ILE:N	2.34	0.42
1:B:195:LEU:HD12	1:B:230:ILE:HD13	2.00	0.42
1:B:323:ASN:C	1:B:324:THR:HG22	2.44	0.42
1:B:719:LYS:C	1:B:721:SER:N	2.76	0.42
1:C:297:LYS:CB	1:C:297:LYS:HZ2	2.32	0.42
1:C:368:GLN:HG3	1:C:383:GLY:HA3	2.01	0.42
1:C:495:PHE:CD1	1:C:495:PHE:O	2.72	0.42
1:C:581:GLN:O	1:C:629:ASN:HA	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:108:ASP:HB3	1:D:109:ILE:H	1.72	0.42
1:D:456:LYS:HB3	1:D:471:TRP:N	2.34	0.42
1:E:353:LYS:H	1:E:368:GLN:HE22	1.67	0.42
1:E:609:GLU:O	1:E:613:ARG:N	2.34	0.42
1:E:639:ASN:HD22	1:E:639:ASN:N	2.13	0.42
1:F:182:ILE:C	1:F:187:SER:HB2	2.43	0.42
1:F:635:ILE:HD12	1:F:635:ILE:N	2.34	0.42
1:F:719:LYS:C	1:F:721:SER:N	2.75	0.42
2:O:52:ILE:C	2:O:54:GLU:N	2.77	0.42
2:P:12:PHE:HZ	2:P:76:MET:HE3	1.83	0.42
2:P:84:GLU:N	2:P:84:GLU:OE2	2.53	0.42
2:P:97:ASN:ND2	2:P:97:ASN:C	2.76	0.42
2:Q:18:LEU:HB3	2:Q:19:PHE:HD1	1.84	0.42
2:R:55:VAL:HB	2:R:67:GLU:OE1	2.19	0.42
2:S:16:PHE:CZ	2:S:27:ILE:HG23	2.54	0.42
2:S:55:VAL:HB	2:S:67:GLU:OE1	2.19	0.42
2:T:84:GLU:N	2:T:84:GLU:OE2	2.52	0.42
1:A:83:GLN:O	1:A:86:LEU:N	2.52	0.42
1:A:231:LYS:O	1:A:232:GLU:C	2.62	0.42
1:A:307:LEU:HD12	1:A:331:VAL:CG2	2.49	0.42
1:A:338:LEU:HD21	1:A:409:ARG:NH1	2.33	0.42
1:A:353:LYS:H	1:A:368:GLN:HE22	1.68	0.42
1:A:656:THR:HG22	1:A:657:ILE:O	2.19	0.42
1:B:88:LYS:NZ	1:B:172:GLU:CD	2.75	0.42
1:B:96:ILE:O	1:B:100:LEU:HG	2.18	0.42
1:B:254:ARG:HB3	1:B:254:ARG:HH11	1.83	0.42
1:B:462:ILE:CD1	1:B:466:GLY:HA2	2.50	0.42
1:B:514:ASP:O	1:B:516:VAL:N	2.52	0.42
1:B:530:THR:O	1:B:534:ILE:HG13	2.19	0.42
1:B:694:VAL:O	1:B:695:LYS:C	2.61	0.42
1:C:97:TYR:O	1:C:99:GLU:N	2.50	0.42
1:C:115:LYS:HE2	1:C:116:GLU:HG2	1.97	0.42
1:C:186:LYS:HD2	1:C:234:LEU:HD12	2.00	0.42
1:C:305:SER:C	1:C:307:LEU:N	2.76	0.42
1:C:469:PHE:CD1	1:C:469:PHE:C	2.98	0.42
1:C:556:MET:HE2	1:C:556:MET:HB2	1.86	0.42
1:C:719:LYS:O	1:C:720:ILE:C	2.61	0.42
1:D:130:SER:CB	1:D:170:TYR:CE2	3.02	0.42
1:D:234:LEU:CD2	1:D:235:THR:HG23	2.48	0.42
1:D:794:GLN:HE22	1:D:795:LYS:HG3	1.84	0.42
1:E:201:ASP:CG	1:E:225:ILE:HD12	2.44	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:217:LYS:HB2	1:E:236:GLU:CD	2.44	0.42
1:E:234:LEU:CD2	1:E:235:THR:HG23	2.49	0.42
1:E:530:THR:O	1:E:534:ILE:HG13	2.18	0.42
1:E:767:GLN:HG2	1:E:768:LYS:H	1.82	0.42
1:E:772:GLU:O	1:E:773:PHE:C	2.62	0.42
1:F:105:TYR:HB2	1:F:153:ILE:HG12	2.01	0.42
1:F:202:ASP:CB	1:F:208:LEU:HD23	2.49	0.42
1:F:333:LYS:O	1:F:335:ALA:N	2.49	0.42
1:F:420:LEU:C	1:F:420:LEU:HD13	2.45	0.42
2:P:7:GLU:C	2:P:9:ILE:H	2.26	0.42
2:P:97:ASN:C	2:P:99:TYR:N	2.76	0.42
2:P:99:TYR:CD2	2:P:135:GLN:NE2	2.88	0.42
2:Q:89:PHE:HB2	2:Q:141:PHE:CD2	2.54	0.42
2:Q:94:LYS:HE3	2:Q:107:HIS:HD2	1.82	0.42
2:R:109:MET:O	2:R:114:GLU:HB3	2.19	0.42
1:A:255:THR:HG22	1:A:258:GLU:OE2	2.19	0.42
1:A:365:PRO:C	1:A:367:ASP:N	2.75	0.42
1:A:554:LYS:O	1:A:557:LEU:N	2.52	0.42
1:B:123:GLU:O	1:B:146:LYS:NZ	2.53	0.42
1:B:140:ARG:HA	1:B:140:ARG:NE	2.33	0.42
1:C:100:LEU:CD1	1:C:182:ILE:HG21	2.49	0.42
1:C:195:LEU:HD12	1:C:230:ILE:HD13	2.01	0.42
1:C:240:ALA:C	1:C:242:SER:N	2.76	0.42
1:C:270:LYS:HA	1:C:273:LYS:CG	2.49	0.42
1:C:316:LYS:HG3	1:C:600:GLY:HA2	2.01	0.42
1:C:517:VAL:CB	1:C:525:LYS:HZ1	2.22	0.42
1:D:173:ILE:O	1:D:174:GLY:C	2.63	0.42
1:D:581:GLN:O	1:D:629:ASN:HA	2.20	0.42
1:D:667:LEU:C	1:D:669:SER:N	2.74	0.42
1:E:83:GLN:O	1:E:86:LEU:N	2.52	0.42
1:E:661:ALA:HB2	2:S:38:SER:O	2.19	0.42
1:F:254:ARG:HB3	1:F:254:ARG:HH11	1.84	0.42
1:F:372:LYS:CG	1:F:373:LYS:N	2.77	0.42
1:F:453:VAL:HG12	1:F:454:GLN:N	2.32	0.42
1:F:514:ASP:O	1:F:516:VAL:N	2.52	0.42
1:F:561:ASN:OD1	1:F:574:VAL:N	2.45	0.42
1:F:735:VAL:HG12	1:F:741:ILE:HD13	2.00	0.42
2:P:65:PHE:H	2:P:65:PHE:HD1	1.64	0.42
2:P:104:GLU:O	2:P:108:VAL:HG23	2.20	0.42
2:P:137:ASN:H	2:P:140:GLU:HG3	1.83	0.42
2:T:7:GLU:C	2:T:9:ILE:H	2.27	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:520:PRO:HG2	1:A:521:ASN:H	1.84	0.42
1:A:686:ASP:O	1:A:689:ALA:HB3	2.19	0.42
1:B:186:LYS:HD2	1:B:234:LEU:HD12	2.01	0.42
1:B:202:ASP:CB	1:B:208:LEU:HD23	2.48	0.42
1:B:297:LYS:HZ1	1:B:601:GLU:HB3	1.83	0.42
1:B:450:ASN:ND2	1:B:452:GLU:HG3	2.35	0.42
1:B:514:ASP:C	1:B:516:VAL:N	2.77	0.42
1:B:609:GLU:O	1:B:610:MET:C	2.62	0.42
1:C:337:ASN:C	1:C:339:ILE:N	2.72	0.42
1:C:353:LYS:CB	1:C:368:GLN:HE22	2.32	0.42
1:D:186:LYS:HD2	1:D:234:LEU:HD12	2.01	0.42
1:D:250:ALA:CB	1:D:253:HIS:CE1	3.02	0.42
1:D:333:LYS:C	1:D:335:ALA:N	2.78	0.42
1:D:333:LYS:HA	1:D:336:THR:OG1	2.18	0.42
1:D:360:VAL:O	1:D:361:ALA:C	2.61	0.42
1:D:368:GLN:C	1:D:370:LEU:N	2.77	0.42
1:D:412:GLU:O	1:D:416:ASN:HB2	2.20	0.42
1:E:320:ARG:HG2	1:E:599:GLU:HA	2.00	0.42
1:E:532:LEU:HA	1:E:532:LEU:HD23	1.78	0.42
1:E:691:LYS:NZ	2:S:20:ASP:O	2.47	0.42
1:F:64:ASN:HD22	1:F:64:ASN:N	2.18	0.42
1:F:165:GLN:HE22	1:F:252:ASP:HB3	1.81	0.42
1:F:297:LYS:NZ	1:F:297:LYS:CB	2.79	0.42
1:F:615:ILE:HD12	1:F:645:TRP:HH2	1.85	0.42
1:F:764:LEU:HD23	1:F:764:LEU:HA	1.88	0.42
1:F:772:GLU:O	1:F:773:PHE:C	2.62	0.42
2:S:97:ASN:O	2:S:99:TYR:HD1	2.03	0.42
1:A:109:ILE:O	1:A:109:ILE:HG22	2.19	0.42
1:A:115:LYS:HE2	1:A:116:GLU:HG2	1.98	0.42
1:A:210:PHE:C	1:A:212:GLN:N	2.72	0.42
1:A:254:ARG:HB3	1:A:254:ARG:HH11	1.83	0.42
1:A:255:THR:C	1:A:257:LEU:N	2.75	0.42
1:A:295:VAL:HB	1:A:603:ILE:HG23	2.01	0.42
1:A:412:GLU:O	1:A:416:ASN:HB2	2.20	0.42
1:A:469:PHE:C	1:A:469:PHE:CD1	2.98	0.42
1:A:694:VAL:O	1:A:695:LYS:C	2.61	0.42
1:A:719:LYS:C	1:A:721:SER:N	2.77	0.42
1:A:753:LYS:O	1:A:754:GLU:C	2.63	0.42
1:B:97:TYR:CE1	1:B:178:SER:HB2	2.54	0.42
1:B:527:LYS:HG3	2:P:145:MET:HA	2.02	0.42
1:B:670:ILE:O	1:B:671:ARG:C	2.63	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:691:LYS:NZ	2:P:20:ASP:O	2.48	0.42
1:C:202:ASP:CB	1:C:208:LEU:HD23	2.48	0.42
1:C:306:GLY:O	1:C:307:LEU:C	2.63	0.42
1:C:529:VAL:O	1:C:532:LEU:HB2	2.20	0.42
1:D:97:TYR:CE1	1:D:178:SER:HB2	2.54	0.42
1:D:323:ASN:ND2	1:D:598:PRO:CB	2.82	0.42
1:D:325:TYR:CD1	1:D:598:PRO:HD3	2.55	0.42
1:D:387:ASN:ND2	1:D:387:ASN:H	2.17	0.42
1:E:88:LYS:NZ	1:E:172:GLU:CD	2.73	0.42
1:E:196:ILE:O	1:E:199:LEU:N	2.53	0.42
1:E:288:VAL:O	1:E:290:LYS:N	2.52	0.42
1:F:184:LYS:HE2	1:F:193:LEU:CD1	2.31	0.42
2:O:99:TYR:CD2	2:O:135:GLN:NE2	2.88	0.42
2:Q:65:PHE:HB2	2:Q:66:PRO:CD	2.48	0.42
2:Q:68:PHE:C	2:Q:70:THR:H	2.27	0.42
2:Q:137:ASN:O	2:Q:138:TYR:C	2.61	0.42
2:R:101:SER:OG	2:R:104:GLU:HG2	2.20	0.42
2:S:68:PHE:C	2:S:70:THR:H	2.26	0.42
2:S:93:ASP:OD1	2:S:97:ASN:ND2	2.51	0.42
2:S:144:MET:HE1	2:S:145:MET:HE1	2.01	0.42
1:A:320:ARG:HG2	1:A:599:GLU:HA	2.01	0.42
1:A:353:LYS:CB	1:A:368:GLN:HE22	2.31	0.42
1:A:530:THR:HG21	2:O:145:MET:CE	2.50	0.42
1:A:581:GLN:O	1:A:629:ASN:HA	2.20	0.42
1:B:180:ASP:O	1:B:183:SER:N	2.37	0.42
1:B:294:ASP:O	1:B:610:MET:SD	2.78	0.42
1:B:469:PHE:CD1	1:B:469:PHE:C	2.98	0.42
1:C:234:LEU:HD21	1:C:235:THR:HG23	2.02	0.42
1:C:247:TYR:HB3	1:C:257:LEU:HB2	2.02	0.42
1:C:661:ALA:HB2	2:Q:38:SER:O	2.18	0.42
1:C:739:LYS:HG2	1:C:740:GLN:N	2.31	0.42
1:C:741:ILE:O	1:C:742:ALA:C	2.63	0.42
1:D:131:ARG:H	1:D:170:TYR:HE2	1.67	0.42
1:E:172:GLU:CB	1:E:246:SER:HA	2.47	0.42
1:F:323:ASN:ND2	1:F:598:PRO:CB	2.83	0.42
1:F:661:ALA:HB2	2:T:38:SER:O	2.19	0.42
2:O:7:GLU:C	2:O:9:ILE:H	2.27	0.42
2:O:101:SER:OG	2:O:104:GLU:HG2	2.19	0.42
2:P:97:ASN:HD22	2:P:97:ASN:N	2.17	0.42
2:Q:99:TYR:CD2	2:Q:135:GLN:NE2	2.88	0.42
2:R:24:ASP:OD1	2:R:25:GLY:N	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:S:89:PHE:HB2	2:S:141:PHE:CD2	2.54	0.42
2:T:18:LEU:HB3	2:T:19:PHE:HD1	1.85	0.42
1:A:186:LYS:HD2	1:A:234:LEU:HD12	2.01	0.42
1:A:247:TYR:HB3	1:A:257:LEU:HB2	2.02	0.42
1:A:387:ASN:ND2	1:A:387:ASN:H	2.17	0.42
1:A:687:GLU:HA	1:A:687:GLU:OE2	2.20	0.42
1:A:764:LEU:HD23	1:A:764:LEU:HA	1.87	0.42
1:B:105:TYR:HB2	1:B:153:ILE:HG12	2.01	0.42
1:B:175:LYS:HB2	1:B:175:LYS:HZ2	1.84	0.42
1:B:217:LYS:HB2	1:B:236:GLU:CD	2.44	0.42
1:B:353:LYS:H	1:B:368:GLN:HE22	1.68	0.42
1:B:412:GLU:C	1:B:414:LYS:N	2.78	0.42
1:B:687:GLU:O	1:B:690:LYS:HB2	2.20	0.42
1:C:207:ASP:CB	1:C:262:PRO:HG3	2.48	0.42
1:C:234:LEU:CD2	1:C:235:THR:HG23	2.49	0.42
1:C:365:PRO:C	1:C:367:ASP:N	2.76	0.42
1:C:435:LEU:HD23	1:C:435:LEU:HA	1.83	0.42
1:C:687:GLU:HA	1:C:687:GLU:OE2	2.20	0.42
1:C:794:GLN:HE22	1:C:795:LYS:HG3	1.84	0.42
1:D:609:GLU:O	1:D:610:MET:C	2.62	0.42
1:D:615:ILE:HD12	1:D:645:TRP:HH2	1.84	0.42
1:D:661:ALA:HB2	2:R:38:SER:O	2.20	0.42
1:D:772:GLU:O	1:D:773:PHE:C	2.63	0.42
1:E:353:LYS:CB	1:E:368:GLN:HE22	2.32	0.42
1:F:353:LYS:CB	1:F:368:GLN:HE22	2.32	0.42
1:F:368:GLN:HG3	1:F:383:GLY:HA3	2.01	0.42
1:F:438:ASN:HD22	1:F:438:ASN:HA	1.59	0.42
1:F:527:LYS:HG3	2:T:145:MET:HA	2.01	0.42
1:F:687:GLU:HA	1:F:687:GLU:OE2	2.20	0.42
2:P:13:LYS:HD2	2:P:13:LYS:HA	1.87	0.42
2:Q:97:ASN:ND2	2:Q:97:ASN:C	2.76	0.42
2:Q:102:ALA:HB1	2:Q:121:VAL:HG12	2.01	0.42
2:R:16:PHE:CE1	2:R:27:ILE:CG1	3.03	0.42
2:S:109:MET:O	2:S:114:GLU:HB3	2.19	0.42
2:T:97:ASN:O	2:T:99:TYR:CD1	2.73	0.42
1:A:514:ASP:C	1:A:516:VAL:N	2.78	0.42
1:A:609:GLU:O	1:A:610:MET:C	2.63	0.42
1:A:628:PHE:O	1:A:629:ASN:C	2.63	0.42
1:A:687:GLU:O	1:A:690:LYS:HB2	2.19	0.42
1:A:767:GLN:HG2	1:A:768:LYS:H	1.82	0.42
1:A:794:GLN:HE22	1:A:795:LYS:HG3	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:173:ILE:O	1:B:174:GLY:C	2.63	0.42
1:B:412:GLU:O	1:B:416:ASN:HB2	2.20	0.42
1:B:462:ILE:CG1	1:B:463:THR:N	2.74	0.42
1:B:615:ILE:HD12	1:B:645:TRP:HH2	1.85	0.42
1:C:197:LYS:CA	1:C:197:LYS:HZ3	2.33	0.42
1:C:254:ARG:HB3	1:C:254:ARG:HH11	1.83	0.42
1:C:464:VAL:HG23	1:C:465:LEU:N	2.35	0.42
1:C:667:LEU:HD13	1:C:678:VAL:HG21	2.02	0.42
1:D:462:ILE:CD1	1:D:466:GLY:HA2	2.49	0.42
1:E:250:ALA:CB	1:E:253:HIS:CE1	3.03	0.42
1:E:270:LYS:HA	1:E:273:LYS:CG	2.50	0.42
1:E:356:ASP:OD2	1:E:356:ASP:N	2.53	0.42
1:F:105:TYR:O	1:F:154:ILE:N	2.35	0.42
1:F:556:MET:HE2	1:F:556:MET:HB2	1.83	0.42
1:F:776:LEU:HD23	1:F:780:LEU:HD21	2.02	0.42
2:O:18:LEU:HB3	2:O:19:PHE:HD1	1.85	0.42
2:P:102:ALA:HB1	2:P:121:VAL:HG12	2.02	0.42
2:Q:16:PHE:CZ	2:Q:27:ILE:HG23	2.55	0.42
2:R:97:ASN:HD22	2:R:97:ASN:N	2.17	0.42
2:T:77:LYS:HG3	2:T:77:LYS:O	2.20	0.42
1:A:130:SER:O	1:A:131:ARG:C	2.62	0.42
1:A:615:ILE:HD12	1:A:645:TRP:CH2	2.55	0.42
1:A:711:ILE:HG13	1:A:712:PHE:CE2	2.55	0.42
1:B:100:LEU:CD2	1:B:182:ILE:HG21	2.50	0.42
1:B:297:LYS:CE	1:B:601:GLU:OE1	2.67	0.42
1:B:464:VAL:HG23	1:B:465:LEU:N	2.35	0.42
1:B:694:VAL:C	1:B:696:LYS:N	2.76	0.42
1:C:109:ILE:HD13	1:C:157:LYS:HZ3	1.84	0.42
1:C:153:ILE:C	1:C:154:ILE:CG1	2.93	0.42
1:D:89:ILE:H	1:D:89:ILE:HG13	1.49	0.42
1:D:153:ILE:C	1:D:154:ILE:CG1	2.92	0.42
1:D:172:GLU:CB	1:D:246:SER:HA	2.48	0.42
1:D:173:ILE:O	1:D:175:LYS:N	2.53	0.42
1:D:190:PRO:HB2	1:D:195:LEU:CD1	2.45	0.42
1:D:336:THR:O	1:D:339:ILE:HB	2.20	0.42
1:D:357:TRP:HA	1:D:418:ILE:HG23	2.02	0.42
1:D:420:LEU:HD13	1:D:420:LEU:C	2.45	0.42
1:D:469:PHE:C	1:D:469:PHE:CD1	2.98	0.42
1:D:656:THR:HG22	1:D:657:ILE:O	2.20	0.42
1:E:130:SER:O	1:E:131:ARG:C	2.62	0.42
1:E:202:ASP:CB	1:E:208:LEU:HD23	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:295:VAL:HB	1:E:603:ILE:HG23	2.02	0.42
1:E:776:LEU:HD23	1:E:780:LEU:HD21	2.02	0.42
1:F:84:ASP:HA	1:F:87:LYS:HG2	2.02	0.42
1:F:353:LYS:CG	1:F:368:GLN:HE22	2.33	0.42
1:F:629:ASN:CB	1:F:632:TYR:CE1	2.98	0.42
1:F:639:ASN:HD22	1:F:639:ASN:N	2.14	0.42
2:P:21:LYS:O	2:P:23:GLY:N	2.46	0.42
2:P:55:VAL:HB	2:P:67:GLU:OE1	2.20	0.42
2:P:77:LYS:HG3	2:P:77:LYS:O	2.19	0.42
2:P:89:PHE:HB2	2:P:141:PHE:CD2	2.55	0.42
2:P:97:ASN:O	2:P:99:TYR:CD1	2.73	0.42
2:R:131:ASP:OD1	2:R:131:ASP:N	2.53	0.42
2:S:137:ASN:H	2:S:140:GLU:HG3	1.84	0.42
2:T:48:LEU:HA	2:T:51:MET:CE	2.41	0.42
2:T:89:PHE:HD1	2:T:141:PHE:CD2	2.38	0.42
1:A:97:TYR:CE2	1:A:102:GLY:HA3	2.55	0.41
1:A:420:LEU:HD13	1:A:420:LEU:C	2.45	0.41
1:B:602:PHE:C	1:B:603:ILE:HG13	2.45	0.41
1:B:636:ALA:HA	1:B:637:PRO:HD3	1.87	0.41
1:C:311:HIS:O	1:C:312:ALA:C	2.63	0.41
1:C:357:TRP:HA	1:C:418:ILE:HG23	2.02	0.41
1:C:628:PHE:O	1:C:629:ASN:C	2.63	0.41
1:C:706:ASN:O	2:Q:130:ILE:HG23	2.20	0.41
1:C:747:ASN:O	1:C:748:TYR:C	2.61	0.41
1:D:231:LYS:O	1:D:232:GLU:C	2.63	0.41
1:D:234:LEU:HD21	1:D:235:THR:HG23	2.01	0.41
1:D:240:ALA:C	1:D:242:SER:N	2.77	0.41
1:D:254:ARG:HB3	1:D:254:ARG:HH11	1.83	0.41
1:D:397:GLU:O	1:D:480:ASN:N	2.51	0.41
1:D:450:ASN:ND2	1:D:452:GLU:OE1	2.53	0.41
