



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 10:50 AM UTC

PDB ID : 3QT1 / pdb_00003qt1
Title : RNA polymerase II variant containing A Chimeric RPB9-C11 subunit
Authors : Ruan, W.; Lehmann, E.; Thomm, M.; Kostrewa, D.; Cramer, P.
Deposited on : 2011-02-22
Resolution : 4.30 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:

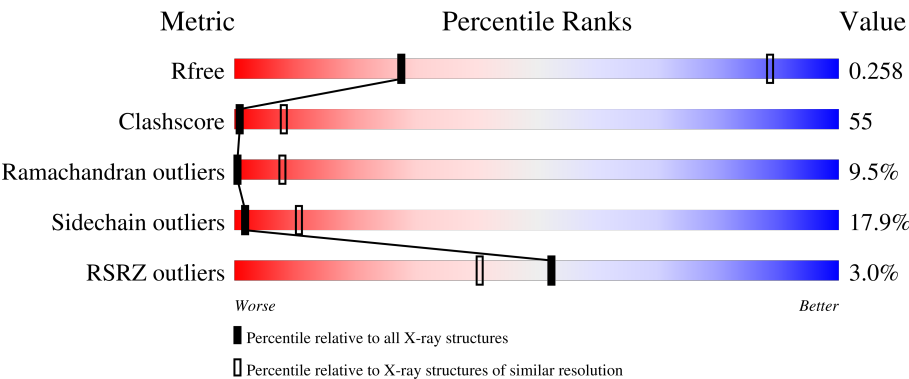
MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
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 4.30 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	1052 (4.70-3.90)
Clashscore	190562	1097 (4.70-3.90)
Ramachandran outliers	187476	1001 (4.70-3.90)
Sidechain outliers	187428	1007 (4.72-3.88)
RSRZ outliers	180081	1049 (4.70-3.90)

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	1733	2% 20% 44% 16% • 18%
2	B	1224	3% 24% 49% 16% • 10%
3	C	318	2% 25% 44% 14% • 16%
4	D	219	3% 20% 45% 14% • 19%
5	E	215	2% 24% 62% 13% •

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Mol	Chain	Length	Quality of chain
6	F	155	%12%32%12%.44%
7	G	171	%29%56%15%. .
8	H	146	7%32%44%13%.8%
9	I	133	2%14%17%. .65%
10	J	70	3%20%60%11%.7%
11	K	120	2%31%44%18%.5%
12	L	70	4%11%27%20%7%34%

2 Entry composition [i](#)

There are 14 unique types of molecules in this entry. The entry contains 30535 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 DNA-directed RNA polymerase II subunit RPB1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	1416	Total	C	N	O	S	0	0	0
			11143	7021	1949	2111	62			

- Molecule 2 is a protein called DNA-directed RNA polymerase II subunit RPB2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	1103	Total	C	N	O	S	0	0	0
			8770	5554	1535	1626	55			

- Molecule 3 is a protein called DNA-directed RNA polymerase II subunit RPB3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	266	Total	C	N	O	S	0	0	0
			2095	1317	348	417	13			

- Molecule 4 is a protein called DNA-directed RNA polymerase II subunit RPB4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	178	Total	C	N	O	S	0	0	0
			1434	887	257	288	2			

- Molecule 5 is a protein called DNA-directed RNA polymerases I, II, and III subunit RPABC1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	E	214	Total	C	N	O	S	0	0	0
			1752	1111	309	321	11			

- Molecule 6 is a protein called DNA-directed RNA polymerases I, II, and III subunit RPABC2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	F	87	Total	C	N	O	S	0	0	0
			705	451	119	132	3			

- Molecule 7 is a protein called DNA-directed RNA polymerase II subunit RPB7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	G	171	Total	C	N	O	S	0	0	0
			1340	861	222	249	8			

- Molecule 8 is a protein called DNA-directed RNA polymerases I, II, and III subunit RPABC3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
8	H	134	Total	C	N	O	S	0	0	0
			1076	677	182	213	4			

- Molecule 9 is a protein called DNA-directed RNA polymerase II subunit RPB9, DNA-directed RNA polymerase III subunit RPC10.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
9	I	47	Total	C	N	O	S	0	0	0
			398	246	72	75	5			

- Molecule 10 is a protein called DNA-directed RNA polymerases I, II, and III subunit RPABC5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
10	J	65	Total	C	N	O	S	0	0	0
			532	339	93	94	6			

- Molecule 11 is a protein called DNA-directed RNA polymerase II subunit RPB11.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
11	K	114	Total	C	N	O	S	0	0	0
			919	590	156	171	2			

- Molecule 12 is a protein called DNA-directed RNA polymerases I, II, and III subunit RPABC4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
12	L	46	Total	C	N	O	S	0	0	0
			363	224	72	63	4			

- Molecule 13 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
13	A	2	Total	Zn	0	0
			2	2		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
13	B	1	Total 1	Zn 1	0	0
13	C	1	Total 1	Zn 1	0	0
13	I	1	Total 1	Zn 1	0	0
13	J	1	Total 1	Zn 1	0	0
13	L	1	Total 1	Zn 1	0	0

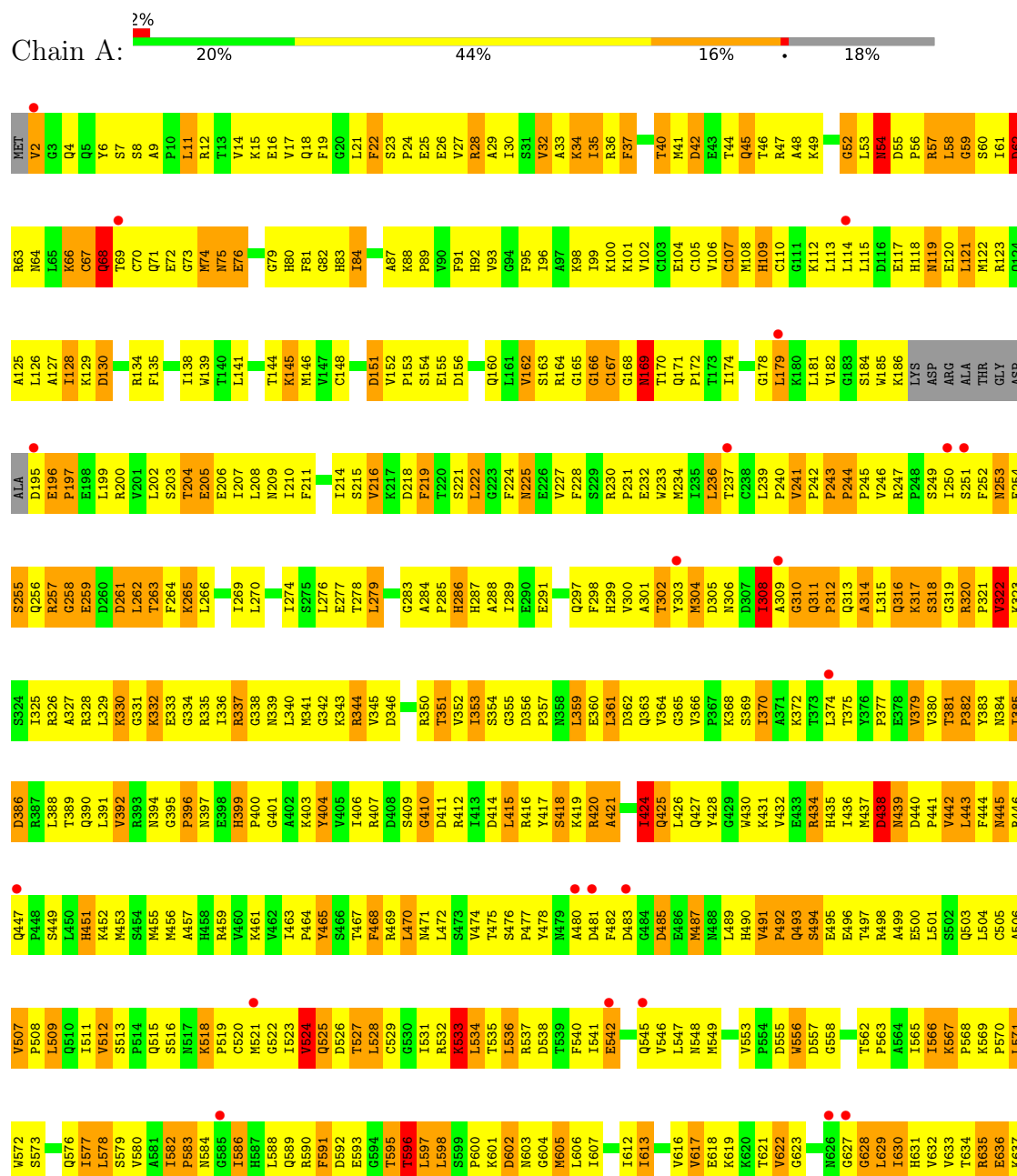
- Molecule 14 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
14	A	1	Total 1	Mg 1	0	0

3 Residue-property plots

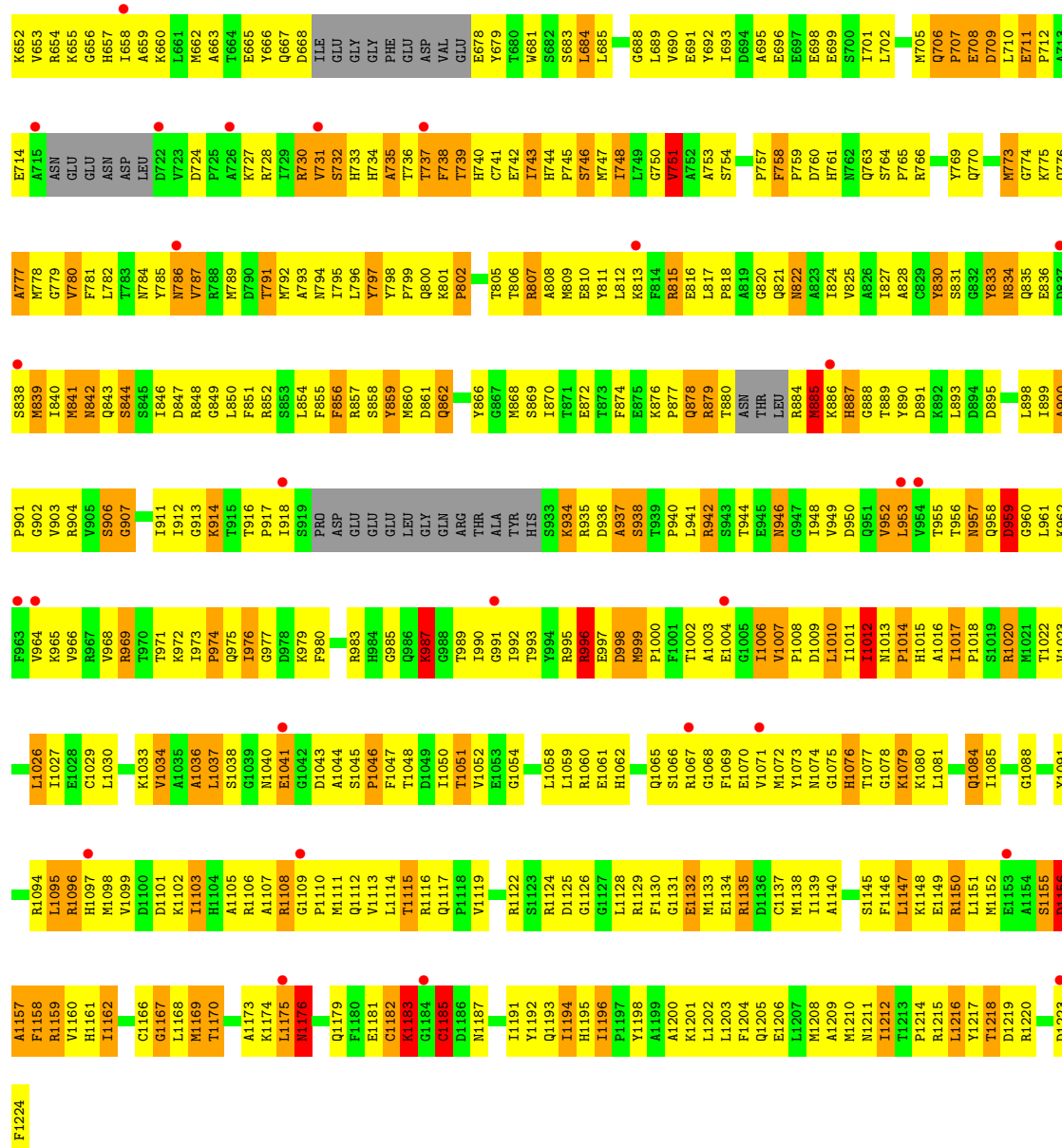
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density.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

- Molecule 1: DNA-directed RNA polymerase II subunit RPB1

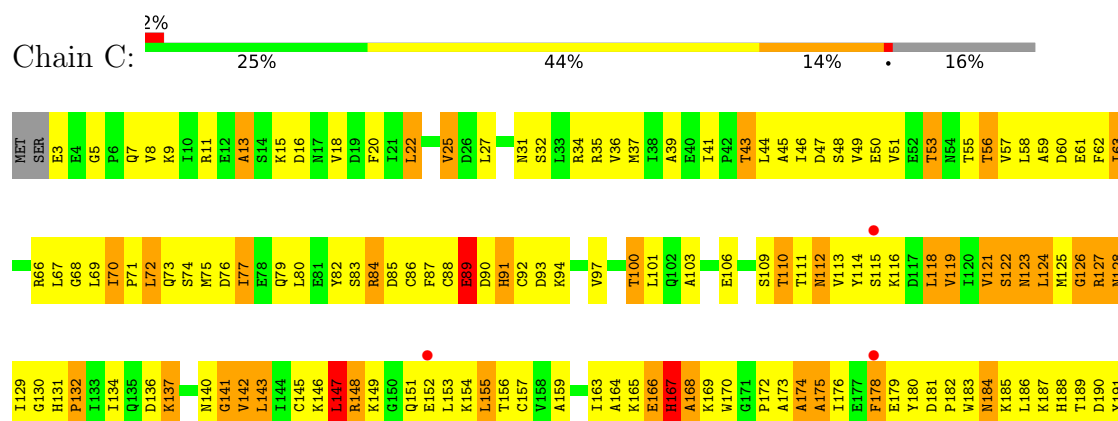


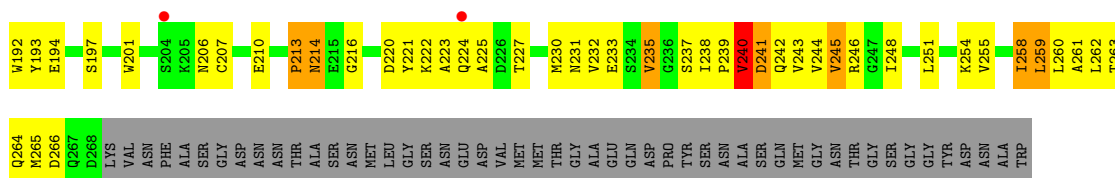




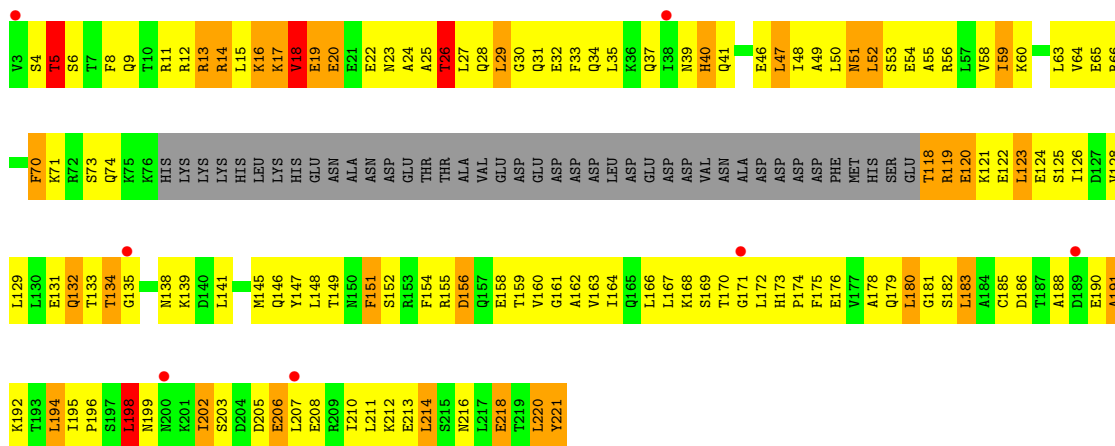
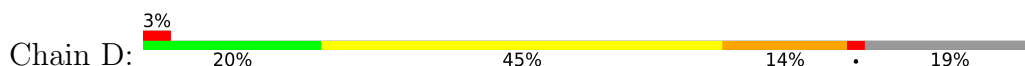


• Molecule 3: DNA-directed RNA polymerase II subunit RPB3

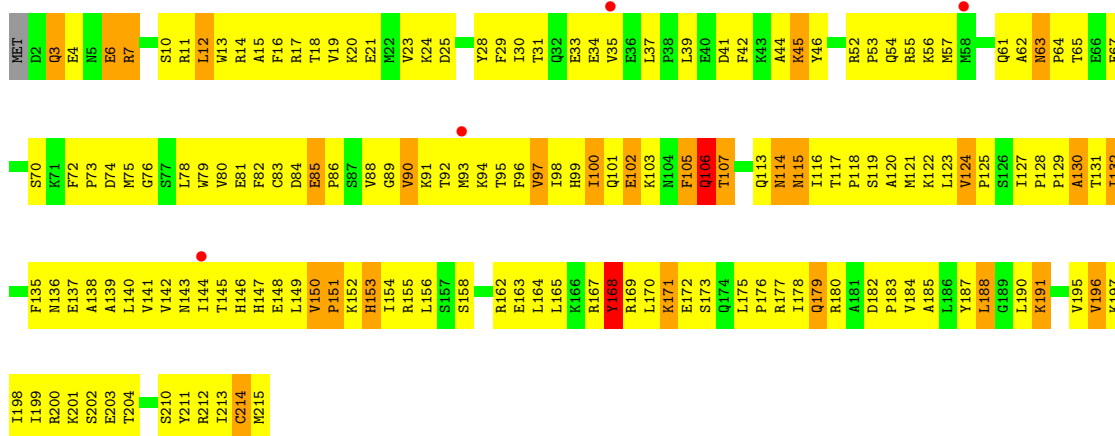




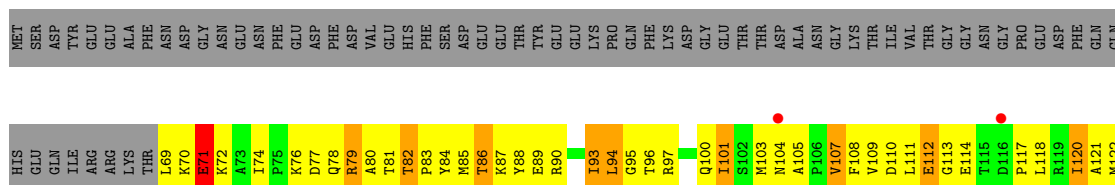
• Molecule 4: DNA-directed RNA polymerase II subunit RPB4