1:D:628:PHE:O	1:D:629:ASN:C	2.63	0.41
1:D:712:PHE:CD1	1:D:716:LYS:HG2	2.55	0.41
1:E:78:LYS:O	1:E:81:GLN:HB3	2.20	0.41
1:E:84:ASP:HA	1:E:87:LYS:HG2	2.02	0.41
1:E:225:ILE:O	1:E:225:ILE:HG22	2.20	0.41
1:E:305:SER:C	1:E:307:LEU:N	2.75	0.41
1:E:480:ASN:C	1:E:481:VAL:CG2	2.93	0.41
1:E:753:LYS:O	1:E:754:GLU:C	2.63	0.41
1:E:767:GLN:CG	1:E:768:LYS:HG2	2.36	0.41
1:F:71:PHE:C	1:F:73:ASN:H	2.28	0.41
1:F:186:LYS:HD2	1:F:234:LEU:HD12	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:197:LYS:CA	1:F:197:LYS:HZ3	2.33	0.41
1:F:450:ASN:ND2	1:F:452:GLU:OE1	2.54	0.41
2:O:9:ILE:HD12	2:O:69:LEU:CD1	2.47	0.41
2:O:97:ASN:HD22	2:O:97:ASN:N	2.17	0.41
2:O:97:ASN:O	2:O:99:TYR:HD1	2.02	0.41
2:Q:32:LEU:HG	2:Q:36:MET:HE2	2.02	0.41
2:Q:71:MET:O	2:Q:71:MET:HG2	2.20	0.41
2:Q:97:ASN:O	2:Q:99:TYR:HD1	2.03	0.41
2:R:16:PHE:CZ	2:R:27:ILE:HG23	2.55	0.41
2:S:104:GLU:O	2:S:108:VAL:HG23	2.19	0.41
1:A:71:PHE:C	1:A:73:ASN:H	2.28	0.41
1:A:173:ILE:O	1:A:174:GLY:C	2.63	0.41
1:A:288:VAL:O	1:A:290:LYS:N	2.53	0.41
1:B:153:ILE:C	1:B:154:ILE:CG1	2.93	0.41
1:B:225:ILE:O	1:B:225:ILE:HG22	2.20	0.41
1:B:307:LEU:HD12	1:B:331:VAL:CG2	2.49	0.41
1:B:333:LYS:C	1:B:335:ALA:N	2.78	0.41
1:B:767:GLN:HG2	1:B:768:LYS:H	1.82	0.41
1:C:225:ILE:O	1:C:225:ILE:HG22	2.20	0.41
1:C:250:ALA:CB	1:C:253:HIS:CE1	3.03	0.41
1:C:514:ASP:O	1:C:516:VAL:N	2.54	0.41
1:D:247:TYR:HB3	1:D:257:LEU:HB2	2.02	0.41
1:D:639:ASN:ND2	1:D:639:ASN:N	2.61	0.41
1:D:660:SER:O	1:D:663:PHE:HB3	2.20	0.41
1:D:711:ILE:HG13	1:D:712:PHE:CE2	2.55	0.41
1:D:719:LYS:O	1:D:722:ILE:N	2.53	0.41
1:E:247:TYR:HB3	1:E:257:LEU:HB2	2.02	0.41
1:E:660:SER:O	1:E:663:PHE:HB3	2.19	0.41
1:F:180:ASP:O	1:F:183:SER:N	2.36	0.41
1:F:320:ARG:HG2	1:F:599:GLU:HA	2.00	0.41
1:F:323:ASN:C	1:F:324:THR:HG22	2.45	0.41
1:F:360:VAL:HG21	1:F:365:PRO:HB3	2.01	0.41
1:F:462:ILE:CG1	1:F:463:THR:N	2.74	0.41
1:F:469:PHE:CD1	1:F:469:PHE:C	2.98	0.41
1:F:602:PHE:C	1:F:603:ILE:HG13	2.45	0.41
1:F:688:PHE:O	1:F:689:ALA:C	2.62	0.41
1:F:694:VAL:C	1:F:696:LYS:N	2.76	0.41
2:P:16:PHE:CZ	2:P:27:ILE:HG23	2.54	0.41
2:Q:97:ASN:ND2	2:Q:97:ASN:N	2.65	0.41
2:Q:137:ASN:H	2:Q:140:GLU:HG3	1.84	0.41
2:R:12:PHE:CE1	2:R:72:MET:CE	2.94	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:S:12:PHE:CE1	2:S:72:MET:CE	2.93	0.41
2:S:71:MET:HG2	2:S:71:MET:O	2.20	0.41
2:S:94:LYS:HB3	2:S:94:LYS:HZ3	1.81	0.41
2:S:99:TYR:CD2	2:S:135:GLN:NE2	2.88	0.41
2:T:55:VAL:HB	2:T:67:GLU:OE1	2.20	0.41
2:T:65:PHE:H	2:T:65:PHE:HD1	1.65	0.41
1:A:86:LEU:C	1:A:88:LYS:N	2.78	0.41
1:A:153:ILE:C	1:A:154:ILE:CG1	2.92	0.41
1:A:353:LYS:CG	1:A:368:GLN:HE22	2.33	0.41
1:B:247:TYR:HB3	1:B:257:LEU:HB2	2.02	0.41
1:B:316:LYS:HG3	1:B:600:GLY:HA2	2.02	0.41
1:B:336:THR:O	1:B:339:ILE:HB	2.20	0.41
1:B:456:LYS:HB3	1:B:471:TRP:N	2.36	0.41
1:B:719:LYS:O	1:B:720:ILE:C	2.62	0.41
1:B:747:ASN:HD22	1:B:747:ASN:N	2.17	0.41
1:B:765:THR:HA	1:B:769:SER:CB	2.50	0.41
1:C:412:GLU:O	1:C:416:ASN:HB2	2.20	0.41
1:C:719:LYS:C	1:C:721:SER:N	2.76	0.41
1:D:413:LEU:CB	1:D:419:ILE:HG12	2.44	0.41
1:D:602:PHE:C	1:D:603:ILE:HG13	2.45	0.41
1:E:83:GLN:O	1:E:85:LEU:N	2.53	0.41
1:E:100:LEU:CD1	1:E:182:ILE:HG21	2.51	0.41
1:E:255:THR:C	1:E:257:LEU:N	2.75	0.41
1:E:297:LYS:CE	1:E:601:GLU:OE1	2.68	0.41
1:E:307:LEU:HD12	1:E:331:VAL:CG2	2.50	0.41
1:E:469:PHE:CD1	1:E:469:PHE:C	2.98	0.41
1:F:225:ILE:O	1:F:225:ILE:HG22	2.20	0.41
1:F:250:ALA:CB	1:F:253:HIS:CE1	3.04	0.41
1:F:356:ASP:OD2	1:F:356:ASP:N	2.53	0.41
1:F:376:GLN:HB3	1:F:379:ALA:CB	2.48	0.41
1:F:387:ASN:ND2	1:F:387:ASN:H	2.18	0.41
1:F:409:ARG:CZ	1:F:413:LEU:CD2	2.95	0.41
1:F:462:ILE:CD1	1:F:466:GLY:HA2	2.49	0.41
1:F:480:ASN:C	1:F:481:VAL:CG2	2.92	0.41
2:O:65:PHE:CB	2:O:66:PRO:HD3	2.47	0.41
2:R:138:TYR:CZ	2:R:142:VAL:CG2	3.04	0.41
2:S:59:GLY:H	2:S:62:THR:CG2	2.33	0.41
2:S:97:ASN:ND2	2:S:97:ASN:C	2.75	0.41
1:A:250:ALA:CB	1:A:253:HIS:CE1	3.03	0.41
1:A:297:LYS:CE	1:A:601:GLU:OE1	2.69	0.41
1:A:306:GLY:O	1:A:307:LEU:C	2.64	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:357:TRP:HA	1:A:418:ILE:HG23	2.02	0.41
1:A:514:ASP:O	1:A:516:VAL:N	2.53	0.41
1:A:765:THR:HA	1:A:769:SER:CB	2.50	0.41
1:B:84:ASP:HA	1:B:87:LYS:HG2	2.02	0.41
1:B:108:ASP:HB3	1:B:109:ILE:H	1.71	0.41
1:B:129:ASN:N	1:B:129:ASN:ND2	2.64	0.41
1:B:130:SER:O	1:B:131:ARG:C	2.63	0.41
1:B:234:LEU:CD2	1:B:235:THR:HG23	2.50	0.41
1:B:581:GLN:O	1:B:629:ASN:HA	2.20	0.41
1:B:686:ASP:O	1:B:689:ALA:HB3	2.20	0.41
1:B:712:PHE:CD1	1:B:716:LYS:HG2	2.55	0.41
1:C:320:ARG:HA	1:C:598:PRO:O	2.21	0.41
1:C:323:ASN:ND2	1:C:598:PRO:CB	2.83	0.41
1:C:456:LYS:HB3	1:C:471:TRP:N	2.35	0.41
1:C:660:SER:O	1:C:663:PHE:HB3	2.20	0.41
1:E:97:TYR:CE2	1:E:102:GLY:HA3	2.55	0.41
1:E:173:ILE:HG13	1:E:242:SER:CB	2.35	0.41
1:E:336:THR:O	1:E:339:ILE:HB	2.20	0.41
1:E:353:LYS:CG	1:E:368:GLN:HE22	2.34	0.41
1:E:615:ILE:HD12	1:E:645:TRP:CH2	2.56	0.41
1:E:694:VAL:HG23	1:E:695:LYS:N	2.34	0.41
1:F:97:TYR:CE2	1:F:102:GLY:HA3	2.55	0.41
1:F:129:ASN:N	1:F:129:ASN:ND2	2.64	0.41
1:F:153:ILE:C	1:F:154:ILE:CG1	2.93	0.41
1:F:164:GLU:O	1:F:166:SER:N	2.54	0.41
1:F:307:LEU:HD12	1:F:331:VAL:CG2	2.50	0.41
1:F:413:LEU:CB	1:F:419:ILE:HG12	2.43	0.41
1:F:662:GLU:OE2	1:F:755:ARG:NH2	2.42	0.41
2:O:65:PHE:HB2	2:O:66:PRO:CD	2.49	0.41
2:P:65:PHE:HB2	2:P:66:PRO:CD	2.49	0.41
2:Q:7:GLU:C	2:Q:9:ILE:H	2.28	0.41
2:Q:104:GLU:O	2:Q:108:VAL:HG23	2.20	0.41
2:R:89:PHE:HB2	2:R:141:PHE:CD2	2.55	0.41
2:S:102:ALA:HB1	2:S:121:VAL:HG12	2.03	0.41
1:A:159:TYR:HD1	1:A:159:TYR:N	2.10	0.41
1:A:313:ASP:O	1:A:316:LYS:HB3	2.20	0.41
1:A:323:ASN:C	1:A:324:THR:HG22	2.45	0.41
1:A:323:ASN:ND2	1:A:598:PRO:CB	2.84	0.41
1:A:336:THR:O	1:A:339:ILE:HB	2.20	0.41
1:A:397:GLU:O	1:A:480:ASN:N	2.51	0.41
1:A:450:ASN:ND2	1:A:452:GLU:OE1	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:495:PHE:CD1	1:A:495:PHE:O	2.73	0.41
1:A:527:LYS:HG3	2:O:145:MET:HA	2.02	0.41
1:B:109:ILE:C	1:B:111:LEU:N	2.76	0.41
1:B:112:VAL:C	1:B:114:HIS:N	2.72	0.41
1:B:196:ILE:C	1:B:198:SER:N	2.77	0.41
1:B:357:TRP:HA	1:B:418:ILE:HG23	2.02	0.41
1:B:420:LEU:C	1:B:420:LEU:HD13	2.44	0.41
1:B:520:PRO:HG2	1:B:521:ASN:H	1.84	0.41
1:B:718:ARG:HH12	1:B:767:GLN:HE21	1.68	0.41
1:B:767:GLN:CG	1:B:768:LYS:HG2	2.36	0.41
1:C:97:TYR:CE1	1:C:178:SER:HB2	2.56	0.41
1:C:313:ASP:O	1:C:316:LYS:HB3	2.21	0.41
1:C:368:GLN:C	1:C:370:LEU:N	2.78	0.41
1:C:469:PHE:HD1	1:C:469:PHE:C	2.29	0.41
1:C:493:ASP:OD2	1:C:577:HIS:HE1	2.01	0.41
1:C:633:ASN:O	1:C:642:TYR:CE1	2.72	0.41
1:D:311:HIS:O	1:D:312:ALA:C	2.64	0.41
1:D:609:GLU:N	1:D:609:GLU:CD	2.79	0.41
1:D:639:ASN:HD22	1:D:639:ASN:N	2.13	0.41
1:E:213:LYS:H	1:E:213:LYS:HG3	1.74	0.41
1:E:306:GLY:O	1:E:307:LEU:C	2.63	0.41
1:E:662:GLU:OE2	1:E:755:ARG:NH2	2.42	0.41
1:F:609:GLU:O	1:F:610:MET:C	2.63	0.41
2:P:48:LEU:HA	2:P:51:MET:CE	2.42	0.41
2:P:144:MET:HE1	2:P:145:MET:CE	2.49	0.41
2:R:68:PHE:C	2:R:70:THR:H	2.28	0.41
2:S:7:GLU:C	2:S:9:ILE:H	2.27	0.41
2:S:65:PHE:H	2:S:65:PHE:HD1	1.65	0.41
2:S:77:LYS:HG3	2:S:77:LYS:O	2.20	0.41
2:T:65:PHE:HB2	2:T:66:PRO:CD	2.49	0.41
2:T:68:PHE:C	2:T:70:THR:H	2.28	0.41
2:T:109:MET:O	2:T:114:GLU:HB3	2.20	0.41
1:A:271:LEU:HA	1:A:275:GLY:HA3	2.02	0.41
1:A:275:GLY:HA2	1:A:278:LYS:CE	2.50	0.41
1:A:480:ASN:HD22	1:A:480:ASN:C	2.29	0.41
1:A:530:THR:HG23	2:O:88:ALA:HB1	2.02	0.41
1:A:602:PHE:C	1:A:603:ILE:HG13	2.45	0.41
1:A:741:ILE:O	1:A:742:ALA:C	2.62	0.41
1:B:71:PHE:C	1:B:73:ASN:H	2.27	0.41
1:B:190:PRO:HB2	1:B:195:LEU:CD1	2.45	0.41
1:B:270:LYS:HA	1:B:273:LYS:CG	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:480:ASN:C	1:B:481:VAL:CG2	2.93	0.41
1:B:527:LYS:HE2	1:B:527:LYS:HB3	1.89	0.41
1:B:615:ILE:HD12	1:B:645:TRP:CH2	2.56	0.41
1:B:736:LEU:HD21	1:B:750:GLN:HE22	1.84	0.41
1:C:164:GLU:O	1:C:166:SER:N	2.54	0.41
1:C:173:ILE:O	1:C:174:GLY:C	2.63	0.41
1:C:196:ILE:O	1:C:199:LEU:N	2.54	0.41
1:C:217:LYS:CB	1:C:236:GLU:CG	2.98	0.41
1:C:387:ASN:ND2	1:C:387:ASN:H	2.18	0.41
1:C:602:PHE:C	1:C:603:ILE:HG13	2.45	0.41
1:C:609:GLU:N	1:C:609:GLU:CD	2.78	0.41
1:C:687:GLU:O	1:C:690:LYS:HB2	2.19	0.41
1:C:772:GLU:O	1:C:773:PHE:C	2.62	0.41
1:D:364:ILE:O	1:D:365:PRO:C	2.64	0.41
1:D:706:ASN:O	2:R:130:ILE:HG23	2.21	0.41
1:E:71:PHE:C	1:E:73:ASN:H	2.28	0.41
1:E:122:GLU:OE1	1:E:147:ARG:HB2	2.21	0.41
1:E:520:PRO:HG2	1:E:521:ASN:H	1.85	0.41
1:E:521:ASN:HB3	1:E:524:GLU:CB	2.51	0.41
1:E:561:ASN:OD1	1:E:574:VAL:N	2.46	0.41
1:E:610:MET:O	1:E:614:PHE:N	2.40	0.41
1:F:297:LYS:CE	1:F:601:GLU:OE1	2.69	0.41
1:F:316:LYS:HG3	1:F:600:GLY:HA2	2.03	0.41
1:F:456:LYS:HB3	1:F:471:TRP:N	2.35	0.41
1:F:741:ILE:H	1:F:741:ILE:HG12	1.60	0.41
2:O:18:LEU:HD23	2:O:18:LEU:HA	1.85	0.41
2:O:24:ASP:OD1	2:O:25:GLY:N	2.48	0.41
2:O:138:TYR:CZ	2:O:142:VAL:CG2	3.03	0.41
2:Q:94:LYS:HB3	2:Q:94:LYS:HZ2	1.86	0.41
2:S:138:TYR:O	2:S:139:GLU:C	2.64	0.41
2:T:59:GLY:H	2:T:62:THR:CG2	2.33	0.41
2:T:102:ALA:HB1	2:T:121:VAL:HG12	2.03	0.41
1:A:234:LEU:CD2	1:A:235:THR:HG23	2.50	0.41
1:A:456:LYS:HB3	1:A:471:TRP:N	2.35	0.41
1:A:772:GLU:O	1:A:773:PHE:C	2.61	0.41
1:B:97:TYR:CE2	1:B:102:GLY:HA3	2.56	0.41
1:B:234:LEU:HD21	1:B:235:THR:HG23	2.02	0.41
1:B:250:ALA:CB	1:B:253:HIS:CE1	3.03	0.41
1:B:255:THR:C	1:B:257:LEU:N	2.75	0.41
1:B:313:ASP:O	1:B:316:LYS:HB3	2.21	0.41
1:B:338:LEU:HD21	1:B:409:ARG:NE	2.36	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:353:LYS:CG	1:B:368:GLN:HE22	2.34	0.41
1:B:561:ASN:O	1:B:562:GLU:C	2.63	0.41
1:B:776:LEU:HD23	1:B:780:LEU:HD21	2.03	0.41
1:C:480:ASN:C	1:C:481:VAL:CG2	2.93	0.41
1:C:520:PRO:HG2	1:C:521:ASN:H	1.85	0.41
1:C:684:ASP:O	1:C:686:ASP:N	2.54	0.41
1:C:794:GLN:O	1:C:795:LYS:C	2.64	0.41
1:D:288:VAL:O	1:D:290:LYS:N	2.54	0.41
1:D:316:LYS:HG3	1:D:600:GLY:HA2	2.02	0.41
1:D:671:ARG:C	1:D:673:SER:N	2.79	0.41
1:E:142:VAL:CG1	1:E:154:ILE:HD12	2.47	0.41
1:E:182:ILE:C	1:E:187:SER:HB2	2.45	0.41
1:E:297:LYS:HB3	1:E:297:LYS:HZ3	1.84	0.41
1:E:316:LYS:HG3	1:E:600:GLY:HA2	2.03	0.41
1:E:462:ILE:CD1	1:E:466:GLY:HA2	2.50	0.41
1:E:546:LYS:CD	1:E:554:LYS:HE2	2.38	0.41
1:F:196:ILE:C	1:F:198:SER:N	2.77	0.41
1:F:254:ARG:H	1:F:254:ARG:CD	2.32	0.41
1:F:271:LEU:HA	1:F:275:GLY:HA3	2.03	0.41
1:F:313:ASP:O	1:F:316:LYS:HB3	2.21	0.41
1:F:712:PHE:CD1	1:F:716:LYS:HG2	2.55	0.41
2:O:32:LEU:HG	2:O:36:MET:HE2	2.03	0.41
2:O:89:PHE:HB2	2:O:141:PHE:CD2	2.56	0.41
2:O:97:ASN:O	2:O:99:TYR:CD1	2.74	0.41
2:Q:55:VAL:HB	2:Q:67:GLU:OE1	2.20	0.41
2:Q:97:ASN:O	2:Q:99:TYR:CD1	2.74	0.41
2:S:32:LEU:HG	2:S:36:MET:HE2	2.02	0.41
1:A:83:GLN:O	1:A:85:LEU:N	2.53	0.41
1:A:323:ASN:ND2	1:A:598:PRO:HB2	2.34	0.41
1:A:412:GLU:C	1:A:414:LYS:N	2.78	0.41
1:A:469:PHE:HD1	1:A:469:PHE:C	2.29	0.41
1:A:609:GLU:N	1:A:609:GLU:CD	2.78	0.41
1:A:671:ARG:C	1:A:673:SER:N	2.79	0.41
1:B:409:ARG:CZ	1:B:413:LEU:CD2	2.94	0.41
1:B:706:ASN:O	2:P:130:ILE:HG23	2.20	0.41
1:B:711:ILE:HG13	1:B:712:PHE:CE2	2.55	0.41
1:C:86:LEU:C	1:C:88:LYS:N	2.78	0.41
1:C:105:TYR:HB2	1:C:153:ILE:HG12	2.00	0.41
1:C:130:SER:CB	1:C:170:TYR:CE2	3.03	0.41
1:C:295:VAL:HB	1:C:603:ILE:HG23	2.02	0.41
1:C:462:ILE:CD1	1:C:466:GLY:HA2	2.49	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:88:LYS:NZ	1:D:172:GLU:CD	2.76	0.41
1:D:450:ASN:ND2	1:D:452:GLU:HG3	2.36	0.41
1:D:521:ASN:HB3	1:D:524:GLU:CB	2.51	0.41
1:D:532:LEU:HD23	1:D:532:LEU:HA	1.78	0.41
1:D:764:LEU:HD23	1:D:764:LEU:HA	1.88	0.41
1:D:765:THR:HA	1:D:769:SER:CB	2.50	0.41
1:E:252:ASP:CG	1:E:253:HIS:H	2.29	0.41
1:E:446:ILE:HG23	1:E:446:ILE:O	2.21	0.41
1:E:464:VAL:HG23	1:E:465:LEU:N	2.35	0.41
1:E:561:ASN:O	1:E:562:GLU:C	2.64	0.41
1:E:602:PHE:C	1:E:603:ILE:HG13	2.45	0.41
1:E:671:ARG:C	1:E:673:SER:N	2.79	0.41
1:F:196:ILE:O	1:F:199:LEU:N	2.54	0.41
1:F:247:TYR:HB3	1:F:257:LEU:HB2	2.03	0.41
1:F:306:GLY:O	1:F:307:LEU:C	2.63	0.41
1:F:311:HIS:O	1:F:312:ALA:C	2.63	0.41
1:F:320:ARG:HA	1:F:598:PRO:O	2.21	0.41
1:F:368:GLN:C	1:F:370:LEU:N	2.77	0.41
1:F:464:VAL:HG23	1:F:465:LEU:N	2.35	0.41
1:F:469:PHE:HD1	1:F:469:PHE:C	2.29	0.41
1:F:615:ILE:HD12	1:F:645:TRP:CH2	2.56	0.41
1:F:794:GLN:HE22	1:F:795:LYS:HG3	1.84	0.41
2:O:59:GLY:H	2:O:62:THR:CG2	2.34	0.41
2:Q:48:LEU:HA	2:Q:51:MET:CE	2.42	0.41
2:S:21:LYS:O	2:S:23:GLY:N	2.45	0.41
1:A:223:LYS:O	1:A:224:SER:C	2.64	0.41
1:A:225:ILE:O	1:A:225:ILE:HG22	2.20	0.41
1:A:316:LYS:HG3	1:A:600:GLY:HA2	2.02	0.41
1:A:464:VAL:HG23	1:A:465:LEU:N	2.36	0.41
1:A:530:THR:O	1:A:534:ILE:HG13	2.20	0.41
1:A:776:LEU:HD23	1:A:780:LEU:HD21	2.02	0.41
1:B:86:LEU:C	1:B:88:LYS:N	2.77	0.41
1:B:173:ILE:O	1:B:175:LYS:N	2.54	0.41
1:B:218:LEU:C	1:B:220:LEU:H	2.25	0.41
1:B:284:LYS:HE3	1:B:284:LYS:HA	2.03	0.41
1:B:469:PHE:HD1	1:B:469:PHE:C	2.29	0.41
1:B:492:TYR:HB2	1:B:575:VAL:CG2	2.51	0.41
1:B:495:PHE:CD1	1:B:495:PHE:O	2.74	0.41
1:B:520:PRO:HG2	1:B:521:ASN:N	2.36	0.41
1:B:656:THR:HG22	1:B:657:ILE:O	2.21	0.41
1:B:660:SER:O	1:B:663:PHE:HB3	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:84:ASP:HA	1:C:87:LYS:HG2	2.02	0.41
1:C:122:GLU:OE1	1:C:147:ARG:HB2	2.21	0.41
1:C:184:LYS:NZ	1:C:191:GLU:HG3	2.36	0.41
1:C:271:LEU:HA	1:C:275:GLY:HA3	2.03	0.41
1:C:275:GLY:HA2	1:C:278:LYS:CE	2.49	0.41
1:C:323:ASN:C	1:C:324:THR:HG22	2.46	0.41
1:C:741:ILE:H	1:C:741:ILE:HG12	1.59	0.41
1:C:765:THR:HA	1:C:769:SER:CB	2.50	0.41
1:C:776:LEU:HD23	1:C:780:LEU:HD21	2.02	0.41
1:D:122:GLU:OE1	1:D:147:ARG:HB2	2.21	0.41
1:D:164:GLU:O	1:D:166:SER:N	2.54	0.41
1:D:205:SER:C	1:D:207:ASP:N	2.79	0.41
1:D:270:LYS:HA	1:D:273:LYS:CG	2.50	0.41
1:D:320:ARG:NE	1:D:599:GLU:O	2.54	0.41
1:D:353:LYS:CG	1:D:368:GLN:HE22	2.34	0.41
1:D:435:LEU:HD23	1:D:435:LEU:HA	1.83	0.41
1:D:469:PHE:HD1	1:D:469:PHE:C	2.29	0.41
1:D:480:ASN:C	1:D:481:VAL:CG2	2.93	0.41
1:D:520:PRO:HG2	1:D:521:ASN:H	1.86	0.41
1:D:561:ASN:OD1	1:D:574:VAL:N	2.46	0.41
1:D:615:ILE:HD12	1:D:645:TRP:CH2	2.55	0.41
1:E:173:ILE:O	1:E:174:GLY:C	2.64	0.41
1:E:254:ARG:H	1:E:254:ARG:CD	2.32	0.41
1:E:294:ASP:O	1:E:610:MET:SD	2.79	0.41
1:E:333:LYS:C	1:E:335:ALA:N	2.78	0.41
1:E:432:TYR:CD2	1:E:447:SER:HA	2.56	0.41
1:E:628:PHE:O	1:E:629:ASN:C	2.64	0.41
1:E:653:LYS:C	1:E:655:ASN:N	2.76	0.41
1:E:687:GLU:OE2	1:E:687:GLU:HA	2.20	0.41
1:E:765:THR:HA	1:E:769:SER:CB	2.50	0.41
1:F:78:LYS:O	1:F:81:GLN:HB3	2.20	0.41
1:F:275:GLY:HA2	1:F:278:LYS:CE	2.50	0.41
1:F:357:TRP:HA	1:F:418:ILE:HG23	2.02	0.41
1:F:403:LEU:CG	1:F:405:LEU:HD11	2.51	0.41
1:F:412:GLU:C	1:F:414:LYS:N	2.78	0.41
1:F:412:GLU:O	1:F:416:ASN:HB2	2.20	0.41
1:F:529:VAL:O	1:F:532:LEU:HB2	2.20	0.41
1:F:530:THR:HG21	2:T:145:MET:CE	2.51	0.41
1:F:656:THR:HG22	1:F:657:ILE:O	2.21	0.41
1:F:706:ASN:O	2:T:130:ILE:HG23	2.20	0.41
2:O:91:VAL:HG12	2:O:92:PHE:N	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:P:71:MET:O	2:P:71:MET:HG2	2.20	0.41
2:Q:15:ALA:O	2:Q:35:VAL:HG13	2.21	0.41
2:Q:32:LEU:O	2:Q:33:GLY:C	2.64	0.41
2:Q:77:LYS:O	2:Q:77:LYS:HG3	2.20	0.41
2:R:63:ILE:HG13	2:R:67:GLU:HB2	2.02	0.41
2:R:73:ALA:HB1	2:R:77:LYS:HB2	2.03	0.41
2:S:18:LEU:HB3	2:S:19:PHE:HD1	1.84	0.41
2:S:39:LEU:HD23	2:S:39:LEU:HA	1.85	0.41
2:S:73:ALA:HB1	2:S:77:LYS:HB2	2.03	0.41
2:T:32:LEU:O	2:T:33:GLY:C	2.64	0.41
2:T:94:LYS:HB3	2:T:94:LYS:HZ3	1.85	0.41
2:T:99:TYR:CD2	2:T:135:GLN:NE2	2.89	0.41
2:T:104:GLU:O	2:T:108:VAL:HG23	2.20	0.41
2:T:138:TYR:O	2:T:139:GLU:C	2.64	0.41
1:A:84:ASP:HA	1:A:87:LYS:HG2	2.02	0.41
1:A:122:GLU:OE1	1:A:147:ARG:HB2	2.21	0.41
1:A:173:ILE:O	1:A:175:LYS:N	2.54	0.41
1:A:196:ILE:O	1:A:199:LEU:N	2.54	0.41
1:A:284:LYS:HE3	1:A:284:LYS:HA	2.03	0.41
1:A:435:LEU:HD23	1:A:435:LEU:HA	1.82	0.41
1:A:520:PRO:HG2	1:A:521:ASN:N	2.36	0.41
1:A:712:PHE:CD1	1:A:716:LYS:HG2	2.56	0.41
1:B:255:THR:HG22	1:B:258:GLU:OE2	2.20	0.41
1:B:295:VAL:HB	1:B:603:ILE:HG23	2.02	0.41
1:B:632:TYR:C	1:B:634:LYS:H	2.29	0.41
1:B:700:TYR:HD1	1:B:728:ALA:N	2.19	0.41
1:B:794:GLN:O	1:B:795:LYS:C	2.64	0.41
1:C:173:ILE:O	1:C:175:LYS:N	2.54	0.41
1:C:527:LYS:HG3	2:Q:145:MET:HA	2.02	0.41
1:D:71:PHE:C	1:D:73:ASN:H	2.28	0.41
1:D:83:GLN:O	1:D:85:LEU:N	2.53	0.41
1:D:97:TYR:CE2	1:D:102:GLY:HA3	2.55	0.41
1:D:368:GLN:HG3	1:D:383:GLY:HA3	2.02	0.41
1:D:517:VAL:HB	1:D:525:LYS:NZ	2.25	0.41
1:D:611:THR:HG22	1:D:615:ILE:CD1	2.47	0.41
1:D:687:GLU:O	1:D:690:LYS:HB2	2.20	0.41
1:E:97:TYR:CE1	1:E:178:SER:HB2	2.55	0.41
1:E:164:GLU:O	1:E:166:SER:N	2.54	0.41
1:E:403:LEU:CG	1:E:405:LEU:HD11	2.51	0.41
1:E:469:PHE:HD1	1:E:469:PHE:C	2.29	0.41
1:E:706:ASN:O	2:S:130:ILE:HG23	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:776:LEU:CD2	1:E:780:LEU:HD21	2.51	0.41
1:E:794:GLN:HE22	1:E:795:LYS:HG3	1.85	0.41
1:F:493:ASP:OD2	1:F:577:HIS:HE1	2.01	0.41
1:F:628:PHE:O	1:F:629:ASN:C	2.63	0.41
1:F:699:GLY:O	1:F:700:TYR:C	2.64	0.41
2:Q:140:GLU:O	2:Q:143:GLN:HB2	2.21	0.41
2:R:59:GLY:H	2:R:62:THR:CG2	2.34	0.41
2:S:97:ASN:O	2:S:99:TYR:CD1	2.74	0.41
2:S:138:TYR:CZ	2:S:142:VAL:HG21	2.56	0.41
2:T:16:PHE:CE1	2:T:27:ILE:CG1	3.04	0.41
1:A:109:ILE:C	1:A:111:LEU:N	2.76	0.40
1:A:202:ASP:CB	1:A:208:LEU:HD23	2.48	0.40
1:A:480:ASN:C	1:A:481:VAL:CG2	2.94	0.40
1:B:205:SER:C	1:B:207:ASP:N	2.79	0.40
1:C:297:LYS:CE	1:C:601:GLU:OE1	2.68	0.40
1:C:360:VAL:HG21	1:C:365:PRO:HB3	2.03	0.40
1:C:615:ILE:HD12	1:C:645:TRP:HH2	1.86	0.40
1:D:83:GLN:O	1:D:86:LEU:N	2.53	0.40
1:D:86:LEU:C	1:D:88:LYS:H	2.29	0.40
1:D:295:VAL:HB	1:D:603:ILE:CG2	2.51	0.40
1:D:323:ASN:C	1:D:324:THR:HG22	2.46	0.40
1:D:514:ASP:C	1:D:516:VAL:N	2.77	0.40
1:D:530:THR:HG21	2:R:145:MET:CE	2.51	0.40
1:D:776:LEU:CD2	1:D:780:LEU:HD21	2.51	0.40
1:E:313:ASP:O	1:E:316:LYS:HB3	2.20	0.40
1:E:420:LEU:HD13	1:E:420:LEU:C	2.45	0.40
1:E:632:TYR:C	1:E:634:LYS:H	2.29	0.40
1:E:688:PHE:O	1:E:689:ALA:C	2.63	0.40
1:F:86:LEU:C	1:F:88:LYS:N	2.77	0.40
1:F:217:LYS:HZ1	1:F:233:ASN:CB	2.34	0.40
1:F:252:ASP:CG	1:F:253:HIS:H	2.29	0.40
1:F:270:LYS:HA	1:F:273:LYS:CG	2.51	0.40
1:F:368:GLN:CB	1:F:380:VAL:HG13	2.51	0.40
1:F:520:PRO:HG2	1:F:521:ASN:N	2.36	0.40
1:F:671:ARG:C	1:F:673:SER:N	2.79	0.40
1:F:776:LEU:C	1:F:776:LEU:CD2	2.95	0.40
2:O:71:MET:O	2:O:71:MET:HG2	2.20	0.40
2:O:73:ALA:HB1	2:O:77:LYS:HB2	2.03	0.40
2:P:144:MET:HE1	2:P:145:MET:HE1	2.04	0.40
2:Q:59:GLY:H	2:Q:62:THR:CG2	2.34	0.40
2:Q:65:PHE:H	2:Q:65:PHE:HD1	1.64	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:T:32:LEU:HD21	2:T:71:MET:HE1	2.03	0.40
1:A:270:LYS:HA	1:A:273:LYS:CG	2.51	0.40
1:A:320:ARG:HA	1:A:598:PRO:O	2.22	0.40
1:B:196:ILE:O	1:B:199:LEU:N	2.54	0.40
1:B:223:LYS:O	1:B:224:SER:C	2.64	0.40
1:B:252:ASP:CG	1:B:253:HIS:H	2.29	0.40
1:C:97:TYR:CE2	1:C:102:GLY:HA3	2.55	0.40
1:C:166:SER:O	1:C:169:VAL:HG12	2.21	0.40
1:C:172:GLU:CB	1:C:246:SER:HA	2.48	0.40
1:C:223:LYS:O	1:C:224:SER:C	2.65	0.40
1:C:492:TYR:HB2	1:C:575:VAL:CG2	2.51	0.40
1:C:611:THR:HG22	1:C:615:ILE:CD1	2.47	0.40
1:C:671:ARG:C	1:C:673:SER:N	2.79	0.40
1:D:529:VAL:O	1:D:532:LEU:HB2	2.21	0.40
1:D:561:ASN:O	1:D:562:GLU:C	2.63	0.40
1:D:615:ILE:HG13	1:D:615:ILE:H	1.60	0.40
1:E:323:ASN:C	1:E:324:THR:HG22	2.46	0.40
1:E:368:GLN:CB	1:E:380:VAL:HG13	2.51	0.40
1:E:456:LYS:HG2	1:E:471:TRP:CE2	2.57	0.40
1:E:530:THR:HG21	2:S:145:MET:CE	2.51	0.40
1:E:551:ASN:HD22	1:E:551:ASN:HA	1.65	0.40
1:F:492:TYR:HB2	1:F:575:VAL:CG2	2.51	0.40
1:F:765:THR:HA	1:F:769:SER:CB	2.50	0.40
2:P:73:ALA:HB1	2:P:77:LYS:HB2	2.03	0.40
2:Q:86:ARG:O	2:Q:86:ARG:CG	2.67	0.40
2:R:32:LEU:HG	2:R:36:MET:HE2	2.04	0.40
2:S:16:PHE:HE1	2:S:27:ILE:CD1	2.34	0.40
2:S:65:PHE:HB2	2:S:66:PRO:CD	2.49	0.40
2:S:66:PRO:C	2:S:68:PHE:H	2.29	0.40
1:A:191:GLU:C	1:A:193:LEU:N	2.79	0.40
1:A:234:LEU:HD21	1:A:235:THR:HG23	2.02	0.40
1:A:252:ASP:CG	1:A:253:HIS:H	2.30	0.40
1:A:376:GLN:HB3	1:A:379:ALA:CB	2.48	0.40
1:A:424:LYS:HZ2	1:A:424:LYS:CB	2.33	0.40
1:A:561:ASN:HA	1:A:564:VAL:CG2	2.52	0.40
1:A:776:LEU:CD2	1:A:780:LEU:HD21	2.51	0.40
1:B:275:GLY:HA2	1:B:278:LYS:CE	2.50	0.40
1:B:306:GLY:HA3	1:B:331:VAL:O	2.22	0.40
1:B:611:THR:HG22	1:B:615:ILE:CD1	2.48	0.40
1:B:628:PHE:O	1:B:629:ASN:C	2.64	0.40
1:C:205:SER:C	1:C:207:ASP:N	2.79	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:353:LYS:CG	1:C:368:GLN:HE22	2.34	0.40
1:C:409:ARG:CZ	1:C:413:LEU:CD2	2.94	0.40
1:C:688:PHE:O	1:C:689:ALA:C	2.62	0.40
1:C:732:ILE:HG23	1:C:749:PHE:CD1	2.53	0.40
1:C:793:PHE:O	1:C:794:GLN:C	2.64	0.40
1:D:115:LYS:HE2	1:D:116:GLU:HG2	1.98	0.40
1:D:284:LYS:HE3	1:D:284:LYS:HA	2.03	0.40
1:D:582:ASP:O	1:D:584:GLU:N	2.53	0.40
1:D:632:TYR:C	1:D:634:LYS:H	2.29	0.40
1:D:671:ARG:HD2	2:R:14:GLU:HG3	2.03	0.40
1:D:776:LEU:C	1:D:776:LEU:CD2	2.95	0.40
1:E:217:LYS:HZ1	1:E:233:ASN:CB	2.35	0.40
1:E:234:LEU:HD21	1:E:235:THR:HG23	2.02	0.40
1:E:271:LEU:HA	1:E:275:GLY:HA3	2.04	0.40
1:E:357:TRP:HA	1:E:418:ILE:HG23	2.03	0.40
1:E:667:LEU:HD13	1:E:678:VAL:HG21	2.04	0.40
1:E:687:GLU:O	1:E:690:LYS:HB2	2.21	0.40
1:F:295:VAL:HB	1:F:603:ILE:HG23	2.02	0.40
1:F:776:LEU:CD2	1:F:780:LEU:HD21	2.51	0.40
2:P:16:PHE:HE1	2:P:27:ILE:CD1	2.35	0.40
2:P:59:GLY:H	2:P:62:THR:CG2	2.34	0.40
2:R:102:ALA:HB1	2:R:121:VAL:HG12	2.02	0.40
2:S:16:PHE:CE1	2:S:27:ILE:CG1	3.04	0.40
2:S:32:LEU:HD21	2:S:71:MET:HE1	2.03	0.40
2:S:49:GLN:O	2:S:50:ASP:C	2.65	0.40
2:T:51:MET:H	2:T:51:MET:HG3	1.67	0.40
1:A:73:ASN:HB3	1:A:74:GLU:H	1.74	0.40
1:A:218:LEU:C	1:A:220:LEU:H	2.24	0.40
1:A:306:GLY:HA3	1:A:331:VAL:O	2.21	0.40
1:A:660:SER:O	1:A:663:PHE:HB3	2.22	0.40
1:A:794:GLN:O	1:A:795:LYS:C	2.64	0.40
1:B:271:LEU:HA	1:B:275:GLY:HA3	2.03	0.40
1:B:288:VAL:O	1:B:290:LYS:N	2.54	0.40
1:B:368:GLN:HG3	1:B:383:GLY:HA3	2.02	0.40
1:B:732:ILE:HG23	1:B:749:PHE:CD1	2.53	0.40
1:C:323:ASN:C	1:C:324:THR:CG2	2.95	0.40
1:C:520:PRO:HG2	1:C:521:ASN:N	2.37	0.40
1:C:792:VAL:C	1:C:796:ILE:HG12	2.47	0.40
1:D:74:GLU:CD	1:D:74:GLU:H	2.30	0.40
1:D:186:LYS:O	1:D:187:SER:C	2.63	0.40
1:D:633:ASN:O	1:D:642:TYR:CE1	2.74	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:365:PRO:C	1:E:367:ASP:N	2.76	0.40
1:E:412:GLU:O	1:E:416:ASN:HB2	2.20	0.40
1:F:186:LYS:O	1:F:187:SER:C	2.65	0.40
1:F:191:GLU:C	1:F:193:LEU:N	2.80	0.40
1:F:295:VAL:HB	1:F:603:ILE:CG2	2.51	0.40
1:F:364:ILE:O	1:F:365:PRO:C	2.64	0.40
1:F:655:ASN:O	1:F:759:GLN:NE2	2.55	0.40
1:F:732:ILE:HG23	1:F:749:PHE:CD1	2.53	0.40
2:O:77:LYS:O	2:O:77:LYS:HG3	2.20	0.40
2:Q:41:GLN:HE22	2:Q:76:MET:HE1	1.86	0.40
2:Q:89:PHE:HD1	2:Q:141:PHE:CD2	2.39	0.40
2:Q:138:TYR:O	2:Q:139:GLU:C	2.65	0.40
2:R:18:LEU:HB3	2:R:19:PHE:HD1	1.85	0.40
2:R:77:LYS:O	2:R:77:LYS:HG3	2.21	0.40
2:R:97:ASN:O	2:R:99:TYR:CD1	2.74	0.40
2:R:140:GLU:O	2:R:143:GLN:HB2	2.22	0.40
1:A:295:VAL:HB	1:A:603:ILE:CG2	2.52	0.40
1:A:368:GLN:CB	1:A:380:VAL:HG13	2.52	0.40
1:A:403:LEU:CG	1:A:405:LEU:HD11	2.51	0.40
1:A:456:LYS:HG2	1:A:471:TRP:CE2	2.56	0.40
1:A:633:ASN:O	1:A:642:TYR:CE1	2.74	0.40
1:A:732:ILE:HG23	1:A:749:PHE:CD1	2.53	0.40
1:B:89:ILE:H	1:B:89:ILE:HG13	1.49	0.40
1:B:225:ILE:HG12	1:B:229:PHE:HE2	1.86	0.40
1:B:254:ARG:H	1:B:254:ARG:CD	2.33	0.40
1:B:432:TYR:CD2	1:B:447:SER:HA	2.56	0.40
1:B:518:ASN:ND2	1:B:518:ASN:N	2.70	0.40
1:B:561:ASN:HA	1:B:564:VAL:CG2	2.52	0.40
1:B:609:GLU:N	1:B:609:GLU:CD	2.79	0.40
1:B:793:PHE:O	1:B:794:GLN:C	2.65	0.40
1:C:521:ASN:HB3	1:C:524:GLU:CB	2.50	0.40
1:C:530:THR:HG21	2:Q:145:MET:CE	2.51	0.40
1:C:610:MET:HE3	1:C:610:MET:HB2	1.97	0.40
1:C:656:THR:HG22	1:C:657:ILE:O	2.22	0.40
1:C:712:PHE:CD1	1:C:716:LYS:HG2	2.56	0.40
1:D:196:ILE:O	1:D:199:LEU:N	2.54	0.40
1:D:210:PHE:C	1:D:212:GLN:N	2.74	0.40
1:D:218:LEU:C	1:D:220:LEU:H	2.24	0.40
1:D:320:ARG:HA	1:D:598:PRO:O	2.22	0.40
1:D:700:TYR:HD1	1:D:728:ALA:HA	1.86	0.40
1:E:412:GLU:C	1:E:414:LYS:N	2.79	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:520:PRO:HG2	1:E:521:ASN:N	2.37	0.40
1:E:700:TYR:HD1	1:E:728:ALA:N	2.20	0.40
1:E:794:GLN:O	1:E:795:LYS:C	2.65	0.40
1:F:166:SER:O	1:F:169:VAL:HG12	2.22	0.40
1:F:218:LEU:HD21	1:F:225:ILE:HD13	2.04	0.40
1:F:323:ASN:C	1:F:324:THR:CG2	2.94	0.40
1:F:581:GLN:O	1:F:629:ASN:HA	2.21	0.40
1:F:615:ILE:HG13	1:F:615:ILE:H	1.60	0.40
1:F:687:GLU:O	1:F:690:LYS:HB2	2.22	0.40
2:O:68:PHE:O	2:O:69:LEU:C	2.65	0.40
2:P:68:PHE:C	2:P:70:THR:H	2.28	0.40
2:Q:49:GLN:O	2:Q:50:ASP:C	2.64	0.40
2:Q:109:MET:O	2:Q:114:GLU:HB3	2.20	0.40
2:R:51:MET:H	2:R:51:MET:HG3	1.67	0.40
2:T:12:PHE:CE1	2:T:72:MET:CE	2.93	0.40
2:T:41:GLN:HE22	2:T:76:MET:HE1	1.86	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	733/747 (98%)	503 (69%)	145 (20%)	85 (12%)	0	1
1	B	733/747 (98%)	507 (69%)	142 (19%)	84 (12%)	0	2
1	C	733/747 (98%)	506 (69%)	145 (20%)	82 (11%)	0	2
1	D	733/747 (98%)	504 (69%)	144 (20%)	85 (12%)	0	1
1	E	733/747 (98%)	502 (68%)	146 (20%)	85 (12%)	0	1
1	F	733/747 (98%)	503 (69%)	146 (20%)	84 (12%)	0	2
2	O	144/149 (97%)	90 (62%)	39 (27%)	15 (10%)	0	2