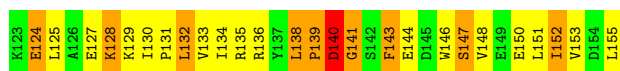


• Molecule 5: DNA-directed RNA polymerases I, II, and III subunit RPABC1

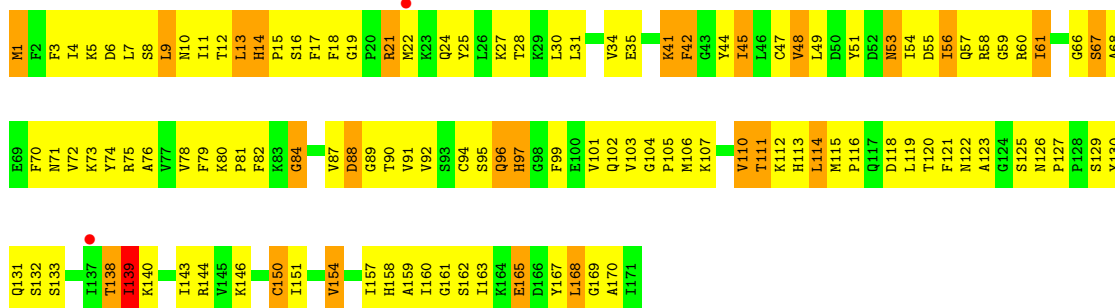


• Molecule 6: DNA-directed RNA polymerases I, II, and III subunit RPABC2

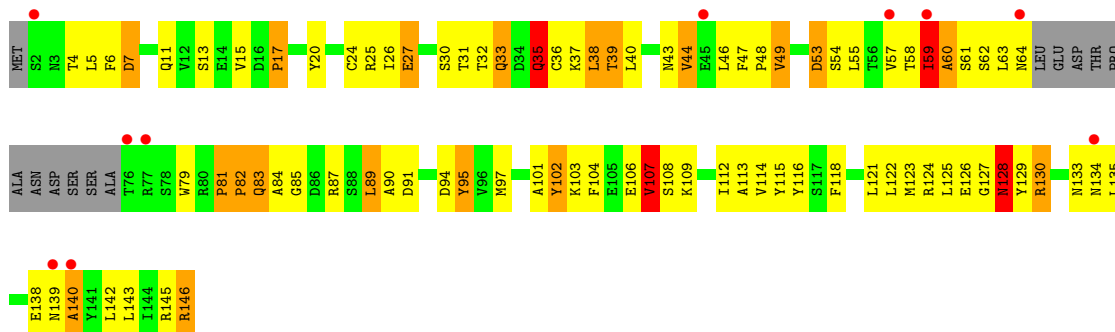




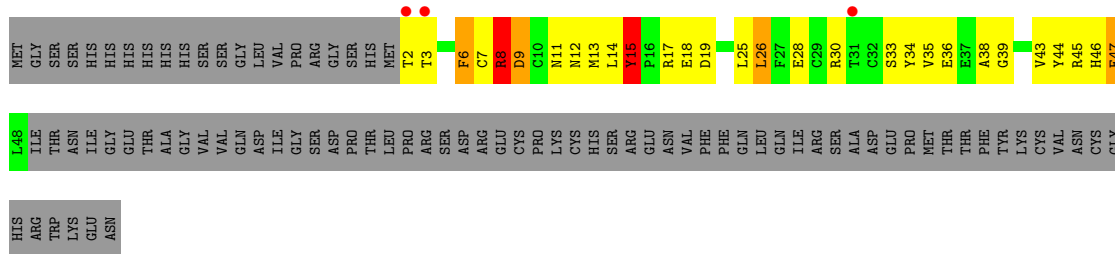
• Molecule 7: DNA-directed RNA polymerase II subunit RPB7



• Molecule 8: DNA-directed RNA polymerases I, II, and III subunit RPABC3

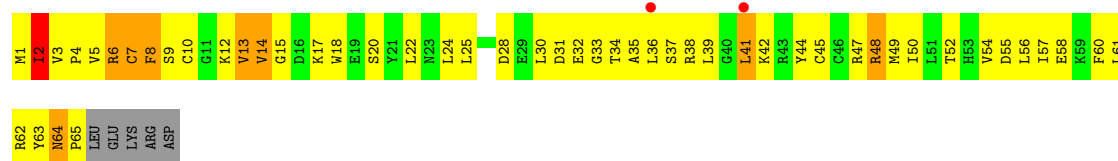


• Molecule 9: DNA-directed RNA polymerase II subunit RPB9, DNA-directed RNA polymerase III subunit RPC10

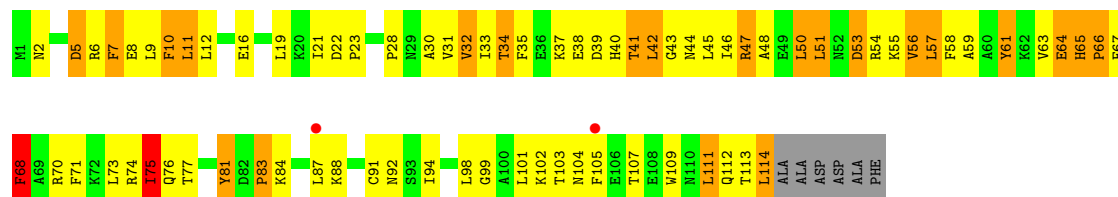


• Molecule 10: DNA-directed RNA polymerases I, II, and III subunit RPABC5

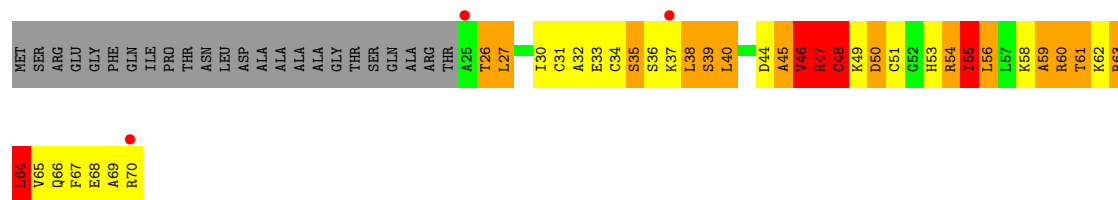




- Molecule 11: DNA-directed RNA polymerase II subunit RPB11



- Molecule 12: DNA-directed RNA polymerases I, II, and III subunit RPABC4



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	222.38Å 393.38Å 281.43Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.60 – 4.30 48.60 – 4.30	Depositor EDS
% Data completeness (in resolution range)	98.5 (48.60-4.30) 98.4 (48.60-4.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.67 (at 4.29Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6.1_357)	Depositor
R, R_{free}	0.235 , 0.281 0.211 , 0.258	Depositor DCC
R_{free} test set	2022 reflections (2.41%)	wwPDB-VP
Wilson B-factor (Å ²)	172.3	Xtriage
Anisotropy	0.553	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 202.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.41$, $\langle L^2 \rangle = 0.24$	Xtriage
Estimated twinning fraction	0.057 for 1/2*h-1/2*k,-3/2*h-1/2*k,-l 0.067 for 1/2*h+1/2*k,3/2*h-1/2*k,-l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	30535	wwPDB-VP
Average B, all atoms (Å ²)	144.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.83% 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: ZN, 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	0.44	0/11342	0.95	45/15337 (0.3%)
2	B	0.42	0/8939	0.89	21/12051 (0.2%)
3	C	0.42	0/2133	0.92	6/2891 (0.2%)
4	D	0.41	0/1444	0.82	1/1935 (0.1%)
5	E	0.39	0/1788	0.92	5/2406 (0.2%)
6	F	0.47	0/717	1.04	6/967 (0.6%)
7	G	0.39	0/1368	0.91	2/1844 (0.1%)
8	H	0.38	0/1094	0.81	1/1481 (0.1%)
9	I	0.47	0/406	0.94	3/546 (0.5%)
10	J	0.41	0/541	0.81	1/727 (0.1%)
11	K	0.43	0/937	0.90	2/1265 (0.2%)
12	L	0.47	0/365	0.87	0/485
All	All	0.42	0/31074	0.91	93/41935 (0.2%)

There are no bond length outliers.

All (93) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	241	VAL	CA-C-N	9.62	130.28	120.38
1	A	241	VAL	C-N-CA	9.62	130.28	120.38
1	A	553	VAL	CA-C-N	9.01	129.59	119.32
1	A	553	VAL	C-N-CA	9.01	129.59	119.32
4	D	26	THR	N-CA-C	-8.62	103.28	112.93
5	E	63	ASN	CA-C-N	7.62	129.37	119.84
5	E	63	ASN	C-N-CA	7.62	129.37	119.84
1	A	518	LYS	N-CA-C	7.52	114.76	108.07
6	F	82	THR	CA-C-N	7.17	126.89	119.64
6	F	82	THR	C-N-CA	7.17	126.89	119.64
5	E	63	ASN	N-CA-C	7.14	116.53	108.25
2	B	1088	GLY	CA-C-N	6.87	128.43	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1088	GLY	C-N-CA	6.87	128.43	119.84
2	B	527	THR	CA-C-N	6.86	126.83	119.76
2	B	527	THR	C-N-CA	6.86	126.83	119.76
1	A	582	ILE	N-CA-C	6.77	114.59	107.76
1	A	524	VAL	N-CA-C	6.76	116.77	108.53
1	A	381	THR	CA-C-N	6.64	128.14	119.84
1	A	381	THR	C-N-CA	6.64	128.14	119.84
2	B	833	TYR	N-CA-C	-6.60	105.76	113.88
11	K	75	ILE	N-CA-C	6.55	117.55	107.78
1	A	395	GLY	CA-C-N	6.50	127.96	119.84
1	A	395	GLY	C-N-CA	6.50	127.96	119.84
1	A	793	SER	CA-C-N	6.48	125.98	119.24
1	A	793	SER	C-N-CA	6.48	125.98	119.24
1	A	279	LEU	N-CA-C	-6.47	106.60	114.75
2	B	427	ASP	N-CA-C	-6.45	105.95	113.88
3	C	84	ARG	N-CA-C	-6.35	105.69	113.50
3	C	194	GLU	N-CA-C	-6.35	106.07	113.88
1	A	553	VAL	N-CA-C	6.28	113.76	107.55
1	A	582	ILE	CA-C-N	6.27	127.68	119.84
1	A	582	ILE	C-N-CA	6.27	127.68	119.84
1	A	954	TRP	N-CA-C	6.25	118.36	108.23
1	A	1299	VAL	N-CA-C	6.22	116.12	108.53
1	A	37	PHE	CA-C-N	6.18	126.72	120.04
1	A	37	PHE	C-N-CA	6.18	126.72	120.04
1	A	470	LEU	N-CA-C	6.14	117.56	108.60
9	I	15	TYR	CA-C-N	6.08	126.10	119.90
9	I	15	TYR	C-N-CA	6.08	126.10	119.90
6	F	71	GLU	N-CA-C	-6.07	105.46	112.92
1	A	320	ARG	CA-C-N	6.01	125.69	119.56
1	A	320	ARG	C-N-CA	6.01	125.69	119.56
2	B	570	VAL	CA-C-N	5.99	127.32	119.84
2	B	570	VAL	C-N-CA	5.99	127.32	119.84
2	B	934	LYS	N-CA-C	5.97	118.50	107.99
6	F	124	GLU	N-CA-C	-5.96	106.34	113.97
1	A	244	PRO	CA-C-N	-5.92	114.79	120.83
1	A	244	PRO	C-N-CA	-5.92	114.79	120.83
5	E	200	ARG	N-CA-C	5.91	117.19	108.86
2	B	706	GLN	CA-C-N	5.85	127.15	119.84
2	B	706	GLN	C-N-CA	5.85	127.15	119.84
9	I	39	GLY	N-CA-C	-5.79	106.63	115.66
1	A	765	VAL	N-CA-C	5.78	116.00	110.74
1	A	230	ARG	CA-C-N	5.71	126.98	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	230	ARG	C-N-CA	5.71	126.98	119.84
3	C	148	ARG	CA-C-O	5.71	121.25	117.94
2	B	900	ALA	CA-C-N	5.67	126.93	119.84
2	B	900	ALA	C-N-CA	5.67	126.93	119.84
1	A	1311	VAL	N-CA-C	5.63	116.11	107.77
11	K	112	GLN	N-CA-C	5.62	117.62	109.07
10	J	50	ILE	CB-CA-C	-5.56	104.87	111.81
2	B	1185	CYS	N-CA-C	-5.48	106.37	113.16
5	E	6	GLU	N-CA-C	-5.42	105.47	112.68
6	F	143	PHE	N-CA-C	5.41	117.48	108.99
1	A	1204	ASP	N-CA-C	-5.39	106.70	113.28
3	C	91	HIS	N-CA-C	5.33	115.28	108.24
2	B	273	LEU	CA-C-N	5.33	125.27	119.78
2	B	273	LEU	C-N-CA	5.33	125.27	119.78
2	B	127	GLY	N-CA-C	5.32	120.00	111.64
1	A	174	ILE	N-CA-C	5.32	115.96	108.36
1	A	243	PRO	CA-C-N	5.25	125.79	120.38
1	A	243	PRO	C-N-CA	5.25	125.79	120.38
7	G	14	HIS	CA-C-N	5.21	126.35	119.84
7	G	14	HIS	C-N-CA	5.21	126.35	119.84
1	A	784	LEU	CA-C-N	5.20	126.34	119.84
1	A	784	LEU	C-N-CA	5.20	126.34	119.84
1	A	769	SER	N-CA-C	5.18	116.95	108.76
1	A	852	TYR	N-CA-C	5.17	119.22	113.01
1	A	638	GLY	CA-C-N	5.17	126.30	119.84
1	A	638	GLY	C-N-CA	5.17	126.30	119.84
6	F	132	LEU	N-CA-C	5.16	117.31	108.90
2	B	254	LEU	N-CA-C	5.12	116.67	107.80
1	A	797	LYS	N-CA-C	-5.11	106.55	112.89
2	B	885	MET	N-CA-C	5.11	114.89	108.45
8	H	35	GLN	N-CA-C	-5.10	107.11	113.38
1	A	1207	LEU	N-CA-C	5.09	115.02	108.34
1	A	244	PRO	N-CA-C	5.09	116.91	110.70
3	C	147	LEU	N-CA-C	5.08	116.79	109.07
1	A	1328	TYR	N-CA-C	5.07	116.57	107.80
2	B	95	ILE	N-CA-C	5.06	116.00	108.46
1	A	487	MET	N-CA-C	5.04	117.33	109.52
2	B	1012	ILE	N-CA-C	5.02	115.81	108.53
3	C	89	GLU	N-CA-C	-5.01	103.22	110.64

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	11143	0	11217	1336	0
2	B	8770	0	8795	1039	0
3	C	2095	0	2051	233	0
4	D	1434	0	1460	172	0
5	E	1752	0	1776	187	0
6	F	705	0	731	108	0
7	G	1340	0	1357	162	0
8	H	1076	0	1046	102	0
9	I	398	0	370	38	0
10	J	532	0	542	93	0
11	K	919	0	929	117	0
12	L	363	0	386	68	0
13	A	2	0	0	0	0
13	B	1	0	0	0	0
13	C	1	0	0	0	0
13	I	1	0	0	0	0
13	J	1	0	0	0	0
13	L	1	0	0	0	0
14	A	1	0	0	0	0
All	All	30535	0	30660	3390	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 55.

All (3390) 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:775:ILE:HB	1:A:797:LYS:O	1.37	1.23
1:A:567:LYS:HB2	1:A:568:PRO:HD2	1.21	1.17
2:B:763:GLN:HG2	2:B:765:PRO:HD2	1.26	1.15
9:I:25:LEU:HB3	9:I:38:ALA:HB2	1.17	1.14
2:B:577:ALA:HB1	2:B:589:VAL:HG11	1.30	1.13
4:D:134:THR:HG22	4:D:135:GLY:H	1.06	1.10
10:J:64:ASN:HB3	10:J:65:PRO:HD3	1.30	1.10
1:A:482:PHE:HB2	2:B:836:GLU:O	1.52	1.09

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:E:94:LYS:HE2	5:E:98:ILE:HD11	1.35	1.07
2:B:824:ILE:HG12	10:J:48:ARG:HH12	1.14	1.03
2:B:1116:ARG:HG3	2:B:1198:TYR:CD2	1.93	1.03
2:B:295:GLY:H	2:B:298:LEU:HD23	1.19	1.03
2:B:510:LYS:HG3	2:B:511:PRO:HD3	1.37	1.02
1:A:855:THR:HG21	1:A:857:ARG:HE	1.23	1.01
11:K:65:HIS:CD2	11:K:67:PHE:H	1.80	0.99
7:G:45:ILE:HA	7:G:78:VAL:HG12	1.44	0.99
4:D:66:ARG:HH22	7:G:48:VAL:HG23	1.26	0.98
2:B:210:LYS:HD3	2:B:482:VAL:HG13	1.41	0.98
4:D:188:ALA:HA	4:D:191:ALA:HB3	1.44	0.98
1:A:1116:LEU:HB3	1:A:1308:THR:HG21	1.44	0.98
2:B:1130:PHE:HE2	2:B:1151:LEU:HD21	1.27	0.97
2:B:706:GLN:HB2	2:B:709:ASP:HB2	1.47	0.97
2:B:273:LEU:HD21	2:B:360:PHE:HD1	1.28	0.96
1:A:845:LEU:HD12	1:A:1069:ALA:HB2	1.47	0.96
1:A:741:ASN:HD21	1:A:743:VAL:HG23	1.26	0.96
3:C:103:ALA:HB3	3:C:153:LEU:HB3	1.47	0.96
1:A:814:PHE:HD1	2:B:519:TRP:HE3	1.03	0.95
1:A:95:PHE:HD1	1:A:234:MET:HG2	1.30	0.95
7:G:13:LEU:HD21	7:G:17:PHE:HD2	1.32	0.95
2:B:430:ARG:HG2	2:B:430:ARG:HH11	1.29	0.95
1:A:903:ASN:HD22	1:A:905:ASP:H	1.01	0.94
2:B:29:ASP:HB3	2:B:658:ILE:HD13	1.45	0.94
1:A:463:ILE:HB	1:A:464:PRO:HD2	1.49	0.94
1:A:361:LEU:HA	1:A:471:ASN:HD22	1.32	0.94
5:E:124:VAL:HG13	5:E:132:ILE:HB	1.47	0.93
2:B:846:ILE:HD13	2:B:974:PRO:HB2	1.51	0.93
2:B:996:ARG:HH22	3:C:175:ALA:H	1.07	0.93
1:A:121:LEU:HD22	1:A:141:LEU:HD21	1.48	0.93
2:B:848:ARG:HH22	2:B:996:ARG:HD2	1.33	0.93
1:A:1156:PRO:HA	1:A:1190:PRO:HB3	1.51	0.93
1:A:1036:ARG:HG2	1:A:1036:ARG:HH11	1.32	0.92
1:A:600:PRO:HG2	1:A:601:LYS:HG3	1.51	0.92
2:B:515:HIS:HD2	2:B:517:THR:H	1.08	0.92
1:A:326:ARG:HG3	1:A:1406:VAL:HG21	1.51	0.91
2:B:1162:ILE:HD11	2:B:1194:ILE:HG21	1.51	0.91
2:B:351:TYR:CE1	2:B:355:ILE:HD11	2.05	0.90
2:B:1034:VAL:HG22	2:B:1059:LEU:HD13	1.51	0.90
1:A:239:LEU:HD12	1:A:240:PRO:HD2	1.53	0.90
1:A:95:PHE:CD1	1:A:234:MET:HG2	2.06	0.90