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	P	144/149 (97%)	91 (63%)	39 (27%)	14 (10%)	0	3
2	Q	144/149 (97%)	90 (62%)	39 (27%)	15 (10%)	0	2
2	R	144/149 (97%)	90 (62%)	39 (27%)	15 (10%)	0	2
2	S	144/149 (97%)	92 (64%)	37 (26%)	15 (10%)	0	2
2	T	144/149 (97%)	92 (64%)	36 (25%)	16 (11%)	0	2
All	All	5262/5376 (98%)	3570 (68%)	1097 (21%)	595 (11%)	0	2

All (595) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	73	ASN
1	A	74	GLU
1	A	80	GLN
1	A	108	ASP
1	A	111	LEU
1	A	129	ASN
1	A	135	VAL
1	A	146	LYS
1	A	180	ASP
1	A	187	SER
1	A	278	LYS
1	A	377	GLN
1	A	485	LEU
1	A	787	THR
1	B	73	ASN
1	B	74	GLU
1	B	80	GLN
1	B	108	ASP
1	B	111	LEU
1	B	129	ASN
1	B	135	VAL
1	B	146	LYS
1	B	180	ASP
1	B	187	SER
1	B	278	LYS
1	B	377	GLN
1	B	485	LEU
1	B	787	THR
1	C	80	GLN
1	C	108	ASP