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:53:LEU:HD23	1:A:54:ASN:H	1.35	0.90
1:A:591:PHE:HD2	1:A:595:THR:HB	1.34	0.90
7:G:139:ILE:HG23	7:G:140:LYS:HG3	1.53	0.90
1:A:216:VAL:O	1:A:219:PHE:HB2	1.71	0.90
1:A:265:LYS:HA	1:A:265:LYS:HE3	1.54	0.90
1:A:671:ALA:H	1:A:676:MET:HE3	1.35	0.90
8:H:130:ARG:HH11	8:H:130:ARG:H	1.16	0.90
1:A:899:VAL:HB	1:A:929:LEU:HD12	1.53	0.90
1:A:1116:LEU:HB3	1:A:1308:THR:CG2	2.01	0.90
2:B:1181:GLU:HA	2:B:1187:ASN:O	1.72	0.89
1:A:254:GLU:HB2	2:B:935:ARG:HH22	1.36	0.89
1:A:265:LYS:HD2	1:A:322:VAL:HG21	1.55	0.89
11:K:65:HIS:HD2	11:K:67:PHE:N	1.70	0.89
1:A:567:LYS:CB	1:A:568:PRO:HD2	2.01	0.89
6:F:90:ARG:HD2	6:F:155:LEU:HD13	1.55	0.89
3:C:63:ILE:HA	3:C:66:ARG:HG3	1.51	0.89
3:C:57:VAL:HG11	10:J:60:PHE:HB3	1.55	0.89
2:B:515:HIS:CD2	2:B:517:THR:H	1.90	0.88
2:B:975:GLN:HG3	2:B:976:ILE:H	1.38	0.88
1:A:1006:ILE:HD11	5:E:163:GLU:HG3	1.56	0.88
2:B:824:ILE:HG12	10:J:48:ARG:NH1	1.88	0.88
11:K:65:HIS:HD2	11:K:67:PHE:H	0.90	0.88
2:B:230:ALA:HB3	2:B:231:PRO:HD3	1.54	0.87
4:D:71:LYS:HG3	4:D:74:GLN:HG3	1.54	0.87
1:A:524:VAL:HG12	1:A:525:GLN:H	1.40	0.87
1:A:666:ILE:HG23	2:B:1026:LEU:HB2	1.57	0.87
1:A:425:GLN:H	1:A:425:GLN:NE2	1.72	0.87
1:A:506:ALA:HB1	1:A:508:PRO:HD2	1.57	0.87
6:F:101:ILE:HD13	6:F:120:ILE:HG21	1.55	0.87
1:A:14:VAL:H	1:A:1432:GLN:HE22	1.22	0.87
2:B:408:LEU:HD11	2:B:545:ILE:HG13	1.55	0.87
4:D:56:ARG:HA	4:D:59:ILE:HD12	1.55	0.87
1:A:814:PHE:CD1	2:B:519:TRP:HE3	1.92	0.86
1:A:506:ALA:HB3	1:A:509:LEU:HG	1.56	0.86
4:D:179:GLN:HE22	7:G:1:MET:HB2	1.38	0.86
11:K:56:VAL:HA	11:K:77:THR:HG22	1.56	0.86
4:D:134:THR:HG22	4:D:135:GLY:N	1.87	0.86
1:A:933:TYR:O	1:A:933:TYR:HD2	1.59	0.85
3:C:115:SER:HB3	3:C:141:GLY:O	1.76	0.85
1:A:1318:THR:HB	5:E:141:VAL:HG11	1.57	0.85
1:A:814:PHE:HD1	2:B:519:TRP:CE3	1.93	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:521:LEU:HD22	2:B:633:VAL:HG12	1.58	0.85
3:C:88:CYS:SG	3:C:91:HIS:HA	2.16	0.85
12:L:39:SER:O	12:L:40:LEU:HG	1.77	0.85
2:B:1130:PHE:CE2	2:B:1151:LEU:HD21	2.12	0.84
1:A:492:PRO:HG3	1:A:501:LEU:HD12	1.59	0.84
1:A:567:LYS:HB2	1:A:568:PRO:CD	2.06	0.84
2:B:810:GLU:HA	2:B:815:ARG:HH22	1.42	0.84
10:J:36:LEU:HB2	10:J:47:ARG:NH1	1.92	0.84
2:B:277:LYS:HE2	2:B:336:ARG:C	2.02	0.84
1:A:19:PHE:HZ	1:A:1397:LEU:HD21	1.41	0.84
1:A:35:ILE:H	1:A:35:ILE:HD12	1.42	0.84
1:A:256:GLN:O	1:A:257:ARG:HB2	1.76	0.84
1:A:1035:TYR:O	1:A:1036:ARG:HB2	1.77	0.84
2:B:65:GLU:OE1	2:B:418:LYS:HE3	1.76	0.84
1:A:110:CYS:HB2	1:A:167:CYS:SG	2.17	0.84
1:A:1094:VAL:HG13	1:A:1113:THR:HB	1.57	0.84
2:B:613:VAL:HG13	2:B:627:PHE:O	1.77	0.84
4:D:134:THR:CG2	4:D:135:GLY:H	1.89	0.84
10:J:64:ASN:HB3	10:J:65:PRO:CD	2.08	0.84
1:A:40:THR:HB	1:A:41:MET:HG2	1.60	0.84
8:H:4:THR:HA	8:H:60:ALA:HB2	1.58	0.83
9:I:35:VAL:HG12	9:I:36:GLU:H	1.43	0.83
1:A:19:PHE:CZ	1:A:1397:LEU:HD21	2.13	0.83
1:A:770:VAL:HG12	1:A:771:GLU:HG3	1.61	0.83
7:G:13:LEU:HD21	7:G:17:PHE:CD2	2.14	0.83
2:B:405:ARG:HB3	2:B:631:GLY:HA3	1.61	0.83
1:A:868:TYR:HE1	1:A:1064:VAL:HG13	1.43	0.83
2:B:408:LEU:O	2:B:412:LEU:HG	1.79	0.83
2:B:241:ARG:HA	2:B:253:THR:HG22	1.58	0.83
1:A:664:THR:HA	1:A:742:ASN:HD22	1.43	0.82
6:F:100:GLN:O	6:F:105:ALA:HB2	1.80	0.82
11:K:21:ILE:HG12	11:K:33:ILE:HG23	1.58	0.82
1:A:591:PHE:CD2	1:A:595:THR:HB	2.14	0.82
1:A:335:ARG:HD3	1:A:339:ASN:HD22	1.43	0.82
1:A:903:ASN:ND2	1:A:905:ASP:H	1.77	0.82
2:B:957:ASN:HD21	2:B:961:LEU:HB2	1.44	0.82
9:I:6:PHE:HA	9:I:14:LEU:HG	1.59	0.82
1:A:954:TRP:CZ3	5:E:203:GLU:HB2	2.15	0.82
1:A:483:ASP:HB2	2:B:987:LYS:HG3	1.62	0.82
7:G:158:HIS:HD2	7:G:159:ALA:H	1.25	0.82
2:B:172:ILE:HD13	2:B:178:ASN:HD22	1.45	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:597:MET:HA	2:B:597:MET:HE3	1.62	0.82
2:B:1159:ARG:HB3	2:B:1159:ARG:HH11	1.45	0.82
1:A:414:ASP:HB3	1:A:417:TYR:HB2	1.60	0.82
10:J:8:PHE:H	10:J:8:PHE:HD2	1.27	0.82
1:A:107:CYS:HA	1:A:171:GLN:HE22	1.45	0.82
2:B:430:ARG:HH11	2:B:430:ARG:CG	1.93	0.82
2:B:570:VAL:HB	2:B:573:GLN:HB3	1.61	0.82
3:C:166:GLU:O	11:K:6:ARG:HG3	1.80	0.82
12:L:26:THR:HG22	12:L:27:LEU:H	1.44	0.82
2:B:202:TYR:H	2:B:202:TYR:HD2	1.27	0.81
1:A:775:ILE:CB	1:A:797:LYS:O	2.26	0.81
1:A:254:GLU:HB2	2:B:935:ARG:NH2	1.95	0.81
2:B:114:PRO:HG3	2:B:181:LEU:HD11	1.62	0.81
2:B:211:VAL:O	2:B:480:SER:HA	1.78	0.81
2:B:406:LEU:HD12	2:B:545:ILE:HD11	1.62	0.81
1:A:342:GLY:HA3	2:B:1131:GLY:HA2	1.59	0.81
1:A:1197:LEU:HD11	1:A:1238:ILE:HD11	1.62	0.81
2:B:345:LYS:HG2	2:B:346:GLU:H	1.44	0.81
2:B:999:MET:HG3	2:B:1000:PRO:HD2	1.63	0.81
12:L:53:HIS:HB3	12:L:55:ILE:CD1	2.10	0.80
1:A:845:LEU:CD1	1:A:1069:ALA:HB2	2.11	0.80
3:C:39:ALA:HA	3:C:164:ALA:HB3	1.64	0.80
1:A:590:ARG:O	1:A:591:PHE:HB2	1.81	0.80
2:B:1076:HIS:HD2	11:K:40:HIS:CE1	1.99	0.80
3:C:148:ARG:H	3:C:151:GLN:HG3	1.46	0.80
10:J:6:ARG:HA	10:J:12:LYS:O	1.81	0.80
1:A:1354:ASN:O	1:A:1358:SER:HB3	1.81	0.80
2:B:899:ILE:HG21	2:B:949:VAL:HG21	1.62	0.80
1:A:332:LYS:HG2	1:A:333:GLU:HG2	1.64	0.80
2:B:801:LYS:O	10:J:52:THR:HG23	1.82	0.80
3:C:118:LEU:HD12	3:C:132:PRO:HG3	1.63	0.80
1:A:362:ASP:HB3	1:A:507:VAL:HG12	1.61	0.80
1:A:868:TYR:CE1	1:A:1064:VAL:HG13	2.17	0.80
2:B:1157:ALA:O	2:B:1158:PHE:HB2	1.82	0.80
7:G:158:HIS:CD2	7:G:159:ALA:H	1.99	0.80
10:J:1:MET:O	10:J:2:ILE:HG22	1.81	0.80
1:A:353:ILE:HG21	1:A:487:MET:HG3	1.64	0.79
2:B:1158:PHE:HE2	2:B:1160:VAL:HG22	1.47	0.79
4:D:54:GLU:O	4:D:58:VAL:HG23	1.81	0.79
11:K:5:ASP:HB2	11:K:7:PHE:CZ	2.17	0.79
2:B:295:GLY:N	2:B:298:LEU:HD23	1.98	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:843:GLN:HB2	2:B:993:THR:HB	1.64	0.79
2:B:996:ARG:NH2	3:C:175:ALA:H	1.78	0.79
1:A:1094:VAL:HG22	1:A:1113:THR:HG21	1.65	0.79
8:H:102:TYR:H	8:H:102:TYR:HD2	1.28	0.79
3:C:259:LEU:HD21	11:K:92:ASN:OD1	1.83	0.79
4:D:53:SER:HB3	4:D:152:SER:HB3	1.65	0.79
2:B:840:ILE:HB	2:B:1011:ILE:HB	1.63	0.79
8:H:13:SER:HB3	8:H:27:GLU:O	1.83	0.79
7:G:111:THR:HG22	7:G:114:LEU:HD22	1.65	0.79
1:A:855:THR:CG2	1:A:857:ARG:HE	1.95	0.78
4:D:25:ALA:CB	4:D:196:PRO:HG2	2.13	0.78
2:B:94:LYS:HG2	2:B:95:ILE:H	1.48	0.78
1:A:391:LEU:HD23	1:A:400:PRO:O	1.84	0.78
3:C:242:GLN:HE21	3:C:246:ARG:HE	1.27	0.78
4:D:138:ASN:HB3	4:D:141:LEU:HB3	1.65	0.78
1:A:537:ARG:HD2	8:H:20:TYR:CE1	2.18	0.78
2:B:261:ARG:HH11	2:B:261:ARG:HB3	1.46	0.78
3:C:184:ASN:HD21	3:C:189:THR:H	1.31	0.78
11:K:113:THR:HG22	11:K:114:LEU:H	1.49	0.78
1:A:852:TYR:CE2	1:A:1060:PRO:HB2	2.19	0.78
2:B:280:ILE:HG23	2:B:284:ILE:HD12	1.66	0.78
7:G:31:LEU:HD22	7:G:48:VAL:HG21	1.64	0.78
1:A:1124:HIS:H	1:A:1124:HIS:CD2	2.02	0.78
2:B:615:MET:HE3	2:B:615:MET:H	1.49	0.78
2:B:955:THR:HG22	2:B:956:THR:N	1.98	0.78
1:A:1102:LYS:HG2	1:A:1106:ASN:HD21	1.48	0.78
6:F:87:LYS:HG3	6:F:88:TYR:CD1	2.19	0.78
2:B:702:LEU:HD23	2:B:738:PHE:H	1.49	0.77
1:A:1205:LYS:O	1:A:1207:LEU:HG	1.84	0.77
2:B:794:ASN:C	2:B:795:ILE:HD12	2.09	0.77
10:J:35:ALA:O	10:J:39:LEU:HD12	1.84	0.77
1:A:134:ARG:HG2	1:A:138:ILE:HD11	1.67	0.77
1:A:946:VAL:HG22	5:E:201:LYS:HD2	1.65	0.77
2:B:174:LEU:HD21	2:B:204:ILE:HD11	1.67	0.77
1:A:518:LYS:HB2	1:A:519:PRO:HD2	1.67	0.77
3:C:70:ILE:O	3:C:72:LEU:HD12	1.83	0.77
1:A:70:CYS:O	1:A:72:GLU:HG2	1.84	0.77
1:A:899:VAL:HG13	1:A:908:LEU:HD21	1.67	0.77
2:B:326:ASP:O	2:B:330:ALA:HB2	1.85	0.77
1:A:66:LYS:HD3	1:A:67:CYS:N	2.00	0.77
1:A:351:THR:HG23	1:A:468:PHE:CE1	2.20	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:34:LYS:HE3	1:A:57:ARG:HH12	1.47	0.77
1:A:108:MET:HB3	1:A:210:ILE:HD13	1.67	0.77
1:A:881:GLN:NE2	1:A:959:ASN:HA	1.99	0.77
2:B:36:ALA:HA	2:B:39:ARG:HD2	1.65	0.77
12:L:32:ALA:HB2	12:L:55:ILE:HG13	1.66	0.77
1:A:336:ILE:HD11	2:B:1203:LEU:CD1	2.15	0.77
2:B:619:ILE:HG22	2:B:620:ARG:HG3	1.66	0.77
3:C:41:ILE:O	3:C:163:ILE:HG22	1.84	0.76
5:E:202:SER:OG	5:E:204:THR:HG22	1.85	0.76
3:C:111:THR:HB	3:C:147:LEU:HG	1.67	0.76
5:E:124:VAL:HB	5:E:125:PRO:HD3	1.68	0.76
3:C:49:VAL:HG22	3:C:157:CYS:SG	2.25	0.76
4:D:56:ARG:HD3	4:D:149:THR:HA	1.65	0.76
7:G:115:MET:HB3	7:G:119:LEU:HD23	1.67	0.76
12:L:32:ALA:CB	12:L:55:ILE:HG13	2.16	0.76
2:B:653:VAL:HG22	2:B:689:LEU:HB3	1.66	0.76
1:A:672:ASP:HB3	1:A:736:ASN:HD21	1.51	0.76
2:B:613:VAL:HG22	2:B:628:THR:HA	1.67	0.76
1:A:1313:LEU:HD23	1:A:1338:VAL:HG21	1.68	0.76
2:B:216:GLU:HA	2:B:406:LEU:HD23	1.68	0.76
4:D:206:GLU:O	4:D:210:ILE:HG13	1.85	0.76
10:J:8:PHE:HD2	10:J:8:PHE:N	1.82	0.76
1:A:93:VAL:HG22	1:A:301:ALA:HA	1.68	0.76
2:B:559:SER:HA	2:B:563:MET:HB3	1.66	0.76
1:A:855:THR:HG21	1:A:857:ARG:NE	1.99	0.76
1:A:1242:VAL:HG12	1:A:1243:VAL:H	1.50	0.76
1:A:755:PHE:HA	1:A:758:ILE:HD12	1.66	0.76
4:D:15:LEU:O	4:D:17:LYS:HG3	1.85	0.76
1:A:600:PRO:HA	8:H:25:ARG:NH1	2.01	0.75
2:B:541:LEU:HB2	2:B:747:MET:HE3	1.68	0.75
7:G:34:VAL:HG12	7:G:45:ILE:HG21	1.67	0.75
1:A:706:HIS:CD2	1:A:1283:VAL:HG22	2.21	0.75
6:F:100:GLN:HB3	7:G:15:PRO:HB3	1.65	0.75
7:G:55:ASP:HB3	7:G:73:LYS:HB2	1.68	0.75
1:A:79:GLY:HA3	1:A:243:PRO:HB2	1.66	0.75
2:B:824:ILE:CG1	10:J:48:ARG:HH12	1.98	0.75
2:B:616:ILE:HD12	2:B:625:LYS:HB2	1.69	0.75
2:B:999:MET:HA	2:B:999:MET:HE3	1.68	0.75
10:J:48:ARG:HG2	10:J:48:ARG:HH11	1.52	0.75
2:B:351:TYR:O	2:B:355:ILE:HG13	1.85	0.75
2:B:387:LEU:HD23	2:B:393:LYS:HG3	1.69	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:53:LEU:HD23	1:A:54:ASN:HD22	1.49	0.75
2:B:796:LEU:HB3	2:B:799:PRO:HG3	1.69	0.75
4:D:25:ALA:HB1	4:D:196:PRO:HG2	1.69	0.75
1:A:606:LEU:HG	1:A:613:ILE:HG13	1.69	0.75
2:B:975:GLN:HG3	2:B:976:ILE:N	2.00	0.75
2:B:1202:LEU:O	2:B:1206:GLU:HG3	1.87	0.74
3:C:56:THR:HG21	3:C:145:CYS:SG	2.27	0.74
4:D:39:ASN:CG	4:D:40:HIS:H	1.94	0.74
1:A:356:ASP:HB3	1:A:359:LEU:HD21	1.68	0.74
1:A:56:PRO:O	1:A:57:ARG:HG3	1.87	0.74
4:D:56:ARG:NH2	4:D:155:ARG:HG2	2.02	0.74
1:A:442:VAL:HB	1:A:489:LEU:HD11	1.68	0.74
1:A:427:GLN:HB2	1:A:430:TRP:CD2	2.21	0.74
2:B:789:MET:HE2	2:B:965:LYS:HB3	1.68	0.74
8:H:101:ALA:HB2	8:H:116:TYR:CE2	2.22	0.74
1:A:852:TYR:CD2	1:A:1060:PRO:HB2	2.23	0.74
1:A:1210:GLY:HA2	1:A:1228:TRP:HE1	1.50	0.74
2:B:1158:PHE:CE2	2:B:1160:VAL:HG22	2.23	0.74
4:D:35:LEU:H	4:D:35:LEU:HD12	1.52	0.74
5:E:44:ALA:O	5:E:45:LYS:HB2	1.87	0.74
1:A:1100:ARG:HH21	1:A:1351:GLU:HG2	1.53	0.74
10:J:64:ASN:CB	10:J:65:PRO:HD3	2.15	0.74
2:B:800:GLN:HB3	10:J:52:THR:HG22	1.70	0.73
4:D:47:LEU:HB3	4:D:174:PRO:HG2	1.70	0.73
1:A:285:PRO:HB2	1:A:288:ALA:HB3	1.70	0.73
2:B:1099:VAL:O	2:B:1103:ILE:HG22	1.88	0.73
1:A:283:GLY:O	1:A:285:PRO:HD3	1.88	0.73
1:A:490:HIS:HB3	2:B:1150:ARG:NH1	2.02	0.73
1:A:1148:ILE:O	9:I:47:GLU:HA	1.89	0.73
2:B:406:LEU:O	2:B:408:LEU:HD12	1.88	0.73
4:D:208:GLU:HG3	4:D:212:LYS:HE3	1.69	0.73
2:B:842:ASN:HD21	2:B:844:SER:HB2	1.54	0.73
6:F:111:LEU:C	6:F:113:GLY:H	1.96	0.73