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Mol	Chain	Res	Type
1	C	111	LEU
1	C	129	ASN
1	C	135	VAL
1	C	146	LYS
1	C	180	ASP
1	C	187	SER
1	C	204	ASP
1	C	278	LYS
1	C	377	GLN
1	C	485	LEU
1	C	787	THR
1	D	73	ASN
1	D	74	GLU
1	D	80	GLN
1	D	108	ASP
1	D	111	LEU
1	D	129	ASN
1	D	135	VAL
1	D	146	LYS
1	D	180	ASP
1	D	187	SER
1	D	204	ASP
1	D	278	LYS
1	D	302	LEU
1	D	377	GLN
1	D	485	LEU
1	D	787	THR
1	E	73	ASN
1	E	74	GLU
1	E	80	GLN
1	E	108	ASP
1	E	111	LEU
1	E	129	ASN
1	E	135	VAL
1	E	146	LYS
1	E	180	ASP
1	E	187	SER
1	E	204	ASP
1	E	278	LYS
1	E	377	GLN
1	E	485	LEU
1	E	787	THR

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Mol	Chain	Res	Type
1	F	73	ASN
1	F	74	GLU
1	F	80	GLN
1	F	108	ASP
1	F	111	LEU
1	F	129	ASN
1	F	135	VAL
1	F	146	LYS
1	F	180	ASP
1	F	187	SER
1	F	278	LYS
1	F	302	LEU
1	F	377	GLN
1	F	485	LEU
1	F	787	THR
2	O	12	PHE
2	O	23	GLY
2	O	50	ASP
2	O	93	ASP
2	P	12	PHE
2	P	23	GLY
2	P	50	ASP
2	Q	12	PHE
2	Q	23	GLY
2	Q	50	ASP
2	Q	93	ASP
2	R	12	PHE
2	R	23	GLY
2	R	50	ASP
2	S	12	PHE
2	S	23	GLY
2	S	50	ASP
2	T	12	PHE
2	T	23	GLY
2	T	50	ASP
2	T	93	ASP
1	A	91	LYS
1	A	112	VAL
1	A	113	GLU
1	A	137	PHE
1	A	138	ALA
1	A	176	GLY