7:G:27:LYS:HD3	7:G:51:TYR:CE2	2.24	0.73
2:B:577:ALA:HB1	2:B:589:VAL:CG1	2.16	0.73
2:B:839:MET:HE1	2:B:980:PHE:HB2	1.70	0.73
2:B:542:MET:SD	2:B:747:MET:HE2	2.29	0.73
2:B:1182:CYS:SG	2:B:1185:CYS:HB2	2.28	0.73
12:L:49:LYS:O	12:L:50:ASP:HB2	1.87	0.73
1:A:507:VAL:HG22	1:A:521:MET:HE1	1.71	0.73
2:B:770:GLN:HB2	2:B:985:GLY:H	1.52	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:369:SER:HB3	11:K:2:ASN:HD21	1.52	0.73
5:E:153:HIS:C	5:E:154:ILE:HD12	2.14	0.73
1:A:115:LEU:HD13	1:A:122:MET:HB2	1.71	0.73
7:G:121:PHE:CZ	7:G:123:ALA:HB2	2.24	0.73
1:A:446:ARG:HB2	1:A:487:MET:HE3	1.71	0.72
6:F:96:THR:O	6:F:100:GLN:HG3	1.89	0.72
7:G:111:THR:HG22	7:G:114:LEU:HD13	1.70	0.72
1:A:635:ARG:NH1	1:A:635:ARG:HA	2.04	0.72
4:D:59:ILE:HG22	4:D:63:LEU:HD12	1.71	0.72
5:E:96:PHE:CE1	5:E:100:ILE:HD11	2.24	0.72
1:A:962:ARG:HA	1:A:965:GLN:HG3	1.71	0.72
2:B:850:LEU:HB2	10:J:8:PHE:HD1	1.55	0.72
7:G:129:SER:HB2	7:G:138:THR:HG23	1.71	0.72
1:A:549:MET:HE1	1:A:656:TRP:CD1	2.24	0.72
1:A:1121:GLU:HG2	1:A:1122:PRO:HD2	1.70	0.72
1:A:1308:THR:HG23	1:A:1310:GLY:H	1.55	0.72
2:B:261:ARG:HB3	2:B:261:ARG:NH1	2.05	0.72
1:A:399:HIS:HB3	1:A:400:PRO:HD3	1.71	0.72
1:A:1269:GLU:O	1:A:1270:ASN:HB2	1.90	0.72
2:B:345:LYS:O	2:B:347:LYS:HG2	1.89	0.72
1:A:1148:ILE:HB	1:A:1196:GLU:HG2	1.72	0.72
4:D:30:GLY:C	4:D:32:GLU:H	1.97	0.72
5:E:16:PHE:CE1	5:E:20:LYS:HE3	2.25	0.72
7:G:81:PRO:HG3	7:G:106:MET:SD	2.29	0.72
11:K:7:PHE:HA	11:K:10:PHE:CE2	2.25	0.72
1:A:726:ARG:HD2	1:A:765:VAL:O	1.89	0.72
1:A:853:ASP:OD1	1:A:855:THR:HG22	1.90	0.72
2:B:25:ILE:HD11	2:B:651:LEU:HD12	1.71	0.72
2:B:777:ALA:HA	2:B:1095:LEU:CB	2.19	0.72
3:C:183:TRP:O	3:C:185:LYS:HG3	1.89	0.72
1:A:706:HIS:HD2	1:A:1283:VAL:HG22	1.53	0.72
2:B:172:ILE:HD13	2:B:178:ASN:ND2	2.04	0.72
1:A:1304:TRP:O	1:A:1305:VAL:HG23	1.89	0.71
4:D:66:ARG:NH2	7:G:48:VAL:HG23	2.03	0.71
7:G:1:MET:HE2	7:G:1:MET:O	1.88	0.71
11:K:5:ASP:OD2	11:K:5:ASP:N	2.21	0.71
1:A:1308:THR:HG23	1:A:1309:ASP:N	2.04	0.71
2:B:165:VAL:HG12	2:B:167:ILE:HD11	1.72	0.71
1:A:1155:ASP:OD2	1:A:1161:THR:HA	1.90	0.71
2:B:457:LEU:O	2:B:461:LEU:HB2	1.90	0.71
7:G:97:HIS:H	7:G:97:HIS:CD2	2.09	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:335:ARG:HA	1:A:339:ASN:HD22	1.55	0.71
7:G:119:LEU:HD12	7:G:131:GLN:O	1.91	0.71
1:A:814:PHE:CE1	2:B:519:TRP:HB2	2.26	0.71
2:B:711:GLU:HB2	2:B:712:PRO:HD3	1.72	0.71
6:F:87:LYS:HG3	6:F:88:TYR:HD1	1.56	0.71
1:A:34:LYS:HE3	1:A:57:ARG:NH1	2.05	0.71
1:A:53:LEU:HD23	1:A:54:ASN:N	2.04	0.71
1:A:427:GLN:HB2	1:A:430:TRP:CG	2.26	0.71
3:C:18:VAL:HG21	11:K:109:TRP:HZ3	1.53	0.71
4:D:51:ASN:HD22	4:D:178:ALA:HA	1.55	0.71
8:H:101:ALA:HB2	8:H:116:TYR:HE2	1.56	0.71
10:J:36:LEU:HB2	10:J:47:ARG:HH12	1.52	0.71
1:A:205:GLU:H	1:A:205:GLU:CD	1.98	0.71
1:A:362:ASP:OD1	1:A:459:ARG:HD3	1.91	0.71
2:B:291:ILE:HG12	2:B:300:HIS:NE2	2.06	0.71
2:B:996:ARG:HH22	3:C:175:ALA:N	1.87	0.71
2:B:1138:MET:HB2	2:B:1147:LEU:HD21	1.72	0.71
12:L:27:LEU:HD22	12:L:37:LYS:HD2	1.72	0.71
1:A:511:ILE:O	1:A:519:PRO:HA	1.91	0.71
1:A:1028:THR:O	1:A:1032:LEU:HD12	1.91	0.71
2:B:1160:VAL:HG12	2:B:1161:HIS:N	2.04	0.71
1:A:21:LEU:HD23	2:B:1212:ILE:HG13	1.73	0.70
1:A:535:THR:HG21	1:A:617:VAL:H	1.55	0.70
1:A:901:LEU:H	1:A:926:GLN:NE2	1.89	0.70
2:B:522:VAL:HG12	2:B:523:CYS:N	2.06	0.70
2:B:800:GLN:HB3	10:J:52:THR:CG2	2.21	0.70
7:G:143:ILE:HG12	7:G:170:ALA:HA	1.71	0.70
1:A:795:GLU:CD	1:A:795:GLU:H	1.99	0.70
1:A:947:PHE:CE1	1:A:954:TRP:CE2	2.79	0.70
2:B:745:PRO:O	2:B:748:ILE:HG12	1.91	0.70
3:C:67:LEU:HD23	3:C:70:ILE:HD11	1.72	0.70
4:D:205:ASP:O	4:D:208:GLU:HB3	1.91	0.70
2:B:313:MET:O	2:B:316:PRO:HD2	1.90	0.70
2:B:705:MET:H	2:B:710:LEU:HD12	1.57	0.70
7:G:56:ILE:HG23	7:G:57:GLN:H	1.57	0.70
11:K:47:ARG:HD3	11:K:59:ALA:O	1.91	0.70
1:A:984:LYS:O	1:A:988:LEU:HB2	1.90	0.70
1:A:451:HIS:CD2	1:A:453:MET:HE2	2.26	0.70
1:A:913:LEU:HD12	1:A:914:GLU:N	2.06	0.70
1:A:59:GLY:C	1:A:244:PRO:HG2	2.16	0.70
1:A:388:LEU:HA	1:A:391:LEU:HD12	1.72	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:69:LEU:HD13	2:B:429:PHE:HD1	1.56	0.70
2:B:973:ILE:CG2	2:B:974:PRO:HD2	2.21	0.70
1:A:265:LYS:O	1:A:269:ILE:HG13	1.91	0.70
2:B:1002:THR:OG1	2:B:1006:ILE:HD12	1.92	0.70
7:G:1:MET:HE3	7:G:80:LYS:O	1.92	0.70
1:A:269:ILE:HG12	1:A:299:HIS:HB3	1.72	0.70
1:A:483:ASP:CB	2:B:987:LYS:HG3	2.22	0.70
2:B:123:THR:OG1	2:B:458:LYS:HE2	1.92	0.70
2:B:573:GLN:O	2:B:575:PRO:HD3	1.91	0.70
2:B:616:ILE:CD1	2:B:625:LYS:HB2	2.21	0.70
3:C:242:GLN:NE2	3:C:246:ARG:HE	1.89	0.70
2:B:190:TYR:CE1	10:J:62:ARG:HD3	2.26	0.70
5:E:138:ALA:HA	5:E:141:VAL:HG23	1.73	0.70
11:K:63:VAL:HG23	11:K:63:VAL:O	1.91	0.69
2:B:273:LEU:HD21	2:B:360:PHE:CD1	2.20	0.69
3:C:121:VAL:HG12	3:C:122:SER:N	2.05	0.69
6:F:101:ILE:HG22	6:F:117:PRO:HB3	1.74	0.69
1:A:705:LYS:HB2	1:A:708:MET:HE3	1.74	0.69
1:A:107:CYS:HA	1:A:171:GLN:NE2	2.07	0.69
4:D:188:ALA:HA	4:D:191:ALA:CB	2.22	0.69
4:D:220:LEU:HD23	4:D:221:TYR:H	1.56	0.69
6:F:109:VAL:HG12	6:F:110:ASP:N	2.05	0.69
1:A:463:ILE:HB	1:A:464:PRO:CD	2.22	0.69
1:A:565:ILE:CG2	1:A:567:LYS:HG2	2.22	0.69
1:A:1118:VAL:HG23	1:A:1306:LEU:HB2	1.75	0.69
1:A:1216:ILE:O	1:A:1219:THR:HG22	1.93	0.69
8:H:112:ILE:CG2	8:H:130:ARG:HH22	2.06	0.69
2:B:54:PHE:CZ	2:B:59:LEU:HD13	2.27	0.69
2:B:398:ARG:CZ	2:B:398:ARG:HB2	2.22	0.69
4:D:12:ARG:HD3	4:D:14:ARG:HG2	1.74	0.69
1:A:382:PRO:HD3	1:A:428:TYR:CE2	2.28	0.69
1:A:1127:ASP:HB3	1:A:1130:GLN:HB3	1.75	0.69
2:B:113:TYR:CD2	2:B:114:PRO:HD2	2.28	0.69
2:B:780:VAL:O	2:B:817:LEU:HD22	1.93	0.69
2:B:1182:CYS:SG	2:B:1182:CYS:O	2.51	0.69
3:C:46:ILE:HD12	3:C:67:LEU:HB3	1.75	0.69
3:C:147:LEU:HB2	3:C:151:GLN:HB2	1.74	0.69
7:G:138:THR:HG22	7:G:139:ILE:H	1.58	0.69
8:H:102:TYR:CE2	8:H:115:TYR:HB3	2.28	0.69
1:A:240:PRO:O	1:A:242:PRO:HD3	1.93	0.69
7:G:31:LEU:CD2	7:G:48:VAL:HG21	2.21	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:298:PHE:CZ	1:A:314:ALA:HB2	2.28	0.68
3:C:37:MET:HE3	3:C:176:ILE:HD13	1.74	0.68
7:G:21:ARG:HD2	7:G:24:GLN:HB3	1.74	0.68
8:H:24:CYS:SG	8:H:44:VAL:HG21	2.33	0.68
1:A:664:THR:HG23	1:A:664:THR:O	1.93	0.68
1:A:997:LEU:HB3	1:A:1053:PHE:CE2	2.28	0.68
1:A:1217:LYS:O	1:A:1221:LYS:HA	1.92	0.68
2:B:1007:VAL:HG23	2:B:1008:PRO:HD2	1.74	0.68
3:C:58:LEU:HD21	10:J:57:ILE:HD13	1.75	0.68
1:A:1036:ARG:HG2	1:A:1036:ARG:NH1	2.06	0.68
1:A:1107:VAL:HG12	1:A:1107:VAL:O	1.94	0.68
2:B:363:HIS:CD2	2:B:585:VAL:HG22	2.28	0.68
4:D:190:GLU:HA	7:G:167:TYR:CD1	2.27	0.68
2:B:237:VAL:HG22	2:B:257:LYS:HA	1.76	0.68
2:B:1073:TYR:CE2	2:B:1080:LYS:HG2	2.29	0.68
4:D:179:GLN:NE2	7:G:1:MET:HB2	2.07	0.68
11:K:51:LEU:HD11	11:K:59:ALA:HB3	1.76	0.68
3:C:67:LEU:O	3:C:70:ILE:HG13	1.94	0.68
3:C:174:ALA:HA	10:J:10:CYS:O	1.94	0.68
12:L:66:GLN:C	12:L:67:PHE:CD1	2.72	0.68
1:A:1402:PHE:C	1:A:1403:GLU:HG2	2.19	0.68
3:C:75:MET:HG3	3:C:76:ASP:OD1	1.92	0.68
3:C:112:ASN:HB3	3:C:114:TYR:CE1	2.29	0.68
8:H:40:LEU:HD13	8:H:123:MET:HG3	1.75	0.68
8:H:82:PRO:C	8:H:84:ALA:H	2.00	0.68
1:A:53:LEU:CD2	1:A:54:ASN:HD22	2.06	0.68
1:A:251:SER:HA	1:A:257:ARG:O	1.94	0.68
1:A:584:ASN:O	1:A:637:LYS:HE2	1.93	0.68
1:A:1101:LEU:HA	1:A:1104:ILE:HD12	1.75	0.68
1:A:1242:VAL:HG12	1:A:1243:VAL:N	2.09	0.68
2:B:731:VAL:O	2:B:732:SER:HB2	1.93	0.68
2:B:797:TYR:HB3	2:B:798:TYR:HD2	1.59	0.68
1:A:1404:GLU:O	1:A:1408:ILE:HG13	1.94	0.68
1:A:434:ARG:NH2	1:A:437:MET:HG3	2.09	0.67
3:C:63:ILE:CA	3:C:66:ARG:HG3	2.23	0.67
8:H:26:ILE:HG22	8:H:40:LEU:O	1.94	0.67
1:A:266:LEU:HD21	1:A:303:TYR:CE1	2.28	0.67
1:A:492:PRO:O	1:A:493:GLN:NE2	2.26	0.67
1:A:869:GLY:O	5:E:204:THR:HG21	1.94	0.67
2:B:785:TYR:HE2	10:J:60:PHE:CE1	2.12	0.67
11:K:46:ILE:O	11:K:50:LEU:HB2	1.94	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:860:MET:HB2	2:B:965:LYS:HG2	1.76	0.67
6:F:111:LEU:HD21	6:F:120:ILE:HD11	1.76	0.67
1:A:34:LYS:CE	1:A:57:ARG:HH12	2.08	0.67
2:B:483:LEU:HD11	2:B:491:THR:HG23	1.76	0.67
2:B:611:PRO:HB3	2:B:685:LEU:HD11	1.76	0.67
11:K:113:THR:HG22	11:K:114:LEU:N	2.09	0.67
1:A:666:ILE:HD12	1:A:667:GLY:H	1.60	0.67
1:A:933:TYR:O	1:A:933:TYR:CD2	2.44	0.67
1:A:1100:ARG:HH12	1:A:1111:MET:HE2	1.58	0.67
2:B:955:THR:HG22	2:B:956:THR:H	1.60	0.67
2:B:1076:HIS:HD2	11:K:40:HIS:NE2	1.91	0.67
6:F:129:LYS:C	6:F:130:ILE:HG13	2.20	0.67
1:A:91:PHE:H	1:A:297:GLN:HE22	1.41	0.67
1:A:330:LYS:O	1:A:334:GLY:HA3	1.95	0.67
1:A:350:ARG:C	1:A:351:THR:HG22	2.18	0.67
1:A:845:LEU:HA	1:A:848:ILE:HD12	1.76	0.67
2:B:973:ILE:HG22	2:B:974:PRO:HD2	1.77	0.67
3:C:169:LYS:NZ	12:L:69:ALA:HB3	2.10	0.67
5:E:15:ALA:HA	5:E:140:LEU:O	1.93	0.67
1:A:49:LYS:HZ1	1:A:61:ILE:HG13	1.58	0.67
1:A:567:LYS:NZ	8:H:46:LEU:HB2	2.10	0.67
5:E:180:ARG:HB2	5:E:215:MET:OXT	1.95	0.67
1:A:219:PHE:CZ	1:A:231:PRO:HD2	2.30	0.67
1:A:1112:LYS:O	1:A:1114:PRO:HD3	1.95	0.67
2:B:102:VAL:HB	2:B:112:LEU:HB2	1.77	0.67
2:B:215:GLN:HB2	2:B:407:ASP:HB2	1.76	0.67
2:B:496:ARG:HH12	2:B:541:LEU:HA	1.60	0.67
3:C:45:ALA:HA	3:C:72:LEU:HD22	1.77	0.67
2:B:817:LEU:N	2:B:818:PRO:HD3	2.09	0.67
2:B:1138:MET:CB	2:B:1147:LEU:HD21	2.25	0.67
5:E:213:ILE:HG12	5:E:214:CYS:N	2.10	0.67
6:F:103:MET:HE3	7:G:66:GLY:H	1.59	0.67
8:H:139:ASN:O	8:H:140:ALA:HB2	1.95	0.67
1:A:649:ILE:O	1:A:653:VAL:HG23	1.95	0.67
1:A:760:GLN:HE21	1:A:765:VAL:HG13	1.59	0.67
1:A:1259:MET:HA	1:A:1262:LYS:HD2	1.76	0.67
4:D:167:LEU:C	4:D:169:SER:H	2.03	0.67
1:A:32:VAL:HG23	1:A:33:ALA:H	1.59	0.66
1:A:69:THR:HG21	2:B:1174:LYS:NZ	2.08	0.66
1:A:589:GLN:HA	1:A:605:MET:O	1.94	0.66
2:B:36:ALA:O	2:B:39:ARG:HB2	1.95	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:J:57:ILE:HG23	10:J:58:GLU:N	2.11	0.66
1:A:262:LEU:HD12	1:A:328:ARG:NH2	2.10	0.66
1:A:963:ILE:HD11	1:A:1048:ASN:HB3	1.77	0.66
1:A:1445:ILE:HD12	1:A:1445:ILE:O	1.94	0.66
2:B:363:HIS:O	2:B:364:ILE:HB	1.95	0.66
7:G:144:ARG:O	7:G:168:LEU:HD22	1.95	0.66
10:J:12:LYS:O	10:J:14:VAL:HG23	1.95	0.66
2:B:559:SER:HA	2:B:563:MET:CB	2.25	0.66
2:B:824:ILE:HB	2:B:1009:ASP:OD2	1.94	0.66
7:G:14:HIS:CD2	7:G:15:PRO:HD2	2.30	0.66
1:A:351:THR:HG23	1:A:468:PHE:CD1	2.31	0.66
6:F:120:ILE:HG22	6:F:121:ALA:N	2.10	0.66
11:K:32:VAL:HG23	11:K:74:ARG:HG3	1.77	0.66
1:A:152:VAL:HG13	1:A:153:PRO:HD2	1.77	0.66
1:A:1138:ILE:O	1:A:1140:HIS:N	2.29	0.66
1:A:361:LEU:CA	1:A:471:ASN:HD22	2.07	0.66
2:B:1114:LEU:CD1	2:B:1202:LEU:HD11	2.26	0.66
4:D:40:HIS:ND1	7:G:6:ASP:HB3	2.10	0.66
1:A:494:SER:O	1:A:497:THR:HB	1.96	0.66
2:B:643:ASP:OD2	2:B:652:LYS:HE3	1.95	0.66
2:B:806:THR:HG22	2:B:808:ALA:H	1.61	0.66
2:B:886:LYS:HE2	2:B:940:PRO:HG3	1.76	0.66
2:B:1107:ALA:O	2:B:1108:ARG:CB	2.44	0.66
12:L:47:ARG:O	12:L:48:CYS:HB2	1.95	0.66
1:A:359:LEU:HD23	1:A:359:LEU:H	1.59	0.66
1:A:1100:ARG:O	1:A:1104:ILE:HG13	1.96	0.66