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Mol	Chain	Res	Type
1	A	192	PHE
1	A	200	SER
1	A	204	ASP
1	A	299	GLU
1	A	302	LEU
1	A	376	GLN
1	A	407	HIS
1	A	423	LYS
1	A	438	ASN
1	A	452	GLU
1	A	510	GLN
1	A	580	GLU
1	A	620	THR
1	A	672	ARG
1	A	711	ILE
1	A	714	GLN
1	A	779	GLN
1	A	783	THR
1	B	91	LYS
1	B	112	VAL
1	B	113	GLU
1	B	137	PHE
1	B	138	ALA
1	B	176	GLY
1	B	192	PHE
1	B	200	SER
1	B	204	ASP
1	B	299	GLU
1	B	302	LEU
1	B	376	GLN
1	B	406	ASP
1	B	407	HIS
1	B	423	LYS
1	B	438	ASN
1	B	452	GLU
1	B	510	GLN
1	B	580	GLU
1	B	620	THR
1	B	672	ARG
1	B	711	ILE
1	B	714	GLN
1	B	779	GLN

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Mol	Chain	Res	Type
1	B	783	THR
1	C	91	LYS
1	C	112	VAL
1	C	113	GLU
1	C	137	PHE
1	C	138	ALA
1	C	176	GLY
1	C	192	PHE
1	C	200	SER
1	C	299	GLU
1	C	302	LEU
1	C	376	GLN
1	C	407	HIS
1	C	423	LYS
1	C	438	ASN
1	C	452	GLU
1	C	510	GLN
1	C	580	GLU
1	C	620	THR
1	C	654	ILE
1	C	672	ARG
1	C	711	ILE
1	C	714	GLN
1	C	779	GLN
1	C	783	THR
1	D	91	LYS
1	D	112	VAL
1	D	113	GLU
1	D	137	PHE
1	D	138	ALA
1	D	176	GLY
1	D	192	PHE
1	D	200	SER
1	D	299	GLU
1	D	376	GLN
1	D	406	ASP
1	D	407	HIS
1	D	423	LYS
1	D	438	ASN
1	D	452	GLU
1	D	510	GLN
1	D	580	GLU

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Mol	Chain	Res	Type
1	D	620	THR
1	D	672	ARG
1	D	711	ILE
1	D	779	GLN
1	D	783	THR
1	E	91	LYS
1	E	112	VAL
1	E	113	GLU
1	E	137	PHE
1	E	138	ALA
1	E	176	GLY
1	E	192	PHE
1	E	200	SER
1	E	299	GLU
1	E	302	LEU
1	E	376	GLN
1	E	406	ASP
1	E	407	HIS
1	E	423	LYS
1	E	438	ASN
1	E	452	GLU
1	E	510	GLN
1	E	580	GLU
1	E	620	THR
1	E	672	ARG
1	E	711	ILE
1	E	779	GLN
1	E	783	THR
1	F	91	LYS
1	F	98	SER
1	F	112	VAL
1	F	113	GLU
1	F	137	PHE
1	F	138	ALA
1	F	176	GLY
1	F	192	PHE
1	F	200	SER
1	F	204	ASP
1	F	299	GLU
1	F	376	GLN
1	F	406	ASP
1	F	407	HIS

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Mol	Chain	Res	Type
1	F	423	LYS
1	F	438	ASN
1	F	452	GLU
1	F	510	GLN
1	F	580	GLU
1	F	620	THR
1	F	654	ILE
1	F	672	ARG
1	F	711	ILE
1	F	714	GLN
1	F	779	GLN
1	F	783	THR
2	O	28	THR
2	O	67	GLU
2	P	28	THR
2	P	67	GLU
2	Q	28	THR
2	Q	67	GLU
2	R	28	THR
2	R	67	GLU
2	S	28	THR
2	S	67	GLU
2	T	28	THR
2	T	67	GLU
1	A	77	ASP
1	A	98	SER
1	A	163	SER
1	A	166	SER
1	A	188	LEU
1	A	206	SER
1	A	232	GLU
1	A	289	GLU
1	A	361	ALA
1	A	372	LYS
1	A	373	LYS
1	A	406	ASP
1	A	654	ILE
1	A	669	SER
1	A	731	GLU
1	A	765	THR
1	A	794	GLN
1	B	77	ASP

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Mol	Chain	Res	Type
1	B	98	SER
1	B	163	SER
1	B	166	SER
1	B	188	LEU
1	B	206	SER
1	B	232	GLU
1	B	289	GLU
1	B	372	LYS
1	B	373	LYS
1	B	515	LYS
1	B	654	ILE
1	B	669	SER
1	B	731	GLU
1	B	765	THR
1	B	794	GLN
1	C	77	ASP
1	C	98	SER
1	C	163	SER
1	C	166	SER
1	C	188	LEU
1	C	206	SER
1	C	232	GLU
1	C	289	GLU
1	C	296	LEU
1	C	307	LEU
1	C	372	LYS
1	C	373	LYS
1	C	406	ASP
1	C	669	SER
1	C	731	GLU
1	C	794	GLN
1	D	77	ASP
1	D	98	SER
1	D	163	SER
1	D	166	SER
1	D	188	LEU
1	D	206	SER
1	D	232	GLU
1	D	289	GLU
1	D	361	ALA
1	D	372	LYS
1	D	373	LYS

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Mol	Chain	Res	Type
1	D	515	LYS
1	D	654	ILE
1	D	669	SER
1	D	714	GLN
1	D	720	ILE
1	D	731	GLU
1	D	794	GLN
1	E	77	ASP
1	E	98	SER
1	E	163	SER
1	E	166	SER
1	E	188	LEU
1	E	206	SER
1	E	232	GLU
1	E	289	GLU
1	E	307	LEU
1	E	361	ALA
1	E	372	LYS
1	E	373	LYS
1	E	515	LYS
1	E	654	ILE
1	E	669	SER
1	E	714	GLN
1	E	720	ILE
1	E	731	GLU
1	E	794	GLN
1	F	77	ASP
1	F	163	SER
1	F	166	SER
1	F	188	LEU
1	F	206	SER
1	F	232	GLU
1	F	289	GLU
1	F	307	LEU
1	F	372	LYS
1	F	373	LYS
1	F	515	LYS
1	F	669	SER
1	F	720	ILE
1	F	731	GLU
1	F	794	GLN
2	O	18	LEU

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Mol	Chain	Res	Type
2	O	25	GLY
2	O	71	MET
2	O	118	ASP
2	P	25	GLY
2	P	71	MET
2	P	118	ASP
2	Q	18	LEU
2	Q	25	GLY
2	Q	71	MET
2	Q	118	ASP
2	R	25	GLY
2	R	71	MET
2	R	118	ASP
2	S	18	LEU
2	S	25	GLY
2	S	71	MET
2	S	118	ASP
2	T	18	LEU
2	T	25	GLY
2	T	71	MET
2	T	118	ASP
1	A	110	ASP
1	A	116	GLU
1	A	165	GLN
1	A	191	GLU
1	A	274	GLY
1	A	290	LYS
1	A	296	LEU
1	A	307	LEU
1	A	357	TRP
1	A	481	VAL
1	A	515	LYS
1	A	568	GLY
1	A	605	THR
1	A	698	ALA
1	A	720	ILE
1	A	755	ARG
1	B	110	ASP
1	B	116	GLU
1	B	130	SER
1	B	165	GLN
1	B	191	GLU

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Mol	Chain	Res	Type
1	B	274	GLY
1	B	290	LYS
1	B	296	LEU
1	B	307	LEU
1	B	357	TRP
1	B	481	VAL
1	B	568	GLY
1	B	605	THR
1	B	698	ALA
1	B	720	ILE
1	B	755	ARG
1	C	75	THR
1	C	110	ASP
1	C	165	GLN
1	C	191	GLU
1	C	274	GLY
1	C	290	LYS
1	C	334	LEU
1	C	357	TRP
1	C	481	VAL
1	C	515	LYS
1	C	568	GLY
1	C	605	THR
1	C	668	SER
1	C	698	ALA
1	C	720	ILE
1	C	755	ARG
1	C	765	THR
1	D	110	ASP
1	D	116	GLU
1	D	165	GLN
1	D	191	GLU
1	D	274	GLY
1	D	290	LYS
1	D	296	LEU
1	D	307	LEU
1	D	334	LEU
1	D	357	TRP
1	D	481	VAL
1	D	568	GLY
1	D	605	THR
1	D	698	ALA

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Mol	Chain	Res	Type
1	D	755	ARG
1	D	765	THR
1	E	110	ASP
1	E	116	GLU
1	E	130	SER
1	E	165	GLN
1	E	191	GLU
1	E	274	GLY
1	E	290	LYS
1	E	296	LEU
1	E	334	LEU
1	E	357	TRP
1	E	481	VAL
1	E	568	GLY
1	E	605	THR
1	E	668	SER
1	E	698	ALA
1	E	765	THR
1	F	110	ASP
1	F	116	GLU
1	F	165	GLN
1	F	191	GLU
1	F	274	GLY
1	F	290	LYS
1	F	296	LEU
1	F	334	LEU
1	F	357	TRP
1	F	481	VAL
1	F	568	GLY
1	F	605	THR
1	F	698	ALA
1	F	755	ARG
1	F	765	THR
2	O	22	ASP
2	O	24	ASP
2	P	18	LEU
2	P	24	ASP
2	Q	24	ASP
2	R	18	LEU
2	R	22	ASP
2	R	24	ASP
2	S	22	ASP

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Mol	Chain	Res	Type
2	S	24	ASP
2	T	24	ASP
1	A	130	SER
1	A	239	HIS
1	A	334	LEU
1	A	484	VAL
1	A	520	PRO
1	A	637	PRO
1	B	239	HIS
1	B	334	LEU
1	B	484	VAL
1	B	520	PRO
1	B	637	PRO
1	B	668	SER
1	C	116	GLU
1	C	130	SER
1	C	239	HIS
1	C	364	ILE
1	C	484	VAL
1	C	520	PRO
1	C	637	PRO
1	D	130	SER
1	D	239	HIS
1	D	364	ILE
1	D	484	VAL
1	D	520	PRO
1	D	637	PRO
1	D	668	SER
1	E	239	HIS
1	E	484	VAL
1	E	520	PRO
1	E	637	PRO
1	E	755	ARG
1	F	130	SER
1	F	239	HIS
1	F	364	ILE
1	F	484	VAL
1	F	520	PRO
1	F	637	PRO
1	F	668	SER
2	O	21	LYS
2	P	21	LYS

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Mol	Chain	Res	Type
2	P	22	ASP
2	Q	21	LYS
2	Q	22	ASP
2	R	21	LYS
2	S	21	LYS
2	T	21	LYS
2	T	22	ASP
1	A	75	THR
1	A	364	ILE
1	A	668	SER
1	B	75	THR
1	B	364	ILE
1	D	75	THR
1	E	75	THR
1	E	364	ILE
1	F	75	THR
1	F	615	ILE
2	O	125	ILE
2	P	125	ILE
2	Q	125	ILE
2	S	125	ILE
2	S	133	ASP
2	T	14	GLU
2	T	125	ILE
1	A	190	PRO
1	B	190	PRO
1	C	615	ILE
1	E	190	PRO
2	R	125	ILE
1	D	190	PRO
1	F	190	PRO
1	A	537	GLY
1	A	615	ILE
1	B	615	ILE
1	C	177	ILE
1	C	190	PRO
1	C	537	GLY
1	D	177	ILE
1	D	615	ILE
1	E	615	ILE
1	F	537	GLY
1	A	177	ILE

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Mol	Chain	Res	Type
1	A	295	VAL
1	B	295	VAL
1	B	537	GLY
1	C	295	VAL
1	D	295	VAL
1	D	537	GLY
1	E	177	ILE
1	F	177	ILE
1	B	177	ILE
1	E	295	VAL
1	E	537	GLY
1	F	295	VAL
2	O	61	GLY
2	P	61	GLY
2	Q	61	GLY
2	R	42	ASN
2	R	61	GLY
2	S	61	GLY
2	T	61	GLY

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	664/675 (98%)	569 (86%)	95 (14%)	3	13
1	B	664/675 (98%)	566 (85%)	98 (15%)	3	12
1	C	664/675 (98%)	568 (86%)	96 (14%)	3	13
1	D	664/675 (98%)	567 (85%)	97 (15%)	3	12
1	E	664/675 (98%)	569 (86%)	95 (14%)	3	13
1	F	664/675 (98%)	568 (86%)	96 (14%)	3	13
2	O	123/127 (97%)	105 (85%)	18 (15%)	3	12
2	P	123/127 (97%)	105 (85%)	18 (15%)	3	12
2	Q	123/127 (97%)	105 (85%)	18 (15%)	3	12

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	R	123/127 (97%)	105 (85%)	18 (15%)	3	12
2	S	123/127 (97%)	105 (85%)	18 (15%)	3	12
2	T	123/127 (97%)	105 (85%)	18 (15%)	3	12
All	All	4722/4812 (98%)	4037 (86%)	685 (14%)	3	13

All (685) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	67	VAL
1	A	78	LYS
1	A	85	LEU
1	A	89	ILE
1	A	91	LYS
1	A	112	VAL
1	A	114	HIS
1	A	115	LYS
1	A	117	LEU
1	A	122	GLU
1	A	141	PHE
1	A	144	GLU
1	A	147	ARG
1	A	148	GLU
1	A	149	THR
1	A	156	ILE
1	A	159	TYR
1	A	172	GLU
1	A	175	LYS
1	A	182	ILE
1	A	188	LEU
1	A	197	LYS
1	A	210	PHE
1	A	212	GLN
1	A	213	LYS
1	A	217	LYS
1	A	254	ARG
1	A	279	ILE
1	A	284	LYS
1	A	288	VAL
1	A	293	ILE
1	A	297	LYS
1	A	324	THR