1:A:1119:TYR:CD2	1:A:1305:VAL:HG21	2.31	0.66
2:B:810:GLU:CA	2:B:815:ARG:HH22	2.08	0.66
2:B:1160:VAL:HG12	2:B:1161:HIS:H	1.59	0.66
5:E:177:ARG:HD3	5:E:215:MET:HE2	1.77	0.66
6:F:89:GLU:O	6:F:93:ILE:HG13	1.96	0.66
2:B:418:LYS:HG2	2:B:422:LYS:HE3	1.78	0.66
7:G:21:ARG:NH1	7:G:24:GLN:HB2	2.10	0.66
1:A:672:ASP:HB3	1:A:736:ASN:ND2	2.10	0.66
2:B:204:ILE:HG23	2:B:207:GLY:O	1.95	0.66
2:B:521:LEU:HB3	2:B:633:VAL:HG11	1.77	0.66
3:C:112:ASN:ND2	3:C:146:LYS:HG2	2.10	0.66
4:D:48:ILE:HG22	4:D:48:ILE:O	1.94	0.66
12:L:55:ILE:HD13	12:L:55:ILE:H	1.60	0.66
1:A:365:GLY:HA3	1:A:469:ARG:HB2	1.78	0.65
1:A:1120:LEU:HD22	1:A:1124:HIS:O	1.95	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:879:ARG:HE	2:B:879:ARG:H	1.44	0.65
4:D:52:LEU:HD12	4:D:182:SER:HB2	1.76	0.65
1:A:350:ARG:N	2:B:1128:LEU:HD11	2.11	0.65
1:A:622:VAL:O	1:A:630:ILE:HD11	1.96	0.65
1:A:1004:ASN:ND2	1:A:1007:ILE:HG13	2.11	0.65
1:A:1255:GLU:CD	1:A:1258:HIS:HB2	2.20	0.65
4:D:66:ARG:O	4:D:70:PHE:HB2	1.96	0.65
6:F:84:TYR:HA	6:F:152:ILE:HD12	1.78	0.65
1:A:58:LEU:CD1	1:A:244:PRO:HD3	2.26	0.65
1:A:445:ASN:CG	1:A:455:MET:HE2	2.21	0.65
2:B:364:ILE:HG12	2:B:585:VAL:HG13	1.77	0.65
2:B:1194:ILE:HG12	2:B:1196:ILE:HG22	1.77	0.65
10:J:5:VAL:HA	10:J:15:GLY:H	1.61	0.65
1:A:27:VAL:C	1:A:29:ALA:H	2.04	0.65
1:A:1453:TYR:CE2	6:F:129:LYS:HA	2.32	0.65
2:B:299:GLU:HB3	2:B:571:PRO:HG2	1.77	0.65
1:A:549:MET:HE1	1:A:656:TRP:HD1	1.60	0.65
1:A:1434:ALA:O	1:A:1436:ILE:N	2.29	0.65
2:B:227:LYS:H	2:B:395:GLN:NE2	1.94	0.65
4:D:46:GLU:HG2	4:D:47:LEU:H	1.62	0.65
9:I:45:ARG:HG3	9:I:46:HIS:N	2.12	0.65
1:A:709:THR:HG22	1:A:710:LEU:N	2.12	0.65
1:A:1030:ARG:HG2	1:A:1034:GLU:OE2	1.95	0.65
1:A:1227:ILE:C	1:A:1228:TRP:HE3	2.04	0.65
1:A:30:ILE:HD11	2:B:1168:LEU:HD13	1.78	0.65
1:A:112:LYS:HG2	1:A:113:LEU:H	1.62	0.65
1:A:569:LYS:HD2	3:C:221:TYR:HB2	1.79	0.65
1:A:714:PHE:O	1:A:718:VAL:HG23	1.96	0.65
1:A:966:ASN:O	1:A:970:THR:HB	1.97	0.65
1:A:1259:MET:HE2	1:A:1263:ILE:HG12	1.78	0.65
1:A:1308:THR:HG21	1:A:1310:GLY:O	1.97	0.65
2:B:1065:GLN:HG3	2:B:1068:GLY:H	1.62	0.65
4:D:8:PHE:CZ	4:D:37:GLN:HB2	2.31	0.65
5:E:198:ILE:HD13	5:E:212:ARG:HH11	1.62	0.65
1:A:864:ILE:HD12	1:A:864:ILE:N	2.11	0.65
1:A:1223:ASP:HA	1:A:1243:VAL:CG2	2.27	0.65
1:A:1349:TYR:O	1:A:1350:LYS:C	2.40	0.65
2:B:488:TYR:CE2	2:B:813:LYS:HB2	2.31	0.65
2:B:488:TYR:HE2	2:B:813:LYS:HB2	1.61	0.65
5:E:78:LEU:HG	5:E:107:THR:HG21	1.77	0.65
6:F:94:LEU:HD21	6:F:122:MET:HG2	1.76	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:7:SER:OG	2:B:1161:HIS:HE1	1.80	0.65
2:B:780:VAL:HG11	10:J:56:LEU:CD1	2.27	0.65
3:C:55:THR:OG1	3:C:152:GLU:N	2.30	0.65
3:C:197:SER:O	3:C:201:TRP:HB2	1.97	0.65
1:A:532:ARG:HD2	1:A:749:ALA:HB2	1.77	0.65
1:A:1027:ALA:O	1:A:1031:VAL:HG23	1.96	0.65
3:C:112:ASN:ND2	3:C:143:LEU:HD21	2.12	0.65
6:F:86:THR:HG23	6:F:89:GLU:CD	2.23	0.65
1:A:1280:GLU:O	1:A:1282:VAL:HG23	1.97	0.64
2:B:758:PHE:CE2	2:B:1044:ALA:HA	2.31	0.64
1:A:2:VAL:O	2:B:1158:PHE:HA	1.97	0.64
1:A:537:ARG:HD2	8:H:20:TYR:CZ	2.32	0.64
2:B:549:THR:HG22	2:B:550:ASP:H	1.61	0.64
2:B:638:PHE:CE1	2:B:743:ILE:HA	2.32	0.64
1:A:467:THR:HG21	2:B:976:ILE:HG22	1.79	0.64
1:A:982:THR:HG22	1:A:984:LYS:H	1.61	0.64
6:F:140:ASP:CG	6:F:141:GLY:N	2.56	0.64
1:A:58:LEU:HD12	1:A:244:PRO:HD3	1.78	0.64
2:B:169:ARG:O	2:B:457:LEU:HD12	1.98	0.64
2:B:430:ARG:HG2	2:B:430:ARG:NH1	2.04	0.64
2:B:807:ARG:HG2	2:B:1045:SER:OG	1.97	0.64
2:B:1169:MET:HE1	2:B:1201:LYS:HA	1.78	0.64
5:E:10:SER:O	5:E:13:TRP:HB3	1.97	0.64
5:E:55:ARG:C	5:E:57:MET:H	2.05	0.64
1:A:642:CYS:O	1:A:645:LEU:HB3	1.97	0.64
1:A:1291:VAL:HG22	1:A:1292:PRO:HD2	1.79	0.64
5:E:61:GLN:HG2	5:E:62:ALA:H	1.63	0.64
5:E:100:ILE:HG23	5:E:105:PHE:CD1	2.32	0.64
1:A:1118:VAL:O	1:A:1305:VAL:HG13	1.97	0.64
5:E:19:VAL:O	5:E:23:VAL:HG23	1.97	0.64
11:K:65:HIS:CD2	11:K:65:HIS:C	2.76	0.64
1:A:135:PHE:HD1	1:A:222:LEU:CD2	2.10	0.64
1:A:441:PRO:HG2	1:A:498:ARG:HB2	1.79	0.64
2:B:1162:ILE:CD1	2:B:1194:ILE:HG21	2.27	0.64
5:E:187:TYR:HD2	5:E:188:LEU:CD2	2.11	0.64
6:F:86:THR:HG23	6:F:89:GLU:OE1	1.98	0.64
1:A:1226:VAL:HG22	1:A:1240:CYS:HB3	1.80	0.64
2:B:213:ILE:CG2	2:B:499:ASN:HB2	2.27	0.64
2:B:798:TYR:CE1	10:J:4:PRO:HB3	2.33	0.64
2:B:955:THR:CG2	2:B:956:THR:N	2.61	0.64
2:B:1096:ARG:HG2	2:B:1097:HIS:H	1.61	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:G:158:HIS:HD2	7:G:159:ALA:N	1.96	0.64
11:K:23:PRO:HA	11:K:31:VAL:HG13	1.80	0.64
1:A:91:PHE:HZ	1:A:207:ILE:HD12	1.62	0.64
1:A:93:VAL:CG2	1:A:301:ALA:HA	2.28	0.64
1:A:332:LYS:HA	1:A:337:ARG:HB3	1.80	0.64
1:A:679:ILE:HG23	1:A:729:ALA:HB1	1.80	0.64
1:A:901:LEU:HD22	1:A:919:ILE:HG23	1.79	0.64
1:A:115:LEU:HB3	1:A:122:MET:HG2	1.79	0.64
1:A:741:ASN:HD22	1:A:741:ASN:C	2.06	0.64
11:K:46:ILE:HG22	11:K:50:LEU:HD12	1.78	0.64
11:K:65:HIS:CD2	11:K:67:PHE:N	2.55	0.64
1:A:1130:GLN:HA	1:A:1133:LEU:HD12	1.80	0.63
1:A:1259:MET:HE3	1:A:1262:LYS:HD2	1.80	0.63
2:B:69:LEU:HB3	2:B:429:PHE:HE1	1.63	0.63
2:B:635:ARG:HB2	2:B:636:PRO:HD2	1.79	0.63
2:B:948:ILE:HG22	2:B:949:VAL:O	1.98	0.63
2:B:1116:ARG:HG3	2:B:1198:TYR:CE2	2.32	0.63
2:B:1195:HIS:C	2:B:1196:ILE:HG22	2.24	0.63
4:D:194:LEU:N	4:D:194:LEU:HD23	2.13	0.63
1:A:1223:ASP:HA	1:A:1243:VAL:HG21	1.80	0.63
3:C:182:PRO:HD2	3:C:210:GLU:OE1	1.98	0.63
3:C:191:TYR:CD2	3:C:201:TRP:CD1	2.87	0.63
1:A:93:VAL:HG21	1:A:304:MET:HB2	1.80	0.63
1:A:336:ILE:HD11	2:B:1203:LEU:HD13	1.79	0.63
1:A:685:GLU:HG3	1:A:686:ALA:N	2.12	0.63
2:B:833:TYR:C	2:B:835:GLN:H	2.05	0.63
3:C:260:LEU:O	3:C:264:GLN:HG3	1.98	0.63
5:E:41:ASP:HA	5:E:44:ALA:HB3	1.79	0.63
1:A:256:GLN:O	1:A:257:ARG:CB	2.47	0.63
1:A:997:LEU:HD22	1:A:1053:PHE:CG	2.33	0.63
1:A:1012:ARG:O	1:A:1013:ASP:C	2.42	0.63
2:B:702:LEU:HD23	2:B:738:PHE:N	2.14	0.63
2:B:1115:THR:O	2:B:1116:ARG:HB2	1.98	0.63
4:D:60:LYS:O	4:D:64:VAL:HG23	1.98	0.63
4:D:159:THR:O	4:D:163:VAL:HG23	1.98	0.63
4:D:163:VAL:HG13	4:D:214:LEU:HD21	1.79	0.63
7:G:42:PHE:O	7:G:80:LYS:HB2	1.99	0.63
7:G:88:ASP:HB3	7:G:144:ARG:HA	1.80	0.63
1:A:836:TYR:CZ	1:A:840:ARG:HD2	2.33	0.63
2:B:619:ILE:HG22	2:B:620:ARG:N	2.13	0.63
2:B:859:TYR:HD1	2:B:859:TYR:H	1.47	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:265:LYS:HE3	1:A:265:LYS:CA	2.27	0.63
2:B:277:LYS:HG3	2:B:336:ARG:HG2	1.79	0.63
2:B:469:GLN:C	2:B:472:ALA:HB3	2.23	0.63
2:B:696:GLU:O	2:B:699:GLU:HB2	1.97	0.63
2:B:842:ASN:C	2:B:842:ASN:HD22	2.06	0.63
1:A:59:GLY:CA	1:A:244:PRO:HG2	2.29	0.63
1:A:634:THR:O	1:A:638:GLY:HA2	1.99	0.63
1:A:999:VAL:HG12	1:A:1000:LEU:HG	1.79	0.63
1:A:1227:ILE:HA	1:A:1228:TRP:CE3	2.34	0.63
10:J:25:LEU:HD21	10:J:32:GLU:HG3	1.81	0.63
11:K:40:HIS:NE2	11:K:63:VAL:HG11	2.13	0.63
12:L:53:HIS:HB3	12:L:55:ILE:HD13	1.80	0.63
1:A:767:GLN:NE2	1:A:774:ARG:HB2	2.14	0.62
2:B:650:GLU:HG3	2:B:651:LEU:N	2.14	0.62
5:E:117:THR:HB	5:E:120:ALA:HB2	1.81	0.62
11:K:45:LEU:O	11:K:48:ALA:HB3	1.99	0.62
12:L:30:ILE:HD11	12:L:59:ALA:HB2	1.81	0.62
1:A:399:HIS:CB	1:A:400:PRO:HD3	2.28	0.62
1:A:535:THR:HG21	1:A:616:VAL:HA	1.80	0.62
2:B:126:SER:OG	2:B:172:ILE:HD11	1.98	0.62
2:B:948:ILE:O	2:B:968:VAL:HG13	1.99	0.62
3:C:44:LEU:HD23	3:C:44:LEU:C	2.24	0.62
7:G:79:PHE:CD1	7:G:79:PHE:C	2.72	0.62
8:H:127:GLY:O	8:H:128:ASN:HB2	1.98	0.62
1:A:1107:VAL:CG2	1:A:1383:SER:HB3	2.29	0.62
2:B:613:VAL:CG2	2:B:628:THR:HA	2.30	0.62
2:B:830:TYR:CE2	2:B:1000:PRO:HD3	2.35	0.62
10:J:48:ARG:HD2	10:J:48:ARG:O	1.99	0.62
1:A:82:GLY:O	1:A:241:VAL:HB	1.99	0.62
1:A:262:LEU:HD12	1:A:328:ARG:CZ	2.30	0.62
1:A:403:LYS:O	1:A:415:LEU:HB2	2.00	0.62
3:C:169:LYS:HZ2	12:L:69:ALA:HB3	1.64	0.62
6:F:128:LYS:HD3	6:F:148:VAL:O	1.98	0.62
7:G:30:LEU:HD13	7:G:72:VAL:HG11	1.81	0.62
9:I:46:HIS:O	9:I:47:GLU:HG2	1.99	0.62
1:A:377:PRO:HB2	1:A:431:LYS:NZ	2.15	0.62
2:B:770:GLN:HG2	2:B:983:ARG:O	2.00	0.62
2:B:770:GLN:CD	2:B:983:ARG:HA	2.24	0.62
4:D:147:TYR:O	4:D:151:PHE:CD1	2.51	0.62
6:F:100:GLN:O	6:F:105:ALA:CB	2.47	0.62
1:A:814:PHE:HE1	2:B:519:TRP:HB2	1.64	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:E:156:LEU:HG	5:E:195:VAL:O	2.00	0.62
7:G:17:PHE:CE2	7:G:25:TYR:HE2	2.16	0.62
4:D:147:TYR:O	4:D:151:PHE:HD1	1.82	0.62
8:H:145:ARG:O	8:H:146:ARG:HB2	1.99	0.62
1:A:49:LYS:NZ	1:A:61:ILE:H	1.98	0.62
1:A:477:PRO:HG2	1:A:521:MET:HE2	1.81	0.62
1:A:1119:TYR:HA	1:A:1305:VAL:HG22	1.80	0.62
1:A:1207:LEU:HD13	1:A:1273:LEU:HD23	1.82	0.62
1:A:1299:VAL:HG12	1:A:1300:LYS:H	1.65	0.62
2:B:638:PHE:HA	2:B:690:VAL:HG12	1.81	0.62
2:B:898:LEU:HD21	2:B:964:VAL:HG21	1.82	0.62
2:B:1198:TYR:CE1	2:B:1201:LYS:HD2	2.34	0.62
5:E:16:PHE:CZ	5:E:20:LYS:HE3	2.34	0.62
7:G:90:THR:HG23	7:G:140:LYS:O	2.00	0.62
1:A:58:LEU:HB3	1:A:80:HIS:O	1.99	0.62
1:A:1208:THR:HB	1:A:1211:GLN:HG3	1.82	0.62
1:A:1210:GLY:HA2	1:A:1228:TRP:NE1	2.14	0.62
3:C:220:ASP:OD2	3:C:223:ALA:HB2	2.00	0.62
5:E:61:GLN:HG2	5:E:62:ALA:N	2.14	0.62
11:K:63:VAL:O	11:K:64:GLU:C	2.42	0.62
2:B:174:LEU:HD21	2:B:204:ILE:CD1	2.29	0.62
2:B:877:PRO:HA	2:B:934:LYS:HZ3	1.62	0.62
2:B:879:ARG:HE	2:B:879:ARG:N	1.97	0.62
11:K:53:ASP:HB3	11:K:56:VAL:HB	1.82	0.62
1:A:604:GLY:O	1:A:605:MET:HB2	2.00	0.61
1:A:1107:VAL:HG23	1:A:1383:SER:HB3	1.80	0.61
2:B:515:HIS:HD2	2:B:517:THR:N	1.90	0.61
5:E:119:SER:O	5:E:122:LYS:HB2	2.00	0.61
1:A:40:THR:HG23	1:A:54:ASN:HD21	1.63	0.61
1:A:506:ALA:CB	1:A:508:PRO:HD2	2.28	0.61
1:A:601:LYS:HD3	1:A:603:ASN:HD21	1.65	0.61
1:A:1154:TYR:CZ	1:A:1156:PRO:HG3	2.34	0.61
9:I:35:VAL:HG12	9:I:36:GLU:N	2.12	0.61
1:A:224:PHE:HD2	1:A:231:PRO:HD3	1.63	0.61
1:A:317:LYS:O	1:A:318:SER:HB3	1.99	0.61
1:A:419:LYS:HG3	1:A:420:ARG:N	2.14	0.61
1:A:621:THR:O	1:A:629:LEU:HB2	2.00	0.61
1:A:1153:TYR:HB2	1:A:1192:LEU:HD23	1.81	0.61
2:B:1166:CYS:C	2:B:1168:LEU:H	2.08	0.61
3:C:237:SER:O	3:C:238:ILE:HG13	1.99	0.61
11:K:63:VAL:O	11:K:65:HIS:N	2.33	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:108:MET:HE3	1:A:210:ILE:HD12	1.83	0.61
2:B:234:ILE:HG12	2:B:257:LYS:HB3	1.83	0.61
2:B:486:TYR:CD1	2:B:1096:ARG:NH2	2.68	0.61
2:B:778:MET:HE2	2:B:1094:ARG:HG2	1.82	0.61
7:G:45:ILE:O	7:G:45:ILE:HG22	1.99	0.61
2:B:68:THR:HG22	2:B:91:SER:HA	1.82	0.61
2:B:105:SER:O	2:B:106:ASP:HB2	1.99	0.61
2:B:530:GLY:C	2:B:531:GLN:HG2	2.25	0.61
3:C:3:GLU:HB3	11:K:104:ASN:HD21	1.65	0.61
4:D:29:LEU:HD23	4:D:33:PHE:HB3	1.81	0.61
8:H:130:ARG:HH11	8:H:130:ARG:N	1.92	0.61
12:L:62:LYS:O	12:L:64:LEU:N	2.32	0.61
1:A:253:ASN:HB2	2:B:884:ARG:NH1	2.14	0.61
1:A:836:TYR:CE2	1:A:840:ARG:HD2	2.36	0.61
1:A:1146:VAL:O	1:A:1197:LEU:HA	2.00	0.61
2:B:129:PHE:CE2	2:B:166:PHE:HB2	2.36	0.61
2:B:309:GLN:O	2:B:312:GLU:HB3	2.00	0.61
2:B:635:ARG:HB2	2:B:636:PRO:CD	2.30	0.61
2:B:955:THR:CG2	2:B:956:THR:H	2.12	0.61
2:B:999:MET:HG2	2:B:1007:VAL:CG2	2.30	0.61