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Mol	Chain	Res	Type
1	A	334	LEU
1	A	349	ASN
1	A	356	ASP
1	A	370	LEU
1	A	377	GLN
1	A	387	ASN
1	A	395	GLU
1	A	397	GLU
1	A	400	LYS
1	A	408	LEU
1	A	413	LEU
1	A	414	LYS
1	A	415	GLU
1	A	433	TYR
1	A	434	LEU
1	A	438	ASN
1	A	446	ILE
1	A	451	ASN
1	A	455	TYR
1	A	456	LYS
1	A	469	PHE
1	A	472	ARG
1	A	473	ASN
1	A	479	LYS
1	A	480	ASN
1	A	481	VAL
1	A	484	VAL
1	A	495	PHE
1	A	500	SER
1	A	501	LEU
1	A	515	LYS
1	A	533	LEU
1	A	541	LYS
1	A	542	PRO
1	A	551	ASN
1	A	556	MET
1	A	625	LEU
1	A	635	ILE
1	A	639	ASN
1	A	646	THR
1	A	659	THR
1	A	665	LYS

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Mol	Chain	Res	Type
1	A	668	SER
1	A	670	ILE
1	A	672	ARG
1	A	675	ASN
1	A	678	VAL
1	A	688	PHE
1	A	707	SER
1	A	709	ASN
1	A	714	GLN
1	A	715	GLU
1	A	716	LYS
1	A	724	ARG
1	A	741	ILE
1	A	744	GLU
1	A	755	ARG
1	A	759	GLN
1	A	766	HIS
1	A	767	GLN
1	A	770	ASN
1	A	794	GLN
1	B	64	ASN
1	B	65	ASN
1	B	67	VAL
1	B	78	LYS
1	B	85	LEU
1	B	89	ILE
1	B	91	LYS
1	B	112	VAL
1	B	114	HIS
1	B	115	LYS
1	B	117	LEU
1	B	122	GLU
1	B	141	PHE
1	B	144	GLU
1	B	147	ARG
1	B	148	GLU
1	B	149	THR
1	B	156	ILE
1	B	159	TYR
1	B	172	GLU
1	B	175	LYS
1	B	182	ILE

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Mol	Chain	Res	Type
1	B	188	LEU
1	B	197	LYS
1	B	210	PHE
1	B	212	GLN
1	B	213	LYS
1	B	217	LYS
1	B	254	ARG
1	B	279	ILE
1	B	284	LYS
1	B	288	VAL
1	B	293	ILE
1	B	297	LYS
1	B	324	THR
1	B	334	LEU
1	B	349	ASN
1	B	356	ASP
1	B	370	LEU
1	B	377	GLN
1	B	387	ASN
1	B	395	GLU
1	B	397	GLU
1	B	400	LYS
1	B	408	LEU
1	B	413	LEU
1	B	414	LYS
1	B	415	GLU
1	B	433	TYR
1	B	434	LEU
1	B	438	ASN
1	B	446	ILE
1	B	451	ASN
1	B	455	TYR
1	B	456	LYS
1	B	469	PHE
1	B	472	ARG
1	B	473	ASN
1	B	479	LYS
1	B	480	ASN
1	B	481	VAL
1	B	484	VAL
1	B	495	PHE
1	B	500	SER

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Mol	Chain	Res	Type
1	B	501	LEU
1	B	515	LYS
1	B	533	LEU
1	B	541	LYS
1	B	542	PRO
1	B	551	ASN
1	B	556	MET
1	B	625	LEU
1	B	635	ILE
1	B	639	ASN
1	B	646	THR
1	B	650	THR
1	B	659	THR
1	B	665	LYS
1	B	668	SER
1	B	670	ILE
1	B	672	ARG
1	B	675	ASN
1	B	678	VAL
1	B	688	PHE
1	B	707	SER
1	B	709	ASN
1	B	714	GLN
1	B	715	GLU
1	B	716	LYS
1	B	724	ARG
1	B	741	ILE
1	B	744	GLU
1	B	755	ARG
1	B	759	GLN
1	B	766	HIS
1	B	767	GLN
1	B	770	ASN
1	B	794	GLN
1	C	65	ASN
1	C	67	VAL
1	C	78	LYS
1	C	85	LEU
1	C	89	ILE
1	C	91	LYS
1	C	112	VAL
1	C	114	HIS

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Mol	Chain	Res	Type
1	C	115	LYS
1	C	117	LEU
1	C	122	GLU
1	C	141	PHE
1	C	144	GLU
1	C	147	ARG
1	C	148	GLU
1	C	149	THR
1	C	156	ILE
1	C	159	TYR
1	C	172	GLU
1	C	175	LYS
1	C	182	ILE
1	C	188	LEU
1	C	197	LYS
1	C	210	PHE
1	C	212	GLN
1	C	213	LYS
1	C	217	LYS
1	C	254	ARG
1	C	279	ILE
1	C	284	LYS
1	C	288	VAL
1	C	293	ILE
1	C	297	LYS
1	C	324	THR
1	C	334	LEU
1	C	349	ASN
1	C	356	ASP
1	C	359	PRO
1	C	370	LEU
1	C	377	GLN
1	C	387	ASN
1	C	395	GLU
1	C	397	GLU
1	C	400	LYS
1	C	408	LEU
1	C	413	LEU
1	C	414	LYS
1	C	415	GLU
1	C	433	TYR
1	C	434	LEU

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Mol	Chain	Res	Type
1	C	438	ASN
1	C	446	ILE
1	C	451	ASN
1	C	455	TYR
1	C	456	LYS
1	C	469	PHE
1	C	472	ARG
1	C	473	ASN
1	C	479	LYS
1	C	480	ASN
1	C	481	VAL
1	C	484	VAL
1	C	495	PHE
1	C	501	LEU
1	C	515	LYS
1	C	533	LEU
1	C	541	LYS
1	C	542	PRO
1	C	551	ASN
1	C	556	MET
1	C	625	LEU
1	C	635	ILE
1	C	639	ASN
1	C	646	THR
1	C	659	THR
1	C	665	LYS
1	C	668	SER
1	C	670	ILE
1	C	672	ARG
1	C	675	ASN
1	C	678	VAL
1	C	688	PHE
1	C	707	SER
1	C	709	ASN
1	C	714	GLN
1	C	715	GLU
1	C	716	LYS
1	C	724	ARG
1	C	741	ILE
1	C	744	GLU
1	C	755	ARG
1	C	759	GLN

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Mol	Chain	Res	Type
1	C	766	HIS
1	C	767	GLN
1	C	770	ASN
1	C	794	GLN
1	D	64	ASN
1	D	67	VAL
1	D	78	LYS
1	D	85	LEU
1	D	89	ILE
1	D	91	LYS
1	D	112	VAL
1	D	114	HIS
1	D	115	LYS
1	D	117	LEU
1	D	122	GLU
1	D	141	PHE
1	D	144	GLU
1	D	147	ARG
1	D	148	GLU
1	D	149	THR
1	D	156	ILE
1	D	159	TYR
1	D	172	GLU
1	D	175	LYS
1	D	182	ILE
1	D	188	LEU
1	D	197	LYS
1	D	210	PHE
1	D	212	GLN
1	D	213	LYS
1	D	217	LYS
1	D	254	ARG
1	D	279	ILE
1	D	284	LYS
1	D	288	VAL
1	D	293	ILE
1	D	297	LYS
1	D	324	THR
1	D	334	LEU
1	D	349	ASN
1	D	356	ASP
1	D	370	LEU

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Mol	Chain	Res	Type
1	D	377	GLN
1	D	387	ASN
1	D	395	GLU
1	D	397	GLU
1	D	400	LYS
1	D	408	LEU
1	D	413	LEU
1	D	414	LYS
1	D	415	GLU
1	D	433	TYR
1	D	434	LEU
1	D	438	ASN
1	D	446	ILE
1	D	451	ASN
1	D	455	TYR
1	D	456	LYS
1	D	469	PHE
1	D	472	ARG
1	D	473	ASN
1	D	479	LYS
1	D	480	ASN
1	D	481	VAL
1	D	484	VAL
1	D	495	PHE
1	D	500	SER
1	D	501	LEU
1	D	515	LYS
1	D	533	LEU
1	D	541	LYS
1	D	542	PRO
1	D	551	ASN
1	D	556	MET
1	D	588	GLU
1	D	625	LEU
1	D	635	ILE
1	D	639	ASN
1	D	646	THR
1	D	659	THR
1	D	665	LYS
1	D	668	SER
1	D	670	ILE
1	D	672	ARG

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Mol	Chain	Res	Type
1	D	675	ASN
1	D	678	VAL
1	D	688	PHE
1	D	707	SER
1	D	709	ASN
1	D	714	GLN
1	D	715	GLU
1	D	716	LYS
1	D	724	ARG
1	D	741	ILE
1	D	744	GLU
1	D	755	ARG
1	D	759	GLN
1	D	766	HIS
1	D	767	GLN
1	D	770	ASN
1	D	794	GLN
1	E	67	VAL
1	E	78	LYS
1	E	85	LEU
1	E	89	ILE
1	E	91	LYS
1	E	112	VAL
1	E	114	HIS
1	E	115	LYS
1	E	117	LEU
1	E	122	GLU
1	E	141	PHE
1	E	144	GLU
1	E	147	ARG
1	E	148	GLU
1	E	149	THR
1	E	156	ILE
1	E	159	TYR
1	E	172	GLU
1	E	175	LYS
1	E	182	ILE
1	E	188	LEU
1	E	197	LYS
1	E	210	PHE
1	E	212	GLN
1	E	213	LYS

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Mol	Chain	Res	Type
1	E	217	LYS
1	E	254	ARG
1	E	279	ILE
1	E	284	LYS
1	E	288	VAL
1	E	293	ILE
1	E	297	LYS
1	E	324	THR
1	E	334	LEU
1	E	349	ASN
1	E	356	ASP
1	E	370	LEU
1	E	377	GLN
1	E	387	ASN
1	E	395	GLU
1	E	397	GLU
1	E	400	LYS
1	E	408	LEU
1	E	413	LEU
1	E	414	LYS
1	E	415	GLU
1	E	433	TYR
1	E	434	LEU
1	E	438	ASN
1	E	446	ILE
1	E	451	ASN
1	E	455	TYR
1	E	456	LYS
1	E	469	PHE
1	E	472	ARG
1	E	473	ASN
1	E	479	LYS
1	E	480	ASN
1	E	481	VAL
1	E	484	VAL
1	E	495	PHE
1	E	500	SER
1	E	501	LEU
1	E	515	LYS
1	E	533	LEU
1	E	541	LYS
1	E	542	PRO

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Mol	Chain	Res	Type
1	E	551	ASN
1	E	556	MET
1	E	625	LEU
1	E	635	ILE
1	E	639	ASN
1	E	646	THR
1	E	659	THR
1	E	665	LYS
1	E	668	SER
1	E	670	ILE
1	E	672	ARG
1	E	675	ASN
1	E	678	VAL
1	E	688	PHE
1	E	707	SER
1	E	709	ASN
1	E	714	GLN
1	E	715	GLU
1	E	716	LYS
1	E	724	ARG
1	E	741	ILE
1	E	744	GLU
1	E	755	ARG
1	E	759	GLN
1	E	766	HIS
1	E	767	GLN
1	E	770	ASN
1	E	794	GLN
1	F	67	VAL
1	F	78	LYS
1	F	85	LEU
1	F	89	ILE
1	F	91	LYS
1	F	112	VAL
1	F	114	HIS
1	F	115	LYS
1	F	117	LEU
1	F	122	GLU
1	F	141	PHE
1	F	144	GLU
1	F	147	ARG
1	F	148	GLU

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Mol	Chain	Res	Type
1	F	149	THR
1	F	156	ILE
1	F	159	TYR
1	F	172	GLU
1	F	175	LYS
1	F	182	ILE
1	F	188	LEU
1	F	197	LYS
1	F	210	PHE
1	F	212	GLN
1	F	213	LYS
1	F	217	LYS
1	F	254	ARG
1	F	279	ILE
1	F	284	LYS
1	F	288	VAL
1	F	293	ILE
1	F	297	LYS
1	F	324	THR
1	F	334	LEU
1	F	349	ASN
1	F	356	ASP
1	F	370	LEU
1	F	377	GLN
1	F	387	ASN
1	F	395	GLU
1	F	397	GLU
1	F	400	LYS
1	F	408	LEU
1	F	413	LEU
1	F	414	LYS
1	F	415	GLU
1	F	433	TYR
1	F	434	LEU
1	F	438	ASN
1	F	446	ILE
1	F	451	ASN
1	F	455	TYR
1	F	456	LYS
1	F	469	PHE
1	F	472	ARG
1	F	473	ASN

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Mol	Chain	Res	Type
1	F	479	LYS
1	F	480	ASN
1	F	481	VAL
1	F	484	VAL
1	F	495	PHE
1	F	500	SER
1	F	501	LEU
1	F	515	LYS
1	F	533	LEU
1	F	541	LYS
1	F	542	PRO
1	F	551	ASN
1	F	556	MET
1	F	625	LEU
1	F	635	ILE
1	F	639	ASN
1	F	646	THR
1	F	650	THR
1	F	659	THR
1	F	665	LYS
1	F	668	SER
1	F	670	ILE
1	F	672	ARG
1	F	675	ASN
1	F	678	VAL
1	F	688	PHE
1	F	707	SER
1	F	709	ASN
1	F	714	GLN
1	F	715	GLU
1	F	716	LYS
1	F	724	ARG
1	F	741	ILE
1	F	744	GLU
1	F	755	ARG
1	F	759	GLN
1	F	766	HIS
1	F	767	GLN
1	F	770	ASN
1	F	794	GLN
2	O	14	GLU
2	O	18	LEU

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Mol	Chain	Res	Type
2	O	49	GLN
2	O	54	GLU
2	O	55	VAL
2	O	62	THR
2	O	63	ILE
2	O	65	PHE
2	O	67	GLU
2	O	72	MET
2	O	83	GLU
2	O	84	GLU
2	O	94	LYS
2	O	97	ASN
2	O	117	THR
2	O	118	ASP
2	O	130	ILE
2	O	140	GLU
2	P	14	GLU
2	P	18	LEU
2	P	49	GLN
2	P	54	GLU
2	P	55	VAL
2	P	62	THR
2	P	63	ILE
2	P	65	PHE
2	P	67	GLU
2	P	72	MET
2	P	83	GLU
2	P	84	GLU
2	P	94	LYS
2	P	97	ASN
2	P	117	THR
2	P	118	ASP
2	P	130	ILE
2	P	140	GLU
2	Q	14	GLU
2	Q	18	LEU
2	Q	49	GLN
2	Q	54	GLU
2	Q	55	VAL
2	Q	62	THR
2	Q	63	ILE
2	Q	65	PHE

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Mol	Chain	Res	Type
2	Q	67	GLU
2	Q	72	MET
2	Q	83	GLU
2	Q	84	GLU
2	Q	94	LYS
2	Q	97	ASN
2	Q	117	THR
2	Q	118	ASP
2	Q	130	ILE
2	Q	140	GLU
2	R	14	GLU
2	R	18	LEU
2	R	49	GLN
2	R	54	GLU
2	R	55	VAL
2	R	62	THR
2	R	63	ILE
2	R	65	PHE
2	R	67	GLU
2	R	72	MET
2	R	83	GLU
2	R	84	GLU
2	R	94	LYS
2	R	97	ASN
2	R	117	THR
2	R	118	ASP
2	R	130	ILE
2	R	140	GLU
2	S	14	GLU
2	S	18	LEU
2	S	49	GLN
2	S	54	GLU
2	S	55	VAL
2	S	62	THR
2	S	63	ILE
2	S	65	PHE
2	S	67	GLU
2	S	72	MET
2	S	83	GLU
2	S	84	GLU
2	S	94	LYS
2	S	97	ASN

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Mol	Chain	Res	Type
2	S	117	THR
2	S	118	ASP
2	S	130	ILE
2	S	140	GLU
2	T	14	GLU
2	T	18	LEU
2	T	49	GLN
2	T	54	GLU
2	T	55	VAL
2	T	62	THR
2	T	63	ILE
2	T	65	PHE
2	T	67	GLU
2	T	72	MET
2	T	83	GLU
2	T	84	GLU
2	T	94	LYS
2	T	97	ASN
2	T	117	THR
2	T	118	ASP
2	T	130	ILE
2	T	140	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (258) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	81	GLN
1	A	83	GLN
1	A	126	ASN
1	A	129	ASN
1	A	155	ASN
1	A	165	GLN
1	A	212	GLN
1	A	239	HIS
1	A	323	ASN
1	A	349	ASN
1	A	351	HIS
1	A	368	GLN
1	A	387	ASN
1	A	438	ASN
1	A	480	ASN
1	A	507	GLN

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Mol	Chain	Res	Type
1	A	510	GLN
1	A	518	ASN
1	A	521	ASN
1	A	551	ASN
1	A	553	GLN
1	A	555	GLN
1	A	577	HIS
1	A	581	GLN
1	A	583	ASN
1	A	618	ASN
1	A	639	ASN
1	A	655	ASN
1	A	666	ASN
1	A	709	ASN
1	A	730	ASN
1	A	747	ASN
1	A	759	GLN
1	A	767	GLN
1	A	770	ASN
1	A	781	ASN
1	A	794	GLN
1	B	81	GLN
1	B	83	GLN
1	B	126	ASN
1	B	129	ASN
1	B	155	ASN
1	B	165	GLN
1	B	212	GLN
1	B	239	HIS
1	B	323	ASN
1	B	349	ASN
1	B	351	HIS
1	B	368	GLN
1	B	387	ASN
1	B	438	ASN
1	B	480	ASN
1	B	507	GLN
1	B	518	ASN
1	B	521	ASN
1	B	551	ASN
1	B	553	GLN
1	B	555	GLN

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Mol	Chain	Res	Type
1	B	577	HIS
1	B	581	GLN
1	B	583	ASN
1	B	618	ASN
1	B	639	ASN
1	B	655	ASN
1	B	666	ASN
1	B	709	ASN
1	B	730	ASN
1	B	747	ASN
1	B	759	GLN
1	B	767	GLN
1	B	770	ASN
1	B	781	ASN
1	B	794	GLN
1	C	81	GLN
1	C	83	GLN
1	C	126	ASN
1	C	129	ASN
1	C	165	GLN
1	C	212	GLN
1	C	239	HIS
1	C	323	ASN
1	C	349	ASN
1	C	351	HIS
1	C	368	GLN
1	C	387	ASN
1	C	438	ASN
1	C	480	ASN
1	C	507	GLN
1	C	510	GLN
1	C	518	ASN
1	C	521	ASN
1	C	551	ASN
1	C	553	GLN
1	C	555	GLN
1	C	577	HIS
1	C	581	GLN
1	C	583	ASN
1	C	618	ASN
1	C	639	ASN
1	C	655	ASN

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Mol	Chain	Res	Type
1	C	666	ASN
1	C	709	ASN
1	C	730	ASN
1	C	747	ASN
1	C	767	GLN
1	C	770	ASN
1	C	781	ASN
1	C	794	GLN
1	D	81	GLN
1	D	83	GLN
1	D	126	ASN
1	D	129	ASN
1	D	155	ASN
1	D	165	GLN
1	D	212	GLN
1	D	239	HIS
1	D	323	ASN
1	D	349	ASN
1	D	351	HIS
1	D	368	GLN
1	D	387	ASN
1	D	438	ASN
1	D	480	ASN
1	D	507	GLN
1	D	510	GLN
1	D	518	ASN
1	D	521	ASN
1	D	551	ASN
1	D	553	GLN
1	D	555	GLN
1	D	577	HIS
1	D	581	GLN
1	D	583	ASN
1	D	618	ASN
1	D	639	ASN
1	D	655	ASN
1	D	666	ASN
1	D	709	ASN
1	D	730	ASN
1	D	747	ASN
1	D	759	GLN
1	D	767	GLN

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Mol	Chain	Res	Type
1	D	770	ASN
1	D	781	ASN
1	D	794	GLN
1	E	81	GLN
1	E	83	GLN
1	E	126	ASN
1	E	129	ASN
1	E	165	GLN
1	E	212	GLN
1	E	239	HIS
1	E	323	ASN
1	E	349	ASN
1	E	351	HIS
1	E	368	GLN
1	E	387	ASN
1	E	438	ASN
1	E	480	ASN
1	E	507	GLN
1	E	510	GLN
1	E	518	ASN
1	E	521	ASN
1	E	551	ASN
1	E	553	GLN
1	E	555	GLN
1	E	577	HIS
1	E	581	GLN
1	E	583	ASN
1	E	618	ASN
1	E	639	ASN
1	E	655	ASN
1	E	666	ASN
1	E	709	ASN
1	E	730	ASN
1	E	747	ASN
1	E	759	GLN
1	E	767	GLN
1	E	770	ASN
1	E	781	ASN
1	E	794	GLN
1	F	81	GLN
1	F	83	GLN
1	F	126	ASN

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Mol	Chain	Res	Type
1	F	129	ASN
1	F	165	GLN
1	F	212	GLN
1	F	239	HIS
1	F	323	ASN
1	F	349	ASN
1	F	351	HIS
1	F	368	GLN
1	F	387	ASN
1	F	438	ASN
1	F	480	ASN
1	F	507	GLN
1	F	510	GLN
1	F	518	ASN
1	F	521	ASN
1	F	551	ASN
1	F	553	GLN
1	F	555	GLN
1	F	577	HIS
1	F	581	GLN
1	F	583	ASN
1	F	618	ASN
1	F	639	ASN
1	F	655	ASN
1	F	666	ASN
1	F	709	ASN
1	F	730	ASN
1	F	747	ASN
1	F	759	GLN
1	F	767	GLN
1	F	770	ASN
1	F	781	ASN
1	F	794	GLN
2	O	41	GLN
2	O	49	GLN
2	O	107	HIS
2	O	111	ASN
2	O	123	GLN
2	O	135	GLN
2	O	143	GLN
2	P	41	GLN
2	P	49	GLN

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Mol	Chain	Res	Type
2	P	107	HIS
2	P	123	GLN
2	P	135	GLN
2	P	143	GLN
2	Q	41	GLN
2	Q	49	GLN
2	Q	107	HIS
2	Q	111	ASN
2	Q	123	GLN
2	Q	135	GLN
2	Q	143	GLN
2	R	41	GLN
2	R	49	GLN
2	R	107	HIS
2	R	111	ASN
2	R	123	GLN
2	R	135	GLN
2	R	143	GLN
2	S	41	GLN
2	S	49	GLN
2	S	107	HIS
2	S	111	ASN
2	S	123	GLN
2	S	135	GLN
2	S	143	GLN
2	T	41	GLN
2	T	49	GLN
2	T	107	HIS
2	T	111	ASN
2	T	123	GLN
2	T	135	GLN
2	T	143	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 24 ligands modelled in this entry, 24 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ > 2			OWAB(Å ²)	Q < 0.9
1	A	735/747 (98%)	-0.22	22 (2%)	52	38	24, 80, 134, 147	0
1	B	735/747 (98%)	-0.21	15 (2%)	65	48	25, 80, 134, 148	0
1	C	735/747 (98%)	-0.22	18 (2%)	59	44	25, 80, 134, 147	0
1	D	735/747 (98%)	-0.22	15 (2%)	65	48	26, 80, 134, 147	0
1	E	735/747 (98%)	-0.23	14 (1%)	66	49	26, 80, 135, 147	0
1	F	735/747 (98%)	-0.22	16 (2%)	62	46	25, 81, 134, 147	0
2	O	146/149 (97%)	-0.31	0	100	100	24, 65, 127, 137	0
2	P	146/149 (97%)	-0.33	0	100	100	24, 65, 126, 137	0
2	Q	146/149 (97%)	-0.33	1 (0%)	84	73	23, 65, 126, 137	0
2	R	146/149 (97%)	-0.30	0	100	100	24, 65, 126, 137	0
2	S	146/149 (97%)	-0.34	0	100	100	24, 66, 126, 137	0
2	T	146/149 (97%)	-0.31	1 (0%)	84	73	23, 65, 126, 137	0
All	All	5286/5376 (98%)	-0.24	102 (1%)	66	49	23, 77, 134, 148	0

All (102) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	171	TYR	6.0
1	F	171	TYR	5.5
1	D	171	TYR	5.3
1	A	171	TYR	5.0
1	B	171	TYR	4.9
1	E	171	TYR	4.6
1	B	237	PHE	4.3
1	D	237	PHE	4.2
1	C	237	PHE	4.1
1	E	237	PHE	4.0
1	A	202	ASP	4.0

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Mol	Chain	Res	Type	RSRZ
1	B	230	ILE	3.9
1	F	237	PHE	3.7
1	C	202	ASP	3.6
1	E	202	ASP	3.6
1	B	205	SER	3.4
1	F	230	ILE	3.4
1	A	203	SER	3.3
1	E	205	SER	3.3
1	C	206	SER	3.3
1	E	162	ASN	3.2
1	F	212	GLN	3.1
1	D	205	SER	3.1
1	D	206	SER	3.1
1	F	185	ASP	3.1
1	F	229	PHE	3.0
1	B	206	SER	3.0
1	D	162	ASN	3.0
1	A	229	PHE	3.0
1	F	162	ASN	3.0
1	C	230	ILE	2.9
1	B	193	LEU	2.9
1	F	72	THR	2.8
1	E	209	LEU	2.8
1	A	162	ASN	2.8
1	E	204	ASP	2.8
1	A	237	PHE	2.8
1	A	204	ASP	2.7
1	C	185	ASP	2.7
1	C	162	ASN	2.7
1	C	193	LEU	2.7
1	A	72	THR	2.7
1	F	213	LYS	2.7
1	A	205	SER	2.7
1	E	206	SER	2.7
1	E	230	ILE	2.6
1	B	229	PHE	2.6
1	D	204	ASP	2.6
1	F	202	ASP	2.6
1	A	193	LEU	2.6
1	C	205	SER	2.6
1	A	213	LYS	2.6
1	C	213	LYS	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	185	ASP	2.5
1	D	193	LEU	2.5
1	C	218	LEU	2.5
1	E	193	LEU	2.5
1	A	212	GLN	2.5
1	D	185	ASP	2.5
1	F	193	LEU	2.4
1	A	186	LYS	2.4
1	E	229	PHE	2.4
1	B	162	ASN	2.4
1	A	230	ILE	2.4
1	E	225	ILE	2.4
1	A	185	ASP	2.4
1	D	202	ASP	2.4
1	D	229	PHE	2.4
1	C	72	THR	2.4
1	D	72	THR	2.4
1	F	205	SER	2.4
1	D	230	ILE	2.3
1	A	209	LEU	2.3
1	A	206	SER	2.3
1	F	118	GLN	2.3
1	A	161	ILE	2.3
1	F	206	SER	2.3
1	B	225	ILE	2.2
1	B	72	THR	2.2
2	T	81	SER	2.2
1	E	118	GLN	2.2
1	B	186	LYS	2.2
1	F	186	LYS	2.2
1	F	204	ASP	2.1
1	C	212	GLN	2.1
1	A	124	GLU	2.1
1	A	118	GLN	2.1
1	B	791	GLU	2.1
1	B	209	LEU	2.1
1	C	220	LEU	2.1
2	Q	52	ILE	2.1
1	E	185	ASP	2.1
1	D	186	LYS	2.0
1	A	657	ILE	2.0
1	C	161	ILE	2.0

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Mol	Chain	Res	Type	RSRZ
1	C	188	LEU	2.0
1	A	566	TYR	2.0
1	C	184	LYS	2.0
1	D	209	LEU	2.0
1	B	204	ASP	2.0
1	D	118	GLN	2.0
1	C	225	ILE	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	MG	B	901	1/1	0.96	0.05	13,13,13,13	0
3	MG	A	900	1/1	0.97	0.09	18,18,18,18	0
3	MG	D	903	1/1	0.97	0.08	19,19,19,19	0
3	MG	E	904	1/1	0.97	0.06	14,14,14,14	0
4	CA	R	707	1/1	0.97	0.04	79,79,79,79	0
4	CA	S	709	1/1	0.97	0.03	78,78,78,78	0
4	CA	S	810	1/1	0.97	0.04	45,45,45,45	0
4	CA	T	711	1/1	0.97	0.04	77,77,77,77	0
4	CA	P	703	1/1	0.98	0.03	71,71,71,71	0
4	CA	P	804	1/1	0.98	0.05	43,43,43,43	0
4	CA	Q	705	1/1	0.98	0.02	79,79,79,79	0
4	CA	Q	806	1/1	0.98	0.04	46,46,46,46	0
3	MG	C	902	1/1	0.98	0.07	14,14,14,14	0
4	CA	R	808	1/1	0.98	0.04	41,41,41,41	0
3	MG	F	905	1/1	0.98	0.08	16,16,16,16	0
4	CA	O	701	1/1	0.98	0.03	83,83,83,83	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	CA	O	802	1/1	0.98	0.04	44,44,44,44	0
4	CA	T	812	1/1	0.98	0.04	43,43,43,43	0
4	CA	O	801	1/1	0.99	0.05	29,29,29,29	0
4	CA	S	809	1/1	0.99	0.04	25,25,25,25	0
4	CA	P	803	1/1	0.99	0.03	33,33,33,33	0
4	CA	R	807	1/1	0.99	0.04	33,33,33,33	0
4	CA	T	811	1/1	0.99	0.04	31,31,31,31	0
4	CA	Q	805	1/1	0.99	0.04	32,32,32,32	0

6.5 Other polymers [i](#)

There are no such residues in this entry.