7:G:115:MET:HB3	7:G:116:PRO:HD2	1.83	0.61
12:L:47:ARG:HH21	12:L:54:ARG:HG2	1.65	0.61
1:A:1079:MET:HG2	1:A:1359:ASP:OD1	2.01	0.61
1:A:1102:LYS:HG2	1:A:1106:ASN:ND2	2.16	0.61
3:C:262:LEU:HD13	11:K:88:LYS:HG2	1.82	0.61
5:E:178:ILE:HB	5:E:212:ARG:HD3	1.82	0.61
6:F:111:LEU:O	6:F:113:GLY:N	2.33	0.61
9:I:17:ARG:HG3	9:I:18:GLU:H	1.65	0.61
12:L:40:LEU:HD13	12:L:44:ASP:HB3	1.82	0.61
1:A:969:GLN:O	1:A:969:GLN:HG2	2.01	0.61
2:B:562:GLY:HA3	2:B:590:HIS:ND1	2.14	0.61
2:B:644:GLU:HG3	2:B:646:LEU:H	1.66	0.61
2:B:656:GLY:O	2:B:660:LYS:HB2	1.99	0.61
7:G:13:LEU:HD23	7:G:14:HIS:N	2.15	0.61
9:I:25:LEU:CB	9:I:38:ALA:HB2	2.11	0.61
10:J:9:SER:HB2	10:J:45:CYS:HB2	1.82	0.61
1:A:321:PRO:O	1:A:322:VAL:HG12	2.00	0.61
1:A:370:ILE:HG22	1:A:374:LEU:CD1	2.30	0.61
1:A:873:MET:O	1:A:1058:VAL:HG23	2.01	0.61
1:A:1199:ARG:HG3	1:A:1236:LEU:HD11	1.81	0.61
1:A:1441:PHE:CZ	6:F:89:GLU:HA	2.35	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:639:ILE:HG22	2:B:641:GLU:HG2	1.81	0.61
3:C:80:LEU:HD12	3:C:94:LYS:O	2.01	0.61
3:C:124:LEU:HD23	3:C:127:ARG:H	1.65	0.61
9:I:7:CYS:HB3	9:I:14:LEU:HD21	1.82	0.61
1:A:399:HIS:ND1	1:A:399:HIS:C	2.58	0.61
1:A:648:ASN:O	1:A:652:VAL:HG23	2.00	0.61
1:A:1259:MET:O	1:A:1263:ILE:HG13	2.01	0.61
2:B:791:THR:HA	2:B:858:SER:HB2	1.82	0.61
2:B:1131:GLY:O	2:B:1132:GLU:C	2.44	0.61
3:C:186:LEU:HD21	3:C:225:ALA:HB2	1.82	0.61
4:D:120:GLU:O	4:D:123:LEU:HB2	2.01	0.61
1:A:709:THR:HG22	1:A:710:LEU:H	1.66	0.60
2:B:841:MET:HE3	2:B:1010:LEU:HG	1.83	0.60
7:G:30:LEU:CD1	7:G:72:VAL:HG11	2.31	0.60
10:J:8:PHE:N	10:J:8:PHE:CD2	2.56	0.60
1:A:42:ASP:HB3	1:A:45:GLN:HA	1.83	0.60
1:A:219:PHE:CE2	1:A:231:PRO:HD2	2.36	0.60
1:A:463:ILE:CB	1:A:464:PRO:HD2	2.27	0.60
1:A:741:ASN:ND2	1:A:743:VAL:N	2.49	0.60
1:A:1263:ILE:HA	1:A:1266:THR:HB	1.83	0.60
2:B:168:GLY:H	2:B:450:ALA:HB1	1.65	0.60
2:B:1096:ARG:CZ	2:B:1096:ARG:HB2	2.30	0.60
8:H:104:PHE:CE2	8:H:114:VAL:HG12	2.36	0.60
2:B:69:LEU:HD22	2:B:429:PHE:CE1	2.37	0.60
2:B:1013:ASN:HD21	2:B:1015:HIS:HD2	1.48	0.60
6:F:100:GLN:NE2	7:G:18:PHE:HE2	1.99	0.60
6:F:139:PRO:O	6:F:140:ASP:C	2.44	0.60
1:A:75:ASN:O	1:A:76:GLU:CB	2.48	0.60
1:A:106:VAL:HG13	1:A:112:LYS:O	2.01	0.60
1:A:640:GLN:CD	1:A:640:GLN:N	2.57	0.60
1:A:902:LEU:HD23	1:A:921:GLY:HA2	1.82	0.60
1:A:1010:ALA:O	1:A:1013:ASP:HB2	2.00	0.60
2:B:493:SER:HA	2:B:751:VAL:HG21	1.82	0.60
2:B:777:ALA:HA	2:B:1095:LEU:HB3	1.82	0.60
7:G:120:THR:OG1	7:G:131:GLN:HB2	2.00	0.60
1:A:102:VAL:HG11	1:A:211:PHE:HE1	1.66	0.60
1:A:359:LEU:HD23	1:A:359:LEU:N	2.15	0.60
2:B:549:THR:HG22	2:B:550:ASP:N	2.17	0.60
9:I:17:ARG:HH21	9:I:30:ARG:NH2	2.00	0.60
10:J:3:VAL:HG21	10:J:18:TRP:HB2	1.83	0.60
1:A:84:ILE:O	1:A:84:ILE:HG22	2.01	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:679:ILE:HD11	1:A:733:ALA:HB2	1.82	0.60
2:B:959:ASP:OD2	2:B:961:LEU:HD12	2.02	0.60
7:G:51:TYR:CD2	7:G:51:TYR:O	2.55	0.60
1:A:567:LYS:HE3	8:H:46:LEU:HD12	1.83	0.60
1:A:744:LYS:O	1:A:748:MET:HB2	2.01	0.60
1:A:873:MET:C	1:A:1058:VAL:HG23	2.26	0.60
2:B:210:LYS:HD2	2:B:480:SER:OG	2.01	0.60
2:B:508:LEU:O	2:B:509:ALA:HB3	2.01	0.60
2:B:841:MET:HG2	2:B:846:ILE:HD11	1.84	0.60
4:D:190:GLU:O	4:D:194:LEU:HG	2.01	0.60
5:E:122:LYS:C	5:E:123:LEU:HG	2.25	0.60
8:H:102:TYR:HD2	8:H:102:TYR:N	1.98	0.60
1:A:1152:ILE:HG23	1:A:1260:LEU:HD23	1.81	0.60
12:L:38:LEU:O	12:L:39:SER:CB	2.49	0.60
12:L:66:GLN:C	12:L:67:PHE:HD1	2.10	0.60
1:A:118:HIS:O	1:A:119:ASN:HB2	2.01	0.60
1:A:528:LEU:HG	1:A:529:CYS:N	2.16	0.60
5:E:14:ARG:HB3	5:E:141:VAL:O	2.02	0.60
5:E:56:LYS:HE2	5:E:84:ASP:HB2	1.84	0.60
1:A:346:ASP:OD1	2:B:1108:ARG:HA	2.02	0.60
1:A:1141:THR:HG23	1:A:1205:LYS:HD3	1.83	0.60
2:B:459:TYR:C	2:B:459:TYR:CD2	2.80	0.60
3:C:121:VAL:HG12	3:C:122:SER:H	1.66	0.60
4:D:8:PHE:HZ	4:D:37:GLN:HB2	1.67	0.60
8:H:59:ILE:O	8:H:60:ALA:HB3	2.01	0.60
1:A:49:LYS:HD3	1:A:54:ASN:O	2.02	0.59
1:A:135:PHE:HA	1:A:138:ILE:HD12	1.83	0.59
1:A:1208:THR:HG23	1:A:1231:ASP:OD2	2.02	0.59
2:B:879:ARG:N	2:B:879:ARG:NE	2.49	0.59
2:B:1023:VAL:HG12	2:B:1027:ILE:HD11	1.82	0.59
3:C:76:ASP:C	3:C:129:ILE:HD11	2.26	0.59
4:D:192:LYS:HG2	4:D:198:LEU:HD12	1.83	0.59
5:E:122:LYS:O	5:E:123:LEU:HG	2.02	0.59
6:F:109:VAL:HG12	6:F:110:ASP:H	1.65	0.59
1:A:1208:THR:HG22	1:A:1210:GLY:H	1.67	0.59
1:A:1291:VAL:HG13	1:A:1292:PRO:HD2	1.83	0.59
2:B:34:ILE:HG23	2:B:542:MET:HE1	1.83	0.59
2:B:126:SER:CB	2:B:172:ILE:HD11	2.32	0.59
2:B:652:LYS:HD3	2:B:688:GLY:O	2.01	0.59
3:C:13:ALA:HB1	3:C:18:VAL:HG22	1.83	0.59
4:D:59:ILE:O	4:D:63:LEU:HB2	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:E:124:VAL:O	5:E:132:ILE:HD13	2.02	0.59
6:F:111:LEU:HD23	6:F:114:GLU:O	2.02	0.59
7:G:158:HIS:CD2	7:G:159:ALA:N	2.69	0.59
2:B:638:PHE:HB3	2:B:651:LEU:HD22	1.83	0.59
3:C:147:LEU:N	3:C:147:LEU:HD23	2.16	0.59
3:C:163:ILE:HD12	3:C:165:LYS:HB2	1.84	0.59
9:I:6:PHE:N	9:I:6:PHE:CD1	2.69	0.59
9:I:25:LEU:HB3	9:I:38:ALA:CB	2.11	0.59
7:G:126:ASN:C	7:G:126:ASN:HD22	2.10	0.59
1:A:66:LYS:HD3	1:A:67:CYS:H	1.67	0.59
1:A:474:VAL:C	1:A:477:PRO:HD2	2.27	0.59
1:A:979:SER:OG	1:A:980:ASP:N	2.35	0.59
1:A:1315:GLU:C	1:A:1317:MET:H	2.09	0.59
3:C:241:ASP:O	3:C:244:VAL:HB	2.02	0.59
5:E:79:TRP:HD1	5:E:96:PHE:HE1	1.49	0.59
1:A:866:PHE:C	1:A:867:ILE:HD12	2.27	0.59
1:A:873:MET:HG2	1:A:957:PRO:HG3	1.83	0.59
1:A:883:LEU:HD23	1:A:1021:LEU:HD13	1.84	0.59
1:A:1227:ILE:HG22	1:A:1228:TRP:N	2.16	0.59
2:B:188:ASP:O	2:B:192:LEU:HD12	2.03	0.59
2:B:446:LEU:O	2:B:447:ALA:HB3	2.03	0.59
2:B:880:THR:HB	2:B:934:LYS:HB2	1.85	0.59
11:K:16:GLU:OE1	11:K:37:LYS:HE3	2.02	0.59
1:A:310:GLY:C	1:A:312:PRO:HD2	2.26	0.59
1:A:875:ALA:HA	1:A:878:ILE:HD12	1.85	0.59
1:A:916:GLY:O	1:A:919:ILE:HG22	2.03	0.59
1:A:923:LEU:O	1:A:927:VAL:HG23	2.03	0.59
1:A:1094:VAL:HG22	1:A:1113:THR:CG2	2.32	0.59
1:A:1344:GLY:O	1:A:1345:ARG:C	2.46	0.59
2:B:596:LEU:HD11	2:B:600:LEU:HD11	1.85	0.59
2:B:734:HIS:O	2:B:735:ALA:CB	2.51	0.59
2:B:872:GLU:OE1	2:B:914:LYS:HE3	2.02	0.59
4:D:39:ASN:CG	4:D:40:HIS:N	2.60	0.59
8:H:63:LEU:HB3	8:H:90:ALA:H	1.68	0.59
11:K:7:PHE:C	11:K:9:LEU:H	2.11	0.59
11:K:53:ASP:CB	11:K:56:VAL:HB	2.33	0.59
1:A:597:LEU:O	1:A:598:LEU:HB2	2.02	0.59
2:B:936:ASP:CG	2:B:937:ALA:N	2.61	0.59
2:B:1110:PRO:HG2	2:B:1125:ASP:O	2.03	0.59
3:C:167:HIS:ND1	3:C:169:LYS:HG2	2.18	0.59
5:E:152:LYS:HG3	5:E:154:ILE:HD11	1.85	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:12:ARG:CB	2:B:1218:THR:HG22	2.33	0.59
1:A:92:HIS:HB3	1:A:95:PHE:HB2	1.85	0.59
2:B:100:PRO:HA	2:B:126:SER:HB3	1.83	0.59
2:B:115:GLN:O	2:B:119:LEU:HD12	2.02	0.59
4:D:73:SER:HB3	7:G:21:ARG:HH21	1.68	0.59
5:E:135:PHE:CD2	5:E:140:LEU:HD21	2.38	0.59
1:A:35:ILE:HG12	1:A:241:VAL:HG11	1.85	0.59
1:A:306:ASN:HD22	1:A:322:VAL:HG12	1.67	0.59
1:A:636:GLU:CD	1:A:966:ASN:HD21	2.10	0.59
1:A:760:GLN:HG2	1:A:765:VAL:HA	1.83	0.59
1:A:871:ASP:OD2	1:A:873:MET:HB2	2.03	0.59
2:B:94:LYS:HG2	2:B:95:ILE:N	2.11	0.59
3:C:34:ARG:HA	3:C:37:MET:HE2	1.84	0.59
3:C:167:HIS:O	3:C:168:ALA:C	2.45	0.59
4:D:53:SER:HB3	4:D:152:SER:CB	2.33	0.59
8:H:128:ASN:C	8:H:128:ASN:HD22	2.11	0.59
1:A:356:ASP:OD2	2:B:833:TYR:HE2	1.86	0.58
1:A:526:ASP:HB2	2:B:835:GLN:CD	2.27	0.58
1:A:577:ILE:HA	1:A:580:VAL:HG23	1.84	0.58
1:A:1343:ALA:HA	5:E:149:LEU:O	2.03	0.58
2:B:205:ILE:N	2:B:205:ILE:HD12	2.18	0.58
2:B:637:LEU:HD12	2:B:693:ILE:HD12	1.85	0.58
2:B:728:ARG:HH11	2:B:730:ARG:NH2	2.01	0.58
2:B:849:GLY:HA2	2:B:852:ARG:HD2	1.85	0.58
2:B:957:ASN:H	2:B:957:ASN:HD22	1.50	0.58
7:G:49:LEU:HG	7:G:76:ALA:HA	1.85	0.58
10:J:36:LEU:O	10:J:41:LEU:HG	2.02	0.58
2:B:815:ARG:HB2	2:B:815:ARG:HH11	1.68	0.58
2:B:1034:VAL:HA	2:B:1037:LEU:CD1	2.32	0.58
2:B:1159:ARG:HB3	2:B:1159:ARG:NH1	2.15	0.58
3:C:112:ASN:HD21	3:C:146:LYS:HE2	1.67	0.58
6:F:100:GLN:NE2	7:G:61:ILE:HD11	2.17	0.58
7:G:9:LEU:HD12	7:G:10:ASN:N	2.19	0.58
1:A:1226:VAL:HG13	1:A:1239:ARG:O	2.03	0.58
1:A:1394:THR:HG22	1:A:1395:GLY:H	1.68	0.58
2:B:119:LEU:HD23	2:B:953:LEU:HD11	1.85	0.58
2:B:308:TRP:CH2	9:I:45:ARG:HG2	2.37	0.58
2:B:681:TRP:CE3	2:B:684:LEU:HD12	2.38	0.58
2:B:901:PRO:HG3	2:B:952:VAL:HG23	1.84	0.58
2:B:1174:LYS:HB2	2:B:1179:GLN:O	2.02	0.58
4:D:147:TYR:CZ	7:G:103:VAL:HG13	2.38	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:392:VAL:HG13	1:A:415:LEU:HD21	1.85	0.58
1:A:849:MET:HE1	1:A:1061:GLY:HA2	1.86	0.58
1:A:1316:VAL:O	1:A:1316:VAL:HG12	2.03	0.58
2:B:69:LEU:HD13	2:B:429:PHE:CD1	2.37	0.58
2:B:663:ALA:O	2:B:667:GLN:HG3	2.04	0.58
3:C:242:GLN:HE21	3:C:246:ARG:NE	1.99	0.58
11:K:70:ARG:O	11:K:70:ARG:HG3	2.02	0.58
1:A:148:CYS:O	1:A:168:GLY:HA2	2.03	0.58
2:B:234:ILE:HG21	2:B:237:VAL:HG23	1.84	0.58
2:B:659:ALA:HA	2:B:662:MET:HE2	1.84	0.58
2:B:912:ILE:HD11	2:B:966:VAL:CG2	2.34	0.58
7:G:51:TYR:O	7:G:51:TYR:HD2	1.86	0.58
8:H:59:ILE:HG22	8:H:60:ALA:N	2.18	0.58
1:A:21:LEU:HD11	1:A:1414:ALA:HA	1.85	0.58
1:A:449:SER:HB2	2:B:1133:MET:HB3	1.85	0.58
1:A:515:GLN:HA	1:A:1367:HIS:NE2	2.19	0.58
1:A:595:THR:O	1:A:596:THR:HG23	2.02	0.58
1:A:973:ILE:HD13	1:A:1038:THR:HG23	1.84	0.58
2:B:168:GLY:N	2:B:450:ALA:HB1	2.18	0.58
5:E:169:ARG:HD3	6:F:140:ASP:OD1	2.04	0.58
1:A:71:GLN:O	1:A:73:GLY:N	2.35	0.58
1:A:391:LEU:O	1:A:394:ASN:HB2	2.03	0.58
1:A:1227:ILE:HG22	1:A:1228:TRP:H	1.67	0.58
3:C:18:VAL:HG21	11:K:109:TRP:CZ3	2.37	0.58
10:J:37:SER:OG	10:J:47:ARG:NH2	2.36	0.58
1:A:1349:TYR:C	1:A:1349:TYR:CD2	2.82	0.58
3:C:8:VAL:HG12	3:C:9:LYS:N	2.18	0.58
4:D:39:ASN:ND2	4:D:40:HIS:H	2.01	0.58
5:E:213:ILE:HG12	5:E:214:CYS:H	1.68	0.58
6:F:77:ASP:O	6:F:78:GLN:HB3	2.02	0.58
10:J:2:ILE:HG23	10:J:3:VAL:O	2.04	0.58
1:A:172:PRO:HG3	1:A:185:TRP:CE2	2.39	0.58
1:A:664:THR:CG2	2:B:1014:PRO:HB3	2.34	0.58
1:A:850:VAL:HG21	1:A:1058:VAL:HG11	1.86	0.58
1:A:1118:VAL:CG2	1:A:1306:LEU:HB2	2.33	0.58
2:B:108:VAL:HG23	2:B:109:THR:H	1.68	0.58
4:D:52:LEU:HB2	4:D:182:SER:HB2	1.84	0.58
5:E:187:TYR:HD2	5:E:188:LEU:HD23	1.69	0.58
8:H:142:LEU:C	8:H:143:LEU:HD12	2.29	0.58
1:A:178:GLY:C	1:A:179:LEU:HD23	2.29	0.58
1:A:827:THR:HG22	1:A:828:ALA:N	2.19	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:858:ASN:HD22	1:A:858:ASN:H	1.52	0.58
1:A:1150:SER:H	9:I:46:HIS:HB3	1.68	0.58
1:A:1187:GLN:HG3	1:A:1188:GLN:H	1.69	0.58
1:A:1350:LYS:O	1:A:1353:TYR:HB3	2.03	0.58
2:B:326:ASP:O	2:B:330:ALA:CB	2.50	0.58
2:B:453:ILE:O	2:B:457:LEU:HG	2.03	0.58
5:E:197:LYS:HE2	5:E:199:ILE:HD11	1.86	0.58
6:F:87:LYS:CE	6:F:88:TYR:HE1	2.16	0.58
1:A:617:VAL:HG12	1:A:622:VAL:HG12	1.85	0.57
1:A:1336:MET:HE2	1:A:1381:LEU:HG	1.85	0.57
1:A:1349:TYR:CE2	1:A:1353:TYR:HB2	2.39	0.57
2:B:975:GLN:CG	2:B:976:ILE:N	2.67	0.57
3:C:45:ALA:CB	3:C:72:LEU:HD13	2.33	0.57
5:E:28:TYR:C	5:E:65:THR:HG23	2.29	0.57
5:E:155:ARG:C	5:E:156:LEU:HD23	2.29	0.57
6:F:130:ILE:O	6:F:132:LEU:N	2.36	0.57
8:H:17:PRO:HB3	8:H:24:CYS:SG	2.43	0.57
1:A:224:PHE:CD2	1:A:231:PRO:HG3	2.38	0.57
1:A:524:VAL:HG12	1:A:525:GLN:N	2.15	0.57
1:A:808:LEU:HB3	1:A:812:GLU:HB2	1.86	0.57
1:A:896:ARG:HD3	1:A:897:TYR:CE1	2.39	0.57
1:A:1406:VAL:HG12	1:A:1410:PHE:CE1	2.39	0.57
2:B:879:ARG:H	2:B:879:ARG:NE	2.02	0.57
2:B:1166:CYS:O	2:B:1168:LEU:N	2.36	0.57
3:C:191:TYR:HD2	3:C:201:TRP:CD1	2.22	0.57
10:J:57:ILE:HG23	10:J:58:GLU:H	1.69	0.57
1:A:6:TYR:CG	1:A:7:SER:N	2.72	0.57
1:A:7:SER:OG	2:B:1161:HIS:CE1	2.58	0.57
1:A:98:LYS:O	1:A:102:VAL:HG23	2.04	0.57
1:A:206:GLU:O	1:A:210:ILE:HG13	2.04	0.57
1:A:1166:ASP:OD2	1:A:1239:ARG:HD2	2.05	0.57
2:B:898:LEU:HD12	12:L:58:LYS:HD2	1.86	0.57
6:F:90:ARG:CD	6:F:155:LEU:HD13	2.32	0.57
1:A:247:ARG:HD2	1:A:263:THR:HG23	1.85	0.57
1:A:447:GLN:HG2	2:B:1134:GLU:OE2	2.04	0.57
1:A:546:VAL:O	1:A:549:MET:HB2	2.05	0.57
1:A:834:THR:HG21	1:A:1077:THR:HG23	1.86	0.57
1:A:975:HIS:O	1:A:976:THR:HG23	2.05	0.57
1:A:1376:THR:O	1:A:1377:THR:C	2.48	0.57
1:A:1444:MET:HG3	7:G:59:GLY:O	2.04	0.57
2:B:365:THR:HG23	2:B:367:LEU:H	1.69	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:655:LYS:HD2	2:B:658:ILE:HG22	1.84	0.57
2:B:777:ALA:HA	2:B:1095:LEU:HB2	1.87	0.57
3:C:47:ASP:HA	12:L:69:ALA:HB3	1.86	0.57
3:C:50:GLU:HB3	3:C:156:THR:HB	1.86	0.57
12:L:58:LYS:O	12:L:59:ALA:O	2.22	0.57
1:A:443:LEU:HD23	1:A:501:LEU:HD22	1.87	0.57
2:B:597:MET:HA	2:B:597:MET:CE	2.32	0.57
6:F:76:LYS:HA	6:F:79:ARG:HD2	1.85	0.57
7:G:115:MET:CB	7:G:119:LEU:HD23	2.35	0.57
10:J:3:VAL:HG21	10:J:18:TRP:CG	2.39	0.57
1:A:409:SER:O	1:A:411:ASP:HB2	2.05	0.57
1:A:640:GLN:CD	1:A:640:GLN:H	2.11	0.57
1:A:1254:ALA:O	1:A:1255:GLU:HB2	2.04	0.57
1:A:1279:ILE:HG23	1:A:1308:THR:OG1	2.04	0.57
2:B:226:PHE:HA	2:B:395:GLN:HG3	1.87	0.57
2:B:345:LYS:HG2	2:B:346:GLU:N	2.17	0.57
2:B:405:ARG:NH1	2:B:632:ARG:HG2	2.19	0.57
3:C:16:ASP:O	3:C:233:GLU:HA	2.04	0.57
6:F:101:ILE:HD13	6:F:120:ILE:CG2	2.29	0.57
10:J:9:SER:CB	10:J:45:CYS:HB2	2.35	0.57
1:A:270:LEU:O	1:A:274:ILE:HG13	2.05	0.57
1:A:355:GLY:N	1:A:482:PHE:CZ	2.72	0.57
1:A:1354:ASN:O	1:A:1358:SER:CB	2.51	0.57
2:B:467:GLY:H	2:B:475:SER:HB2	1.69	0.57
2:B:593:PRO:HB2	2:B:617:ARG:CZ	2.35	0.57
2:B:977:GLY:HA3	2:B:1099:VAL:HG11	1.86	0.57
9:I:15:TYR:N	9:I:15:TYR:CD1	2.73	0.57
10:J:3:VAL:HG21	10:J:18:TRP:CB	2.34	0.57
1:A:850:VAL:HG23	1:A:1064:VAL:HG21	1.87	0.57
1:A:1315:GLU:O	1:A:1317:MET:N	2.37	0.57
2:B:100:PRO:HD2	2:B:180:TYR:CE1	2.40	0.57
2:B:990:ILE:HG22	2:B:992:ILE:H	1.69	0.57
5:E:15:ALA:O	5:E:19:VAL:HG23	2.04	0.57
5:E:185:ALA:O	5:E:190:LEU:HG	2.05	0.57
7:G:111:THR:CG2	7:G:114:LEU:HD13	2.34	0.57
12:L:26:THR:HG22	12:L:27:LEU:N	2.15	0.57
1:A:202:LEU:HB3	1:A:207:ILE:HD11	1.87	0.57
1:A:1356:ILE:HD13	1:A:1363:VAL:HG21	1.86	0.57
2:B:361:LEU:HD21	2:B:377:PHE:HD2	1.69	0.57
2:B:398:ARG:HB2	2:B:398:ARG:NH1	2.20	0.57
6:F:129:LYS:O	6:F:130:ILE:HG13	2.05	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:G:126:ASN:HD22	7:G:127:PRO:N	2.03	0.57
11:K:39:ASP:OD1	11:K:41:THR:HG22	2.05	0.57
1:A:63:ARG:HG2	1:A:74:MET:HE1	1.86	0.57
1:A:403:LYS:HE2	1:A:404:TYR:HE1	1.70	0.57
1:A:438:ASP:O	1:A:439:ASN:C	2.47	0.57
1:A:692:ASP:C	1:A:694:THR:H	2.13	0.57
1:A:933:TYR:O	1:A:937:VAL:HG23	2.04	0.57
1:A:1029:ARG:O	1:A:1033:GLN:HB3	2.04	0.57
2:B:192:LEU:O	2:B:193:LYS:HB2	2.05	0.57
2:B:282:ILE:HD12	2:B:382:ILE:HD13	1.86	0.57
2:B:797:TYR:HB2	2:B:852:ARG:O	2.04	0.57
2:B:1098:MET:O	2:B:1101:ASP:HB2	2.05	0.57
4:D:170:THR:O	4:D:172:LEU:N	2.38	0.57
7:G:92:VAL:HG21	7:G:102:GLN:HB2	1.86	0.57
8:H:89:LEU:C	8:H:91:ASP:H	2.13	0.57
2:B:363:HIS:O	2:B:364:ILE:CB	2.52	0.56
2:B:615:MET:HE3	2:B:615:MET:N	2.17	0.56
2:B:724:ASP:HB3	2:B:727:LYS:HG3	1.86	0.56
2:B:995:ARG:HB3	2:B:997:GLU:OE2	2.05	0.56
11:K:57:LEU:HD13	11:K:76:GLN:NE2	2.18	0.56
1:A:92:HIS:HE1	1:A:1410:PHE:CE2	2.22	0.56
1:A:1351:GLU:O	1:A:1355:VAL:HG23	2.05	0.56
2:B:100:PRO:HG3	2:B:172:ILE:HD12	1.86	0.56
2:B:1135:ARG:O	2:B:1139:ILE:HG13	2.05	0.56
5:E:153:HIS:HB3	5:E:196:VAL:HG21	1.86	0.56
1:A:12:ARG:HB2	2:B:1218:THR:HG22	1.86	0.56
1:A:134:ARG:HD2	1:A:221:SER:O	2.05	0.56
1:A:148:CYS:HB3	1:A:168:GLY:HA2	1.87	0.56
1:A:490:HIS:ND1	2:B:1150:ARG:NH1	2.53	0.56
1:A:903:ASN:HB3	1:A:906:HIS:HB2	1.87	0.56
1:A:1116:LEU:N	1:A:1308:THR:HG22	2.20	0.56
2:B:31:TRP:CE3	2:B:34:ILE:HD12	2.40	0.56
2:B:230:ALA:CB	2:B:231:PRO:HD3	2.33	0.56
2:B:510:LYS:CG	2:B:511:PRO:HD3	2.24	0.56
2:B:792:MET:HA	2:B:856:PHE:O	2.05	0.56
2:B:1110:PRO:O	2:B:1119:VAL:HG13	2.05	0.56
2:B:1157:ALA:O	2:B:1158:PHE:CB	2.52	0.56
5:E:155:ARG:O	5:E:156:LEU:HD23	2.05	0.56
8:H:30:SER:CB	8:H:36:CYS:HB3	2.35	0.56
11:K:53:ASP:C	11:K:55:LYS:H	2.11	0.56
1:A:92:HIS:HE1	1:A:1410:PHE:HE2	1.51	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:185:TRP:CZ3	1:A:200:ARG:HB3	2.41	0.56
1:A:549:MET:CE	1:A:656:TRP:HD1	2.18	0.56
1:A:810:PRO:HD3	2:B:730:ARG:HH21	1.70	0.56
2:B:486:TYR:HD1	2:B:1096:ARG:NH2	2.03	0.56
2:B:759:PRO:C	2:B:761:HIS:H	2.13	0.56
3:C:35:ARG:NH1	11:K:41:THR:HB	2.20	0.56
4:D:24:ALA:HB3	4:D:26:THR:HG23	1.88	0.56
5:E:99:HIS:CE1	5:E:103:LYS:HG3	2.41	0.56
6:F:97:ARG:NH2	6:F:108:PHE:HE1	2.03	0.56
6:F:140:ASP:CG	6:F:141:GLY:H	2.13	0.56
7:G:111:THR:HG22	7:G:114:LEU:CD2	2.34	0.56
1:A:81:PHE:CD2	1:A:243:PRO:HD3	2.41	0.56
1:A:108:MET:O	1:A:109:HIS:HB3	2.04	0.56
1:A:363:GLN:HA	1:A:459:ARG:O	2.05	0.56
1:A:850:VAL:HG12	1:A:1060:PRO:HA	1.87	0.56
1:A:1260:LEU:HA	1:A:1263:ILE:HD12	1.86	0.56
2:B:65:GLU:HG3	2:B:66:ASP:H	1.71	0.56
2:B:792:MET:HG3	2:B:855:PHE:HE1	1.70	0.56
2:B:900:ALA:HB1	12:L:61:THR:OG1	2.05	0.56
5:E:55:ARG:C	5:E:57:MET:N	2.61	0.56
6:F:130:ILE:HG22	6:F:132:LEU:HG	1.88	0.56
1:A:11:LEU:HG	1:A:12:ARG:N	2.21	0.56
1:A:361:LEU:HA	1:A:471:ASN:ND2	2.11	0.56
1:A:370:ILE:C	1:A:372:LYS:H	2.13	0.56
1:A:1362:TYR:HD1	1:A:1363:VAL:N	2.04	0.56
2:B:798:TYR:CD1	10:J:4:PRO:HB3	2.41	0.56
2:B:1065:GLN:HB2	2:B:1069:PHE:O	2.06	0.56
6:F:97:ARG:NH2	6:F:108:PHE:CE1	2.74	0.56
1:A:528:LEU:O	1:A:531:ILE:HG22	2.06	0.56
1:A:547:LEU:HD13	11:K:58:PHE:CD1	2.41	0.56
1:A:617:VAL:HG12	1:A:622:VAL:CG1	2.35	0.56
1:A:868:TYR:HE2	1:A:1366:ARG:HE	1.53	0.56
1:A:1011:GLN:HE22	1:A:1015:VAL:HG21	1.69	0.56
1:A:1152:ILE:CD1	9:I:44:TYR:HD2	2.18	0.56
1:A:1210:GLY:O	1:A:1214:GLU:HG2	2.05	0.56
2:B:515:HIS:CD2	2:B:516:ASN:H	2.24	0.56
2:B:533:CYS:C	2:B:535:LEU:H	2.12	0.56
3:C:31:ASN:O	3:C:35:ARG:HG3	2.05	0.56
3:C:32:SER:O	3:C:36:VAL:HG23	2.05	0.56
4:D:123:LEU:HB3	4:D:124:GLU:OE2	2.06	0.56
4:D:151:PHE:CD1	4:D:151:PHE:N	2.74	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:216:ASN:C	4:D:218:GLU:H	2.13	0.56
8:H:11:GLN:HA	8:H:53:ASP:O	2.06	0.56
1:A:477:PRO:CG	1:A:521:MET:HE2	2.36	0.56
1:A:565:ILE:HG23	1:A:567:LYS:HG2	1.86	0.56
1:A:1110:ASN:H	1:A:1110:ASN:ND2	2.04	0.56
2:B:1079:LYS:HZ2	3:C:188:HIS:CE1	2.23	0.56
3:C:148:ARG:N	3:C:151:GLN:HG3	2.19	0.56
6:F:100:GLN:NE2	7:G:18:PHE:CE2	2.73	0.56
7:G:51:TYR:O	7:G:54:ILE:HG13	2.05	0.56
1:A:664:THR:HG22	2:B:1014:PRO:HB3	1.87	0.56
2:B:120:ARG:HH12	12:L:54:ARG:HH11	1.54	0.56
2:B:220:GLY:HA2	2:B:241:ARG:NH1	2.21	0.56
2:B:334:ILE:HG22	2:B:334:ILE:O	2.05	0.56
3:C:101:LEU:HB3	3:C:155:LEU:HB2	1.86	0.56
4:D:35:LEU:H	4:D:35:LEU:CD1	2.18	0.56
5:E:90:VAL:O	5:E:90:VAL:HG22	2.06	0.56
7:G:110:VAL:HG12	7:G:115:MET:HE2	1.88	0.56
1:A:1394:THR:HG21	1:A:1398:MET:SD	2.46	0.56
1:A:1396:ALA:HA	1:A:1399:ARG:NH2	2.21	0.56
2:B:222:ILE:HD11	2:B:627:PHE:CZ	2.40	0.56
2:B:294:ASP:C	2:B:296:GLU:H	2.14	0.56
7:G:13:LEU:CD2	7:G:17:PHE:HD2	2.12	0.56
1:A:332:LYS:C	1:A:334:GLY:H	2.15	0.55
1:A:541:ILE:HG22	1:A:542:GLU:N	2.21	0.55
1:A:1324:PRO:HB2	5:E:142:VAL:HG11	1.88	0.55
2:B:362:PRO:HB3	2:B:366:GLN:OE1	2.06	0.55
2:B:485:ARG:HH12	2:B:782:LEU:HD21	1.71	0.55
2:B:542:MET:HG2	2:B:747:MET:HB3	1.88	0.55
2:B:828:ALA:HB2	2:B:1085:ILE:HG12	1.88	0.55
3:C:114:TYR:CD2	3:C:140:ASN:HB3	2.41	0.55
4:D:56:ARG:HG3	4:D:145:MET:HE1	1.87	0.55
5:E:148:GLU:HG2	5:E:149:LEU:HD23	1.87	0.55
12:L:34:CYS:O	12:L:35:SER:C	2.48	0.55
2:B:679:TYR:CE1	2:B:683:SER:HB2	2.41	0.55
3:C:43:THR:HG22	3:C:74:SER:OG	2.06	0.55
4:D:208:GLU:O	4:D:212:LYS:HG3	2.06	0.55
8:H:82:PRO:O	8:H:84:ALA:N	2.40	0.55
1:A:145:LYS:HE3	1:A:146:MET:H	1.71	0.55
1:A:172:PRO:HB3	1:A:185:TRP:CD2	2.40	0.55
1:A:341:MET:HE2	1:A:843:LYS:HZ1	1.70	0.55
1:A:718:VAL:O	1:A:721:PHE:HB2	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1191:TRP:HZ3	9:I:43:VAL:HG21	1.72	0.55
2:B:998:ASP:OD1	3:C:35:ARG:NH2	2.40	0.55
3:C:123:ASN:OD1	3:C:125:MET:HG2	2.07	0.55
4:D:156:ASP:C	4:D:158:GLU:H	2.14	0.55
5:E:63:ASN:HB3	5:E:64:PRO:HD2	1.88	0.55
8:H:130:ARG:HB3	8:H:134:ASN:H	1.72	0.55
1:A:1011:GLN:HG3	1:A:1012:ARG:N	2.20	0.55
2:B:522:VAL:HG13	2:B:537:LYS:HB3	1.88	0.55
2:B:805:THR:HG22	2:B:806:THR:O	2.06	0.55
3:C:113:VAL:O	3:C:143:LEU:HD13	2.07	0.55
4:D:134:THR:HG21	4:D:138:ASN:HB2	1.88	0.55
5:E:114:ASN:O	5:E:115:ASN:HB3	2.06	0.55
7:G:160:ILE:HG22	7:G:161:GLY:N	2.21	0.55
11:K:32:VAL:HA	11:K:73:LEU:O	2.06	0.55
1:A:16:GLU:HA	1:A:1419:ASP:O	2.06	0.55
1:A:335:ARG:HD3	1:A:339:ASN:ND2	2.19	0.55
1:A:548:ASN:HD21	11:K:47:ARG:HH21	1.53	0.55
1:A:1291:VAL:HG13	1:A:1292:PRO:CD	2.36	0.55
2:B:120:ARG:NH1	12:L:54:ARG:HH11	2.02	0.55
2:B:390:LEU:O	2:B:391:ASP:C	2.49	0.55
2:B:542:MET:HE2	2:B:747:MET:HG3	1.87	0.55
11:K:83:PRO:O	11:K:84:LYS:C	2.49	0.55
1:A:308:ILE:HG22	1:A:309:ALA:H	1.71	0.55
1:A:353:ILE:HG23	1:A:485:ASP:O	2.07	0.55
1:A:481:ASP:HB2	1:A:483:ASP:OD2	2.07	0.55
1:A:666:ILE:HD11	2:B:1067:ARG:O	2.06	0.55
1:A:670:ILE:CG2	1:A:805:LEU:HD21	2.37	0.55
1:A:679:ILE:O	1:A:683:ILE:HG13	2.06	0.55
2:B:213:ILE:HD11	2:B:481:GLN:CD	2.32	0.55
2:B:416:LEU:O	2:B:420:LEU:HB2	2.07	0.55
2:B:467:GLY:H	2:B:475:SER:CB	2.19	0.55
2:B:785:TYR:CE2	10:J:60:PHE:CE1	2.95	0.55
2:B:876:LYS:NZ	2:B:891:ASP:HA	2.22	0.55
3:C:213:PRO:O	3:C:214:ASN:HB3	2.04	0.55
4:D:39:ASN:HD22	4:D:41:GLN:H	1.53	0.55
5:E:89:GLY:C	5:E:91:LYS:H	2.13	0.55
11:K:37:LYS:O	11:K:38:GLU:HG2	2.05	0.55
11:K:68:PHE:N	11:K:68:PHE:CD2	2.71	0.55
1:A:593:GLU:C	1:A:595:THR:H	2.15	0.55
1:A:711:ARG:O	1:A:714:PHE:HB3	2.06	0.55
1:A:840:ARG:HG2	1:A:1402:PHE:HZ	1.72	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1349:TYR:C	1:A:1349:TYR:HD2	2.14	0.55
2:B:806:THR:HB	2:B:809:MET:HG3	1.87	0.55
2:B:842:ASN:HD21	2:B:844:SER:CB	2.19	0.55
4:D:138:ASN:HD21	7:G:35:GLU:HB3	1.70	0.55
5:E:143:ASN:ND2	5:E:146:HIS:ND1	2.55	0.55
1:A:95:PHE:CE2	1:A:1410:PHE:HD2	2.25	0.55
1:A:853:ASP:OD1	1:A:855:THR:CG2	2.53	0.55
2:B:1174:LYS:HB2	2:B:1179:GLN:H	1.72	0.55
11:K:22:ASP:C	11:K:31:VAL:HG13	2.31	0.55
1:A:556:TRP:O	1:A:558:GLY:N	2.40	0.55
1:A:633:VAL:O	1:A:634:THR:C	2.48	0.55
1:A:666:ILE:HG23	2:B:1026:LEU:CB	2.34	0.55
1:A:1004:ASN:O	1:A:1008:GLN:HG2	2.06	0.55
1:A:1095:THR:HG21	1:A:1112:LYS:HB2	1.89	0.55
1:A:1105:LEU:O	1:A:1384:VAL:HG23	2.07	0.55
1:A:1287:TYR:O	1:A:1302:PRO:HA	2.07	0.55
2:B:185:THR:O	2:B:189:LEU:HG	2.07	0.55
2:B:230:ALA:HB3	2:B:231:PRO:CD	2.32	0.55
2:B:251:ILE:O	2:B:251:ILE:HG22	2.07	0.55
2:B:999:MET:HA	2:B:999:MET:CE	2.35	0.55
2:B:1012:ILE:O	2:B:1014:PRO:HD3	2.07	0.55
3:C:82:TYR:HB2	3:C:85:ASP:OD2	2.06	0.55
1:A:82:GLY:C	1:A:241:VAL:HB	2.32	0.55
1:A:224:PHE:HD2	1:A:231:PRO:CD	2.19	0.55
1:A:909:ASP:O	1:A:911:SER:N	2.40	0.55
3:C:22:LEU:HD22	3:C:230:MET:HE1	1.89	0.55
4:D:13:ARG:C	4:D:15:LEU:H	2.15	0.55
4:D:25:ALA:HB3	4:D:196:PRO:HG2	1.88	0.55
2:B:57:TYR:CD1	2:B:57:TYR:N	2.72	0.54
2:B:179:CYS:SG	2:B:181:LEU:HD12	2.47	0.54
2:B:754:SER:HB2	2:B:812:LEU:HD11	1.89	0.54
2:B:1137:CYS:O	2:B:1140:ALA:HB3	2.07	0.54
4:D:30:GLY:C	4:D:32:GLU:N	2.64	0.54
8:H:130:ARG:H	8:H:130:ARG:NH1	1.96	0.54
10:J:17:LYS:HB3	10:J:39:LEU:CD2	2.38	0.54
1:A:441:PRO:HD2	1:A:498:ARG:CZ	2.37	0.54
1:A:541:ILE:HD12	1:A:577:ILE:HD11	1.87	0.54
1:A:1050:GLU:O	1:A:1053:PHE:HB3	2.07	0.54
1:A:1400:CYS:SG	1:A:1409:LEU:HD21	2.47	0.54
2:B:227:LYS:H	2:B:395:GLN:CD	2.16	0.54
2:B:859:TYR:CD1	2:B:859:TYR:N	2.71	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:88:CYS:SG	3:C:91:HIS:CA	2.93	0.54
4:D:118:THR:HG21	4:D:121:LYS:HD2	1.88	0.54
7:G:159:ALA:C	7:G:160:ILE:HD12	2.32	0.54
11:K:51:LEU:CD1	11:K:59:ALA:HB3	2.37	0.54
11:K:113:THR:CG2	11:K:114:LEU:H	2.20	0.54
1:A:364:VAL:O	1:A:366:VAL:HG23	2.07	0.54
1:A:986:ILE:HG22	1:A:987:VAL:N	2.22	0.54
1:A:1103:GLU:C	1:A:1108:ALA:HB2	2.32	0.54
1:A:1224:LEU:N	1:A:1243:VAL:HG13	2.22	0.54
1:A:1444:MET:HE1	6:F:135:ARG:HB2	1.89	0.54
2:B:856:PHE:CD1	2:B:856:PHE:N	2.74	0.54
2:B:955:THR:HG22	2:B:956:THR:O	2.07	0.54
5:E:177:ARG:HH11	5:E:215:MET:CE	2.20	0.54
5:E:178:ILE:HG22	5:E:213:ILE:O	2.07	0.54
8:H:102:TYR:N	8:H:102:TYR:CD2	2.69	0.54
10:J:30:LEU:HD22	10:J:34:THR:HG21	1.89	0.54
10:J:48:ARG:HH11	10:J:48:ARG:CG	2.20	0.54
1:A:374:LEU:O	1:A:436:ILE:HG12	2.07	0.54
1:A:665:GLY:O	1:A:667:GLY:N	2.40	0.54
1:A:830:LYS:HG3	1:A:1098:VAL:HG11	1.89	0.54
1:A:982:THR:O	1:A:985:ASP:HB2	2.07	0.54
1:A:1313:LEU:O	1:A:1315:GLU:N	2.41	0.54
1:A:1390:ASN:HD21	1:A:1402:PHE:HB3	1.72	0.54
2:B:827:ILE:O	2:B:1085:ILE:HG23	2.07	0.54
2:B:955:THR:HG23	12:L:54:ARG:O	2.08	0.54
2:B:1152:MET:O	2:B:1157:ALA:HB2	2.07	0.54
2:B:1168:LEU:HB2	2:B:1170:THR:OG1	2.07	0.54
4:D:56:ARG:HD2	4:D:122:GLU:OE1	2.08	0.54
5:E:13:TRP:CZ3	5:E:39:LEU:HB2	2.42	0.54
10:J:7:CYS:HB2	10:J:49:MET:HE3	1.90	0.54
1:A:164:ARG:HG2	1:A:165:GLY:N	2.23	0.54
1:A:525:GLN:O	1:A:526:ASP:C	2.50	0.54
1:A:707:GLY:HA2	1:A:1281:ARG:HD3	1.90	0.54
2:B:515:HIS:H	2:B:518:HIS:CD2	2.25	0.54
2:B:516:ASN:H	2:B:516:ASN:HD22	1.55	0.54
1:A:102:VAL:CG1	1:A:211:PHE:HE1	2.20	0.54
1:A:135:PHE:HD1	1:A:222:LEU:HD22	1.71	0.54
1:A:741:ASN:O	1:A:745:GLN:HG3	2.08	0.54
2:B:210:LYS:CD	2:B:482:VAL:HG13	2.28	0.54
2:B:521:LEU:HD22	2:B:633:VAL:CG1	2.35	0.54
2:B:911:ILE:HG21	2:B:966:VAL:HG11	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:H:89:LEU:O	8:H:89:LEU:HD12	2.07	0.54
12:L:55:ILE:O	12:L:56:LEU:HB2	2.08	0.54
1:A:826:ASP:HB2	1:A:830:LYS:HD3	1.90	0.54
2:B:458:LYS:O	2:B:459:TYR:C	2.51	0.54
2:B:581:PHE:O	2:B:582:VAL:HG23	2.08	0.54
4:D:192:LYS:HD2	4:D:199:ASN:H	1.72	0.54
7:G:79:PHE:CD2	7:G:105:PRO:HD2	2.42	0.54
1:A:16:GLU:OE1	4:D:13:ARG:NH2	2.41	0.54
1:A:528:LEU:C	1:A:528:LEU:HD12	2.33	0.54
1:A:1427:ASN:O	1:A:1430:LEU:N	2.40	0.54
3:C:185:LYS:C	3:C:186:LEU:HD12	2.33	0.54
1:A:224:PHE:CD2	1:A:231:PRO:HD3	2.42	0.54
1:A:1173:HIS:CD2	1:A:1227:ILE:HG23	2.43	0.54
2:B:126:SER:HB3	2:B:172:ILE:HD11	1.89	0.54
2:B:213:ILE:HG21	2:B:499:ASN:HB2	1.88	0.54
4:D:190:GLU:HG2	4:D:194:LEU:HD11	1.88	0.54
6:F:121:ALA:HA	6:F:124:GLU:HB2	1.90	0.54
12:L:45:ALA:C	12:L:46:VAL:HG22	2.33	0.54
1:A:500:GLU:O	1:A:504:LEU:HB2	2.08	0.54
1:A:669:THR:O	1:A:762:SER:HB3	2.07	0.54
1:A:741:ASN:ND2	1:A:743:VAL:H	2.07	0.54
1:A:998:LEU:HD23	1:A:1001:ARG:HG3	1.88	0.54
2:B:291:ILE:H	2:B:291:ILE:HD12	1.72	0.54
2:B:611:PRO:O	2:B:692:TYR:HB2	2.08	0.54
2:B:635:ARG:HH22	2:B:742:GLU:CD	2.16	0.54
2:B:662:MET:HA	2:B:665:GLU:HB2	1.88	0.54
2:B:899:ILE:HG23	2:B:903:VAL:HG21	1.89	0.54
3:C:111:THR:O	3:C:147:LEU:HD23	2.08	0.54
6:F:88:TYR:O	6:F:89:GLU:C	2.50	0.54
8:H:62:SER:C	8:H:63:LEU:HD12	2.33	0.54
1:A:196:GLU:HG2	1:A:197:PRO:HD2	1.90	0.53
1:A:404:TYR:CD2	1:A:412:ARG:HD3	2.43	0.53
1:A:425:GLN:H	1:A:425:GLN:CD	2.16	0.53
1:A:427:GLN:HG3	1:A:430:TRP:CE2	2.42	0.53
1:A:540:PHE:HZ	8:H:121:LEU:HD11	1.73	0.53
1:A:567:LYS:HE3	8:H:46:LEU:CD1	2.38	0.53
1:A:908:LEU:HD23	1:A:1029:ARG:NH2	2.23	0.53
1:A:1098:VAL:HB	1:A:1099:PRO:CD	2.38	0.53
1:A:1119:TYR:HA	1:A:1305:VAL:CG2	2.38	0.53
1:A:1362:TYR:CD1	1:A:1362:TYR:C	2.85	0.53
2:B:37:PHE:HE1	2:B:41:LYS:HG3	1.73	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:828:ALA:CB	2:B:1085:ILE:HG12	2.37	0.53
4:D:138:ASN:O	4:D:139:LYS:C	2.51	0.53
1:A:63:ARG:HA	1:A:74:MET:HE2	1.90	0.53
1:A:964:ILE:O	1:A:964:ILE:HG22	2.08	0.53
2:B:508:LEU:O	2:B:509:ALA:CB	2.57	0.53
2:B:917:PRO:O	2:B:918:ILE:HG13	2.07	0.53
2:B:1198:TYR:O	2:B:1201:LYS:HB3	2.09	0.53
4:D:25:ALA:C	4:D:27:LEU:H	2.15	0.53
4:D:39:ASN:ND2	4:D:41:GLN:H	2.05	0.53
4:D:167:LEU:O	4:D:169:SER:N	2.41	0.53
6:F:118:LEU:O	6:F:122:MET:HG3	2.07	0.53
11:K:30:ALA:HA	11:K:75:ILE:O	2.08	0.53
1:A:49:LYS:HZ1	1:A:61:ILE:H	1.56	0.53
1:A:605:MET:SD	1:A:621:THR:HG21	2.49	0.53
2:B:172:ILE:HG21	2:B:178:ASN:HB3	1.90	0.53
2:B:188:ASP:C	2:B:192:LEU:HD12	2.33	0.53
2:B:654:ARG:O	2:B:657:HIS:HB2	2.08	0.53
2:B:744:HIS:CD2	2:B:746:SER:OG	2.60	0.53
3:C:62:PHE:O	3:C:66:ARG:CG	2.57	0.53
4:D:170:THR:O	4:D:172:LEU:HG	2.08	0.53
7:G:17:PHE:C	7:G:19:GLY:H	2.16	0.53
8:H:114:VAL:HG22	8:H:125:LEU:HB3	1.91	0.53
9:I:6:PHE:HB3	9:I:12:ASN:O	2.08	0.53
1:A:236:LEU:N	1:A:236:LEU:HD23	2.23	0.53
1:A:317:LYS:O	1:A:318:SER:CB	2.55	0.53
1:A:360:GLU:O	1:A:361:LEU:C	2.51	0.53
1:A:381:THR:CG2	1:A:382:PRO:HD2	2.38	0.53
1:A:635:ARG:HA	1:A:635:ARG:HH11	1.72	0.53
2:B:284:ILE:HD13	2:B:333:PHE:CD2	2.44	0.53
2:B:323:VAL:O	2:B:324:ILE:HG13	2.08	0.53
2:B:360:PHE:O	2:B:361:LEU:C	2.52	0.53
2:B:763:GLN:O	2:B:766:ARG:HB2	2.07	0.53
2:B:878:GLN:H	2:B:934:LYS:HZ2	1.55	0.53
3:C:61:GLU:O	3:C:62:PHE:C	2.50	0.53
7:G:30:LEU:HD23	7:G:31:LEU:HD23	1.91	0.53
9:I:7:CYS:SG	9:I:8:ARG:O	2.66	0.53
10:J:24:LEU:HD11	10:J:38:ARG:HG2	1.89	0.53
1:A:763:ALA:C	1:A:803:SER:HB3	2.33	0.53
1:A:954:TRP:HZ3	5:E:203:GLU:HB2	1.68	0.53
2:B:522:VAL:CG1	2:B:523:CYS:N	2.70	0.53
2:B:583:ASN:OD1	2:B:628:THR:HG22	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:F:83:PRO:HA	6:F:146:TRP:CZ3	2.43	0.53
8:H:113:ALA:HA	8:H:125:LEU:O	2.08	0.53
11:K:109:TRP:C	11:K:109:TRP:CD1	2.87	0.53
1:A:298:PHE:C	1:A:298:PHE:CD2	2.87	0.53
2:B:324:ILE:HD11	2:B:329:THR:HG22	1.91	0.53
2:B:409:ALA:C	2:B:413:LEU:HD12	2.34	0.53
3:C:244:VAL:HG11	11:K:105:PHE:CE2	2.43	0.53
5:E:158:SER:O	5:E:162:ARG:HG3	2.07	0.53
1:A:535:THR:O	1:A:536:LEU:O	2.27	0.53
1:A:664:THR:HA	1:A:742:ASN:ND2	2.19	0.53
1:A:1005:GLU:O	1:A:1006:ILE:C	2.51	0.53
1:A:1016:THR:O	1:A:1019:CYS:HB2	2.09	0.53
1:A:1144:LYS:HA	1:A:1268:LEU:HD22	1.91	0.53
2:B:64:CYS:HA	2:B:67:SER:OG	2.08	0.53
2:B:825:VAL:HG11	2:B:1012:ILE:HD11	1.90	0.53
2:B:990:ILE:HG22	2:B:991:GLY:N	2.24	0.53
3:C:20:PHE:CE1	3:C:230:MET:HE3	2.44	0.53
1:A:23:SER:HB2	1:A:233:TRP:CE2	2.44	0.53
1:A:381:THR:O	1:A:384:ASN:OD1	2.27	0.53
2:B:124:TYR:CD2	2:B:124:TYR:O	2.61	0.53
2:B:652:LYS:O	2:B:689:LEU:HD22	2.09	0.53
2:B:941:LEU:HG	2:B:942:ARG:N	2.23	0.53
6:F:69:LEU:HB3	6:F:71:GLU:OE1	2.09	0.53
7:G:56:ILE:O	7:G:57:GLN:HB2	2.09	0.53
1:A:386:ASP:OD1	1:A:386:ASP:N	2.38	0.53
1:A:396:PRO:O	1:A:397:ASN:OD1	2.27	0.53
1:A:852:TYR:HD2	1:A:1060:PRO:CB	2.22	0.53
1:A:868:TYR:HE1	1:A:1064:VAL:CG1	2.17	0.53
1:A:1037:LEU:HD13	1:A:1041:ALA:HB1	1.90	0.53
2:B:581:PHE:HA	2:B:585:VAL:O	2.08	0.53
2:B:1096:ARG:CG	2:B:1097:HIS:H	2.22	0.53
2:B:1130:PHE:CE1	2:B:1134:GLU:HB3	2.44	0.53
4:D:56:ARG:HH21	4:D:155:ARG:HG2	1.73	0.53
6:F:87:LYS:HE2	6:F:88:TYR:HE1	1.73	0.53
6:F:133:VAL:HG21	7:G:58:ARG:HE	1.74	0.53
7:G:121:PHE:CE2	7:G:123:ALA:HB2	2.44	0.53
7:G:150:CYS:C	7:G:151:ILE:HD12	2.33	0.53
9:I:8:ARG:HG3	9:I:34:TYR:CE1	2.44	0.53
1:A:15:LYS:HB3	2:B:1220:ARG:NH2	2.23	0.53
1:A:315:LEU:HD22	1:A:319:GLY:O	2.08	0.53
1:A:1098:VAL:HB	1:A:1099:PRO:HD3	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1313:LEU:HD23	1:A:1338:VAL:CG2	2.37	0.53
2:B:120:ARG:HH12	12:L:54:ARG:NH1	2.07	0.53
2:B:493:SER:N	2:B:751:VAL:HG11	2.24	0.53
2:B:553:PRO:O	2:B:557:PHE:HB2	2.09	0.53
2:B:1084:GLN:N	2:B:1084:GLN:HE21	2.07	0.53
5:E:78:LEU:HG	5:E:107:THR:CG2	2.39	0.53
7:G:5:LYS:HG3	7:G:6:ASP:N	2.23	0.53
1:A:970:THR:HG22	1:A:971:PHE:CD1	2.44	0.52
2:B:289:LEU:HD13	2:B:375:ALA:HB2	1.91	0.52
2:B:655:LYS:HD2	2:B:658:ILE:CG2	2.39	0.52
2:B:801:LYS:O	2:B:801:LYS:HG3	2.09	0.52
4:D:185:CYS:SG	4:D:190:GLU:HB3	2.49	0.52
7:G:55:ASP:HB3	7:G:73:LYS:CB	2.39	0.52
8:H:145:ARG:O	8:H:146:ARG:CB	2.56	0.52
9:I:13:MET:HG2	9:I:14:LEU:N	2.23	0.52
11:K:40:HIS:CE1	11:K:63:VAL:HG11	2.43	0.52
12:L:30:ILE:CD1	12:L:59:ALA:HB2	2.40	0.52
1:A:1227:ILE:C	1:A:1228:TRP:CE3	2.86	0.52
1:A:1394:THR:HG22	1:A:1395:GLY:N	2.24	0.52
2:B:259:TYR:HD1	2:B:259:TYR:H	1.56	0.52
2:B:744:HIS:HD2	2:B:746:SER:OG	1.92	0.52
3:C:62:PHE:O	3:C:66:ARG:HG3	2.09	0.52
1:A:34:LYS:HB2	1:A:36:ARG:NH1	2.24	0.52
1:A:122:MET:HA	1:A:141:LEU:HD13	1.91	0.52
1:A:285:PRO:O	1:A:287:HIS:N	2.42	0.52
1:A:370:ILE:C	1:A:372:LYS:N	2.64	0.52
1:A:894:GLU:O	1:A:898:ARG:HB3	2.10	0.52
1:A:1258:HIS:O	1:A:1262:LYS:HE3	2.10	0.52
1:A:1346:ALA:O	1:A:1347:ALA:C	2.52	0.52
2:B:102:VAL:HG21	2:B:122:LEU:HD13	1.91	0.52
2:B:126:SER:O	2:B:169:ARG:HA	2.09	0.52
2:B:732:SER:HB3	2:B:734:HIS:CE1	2.45	0.52
2:B:878:GLN:H	2:B:934:LYS:NZ	2.07	0.52
2:B:899:ILE:HG22	2:B:900:ALA:O	2.09	0.52
2:B:1194:ILE:HD11	2:B:1196:ILE:HG21	1.90	0.52
3:C:46:ILE:HA	3:C:159:ALA:HA	1.91	0.52
3:C:115:SER:CB	3:C:142:VAL:HB	2.39	0.52
4:D:156:ASP:C	4:D:158:GLU:N	2.64	0.52
5:E:85:GLU:OE1	5:E:88:VAL:HA	2.10	0.52
1:A:129:LYS:O	1:A:130:ASP:CB	2.56	0.52
1:A:257:ARG:HG2	1:A:259:GLU:HG2	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:284:ILE:HD13	2:B:333:PHE:HD2	1.75	0.52
2:B:421:PHE:CE1	2:B:424:LEU:HD23	2.43	0.52
2:B:508:LEU:N	2:B:512:ARG:HH21	2.07	0.52
2:B:684:LEU:HA	2:B:689:LEU:HD12	1.92	0.52
2:B:785:TYR:C	2:B:787:VAL:N	2.67	0.52
2:B:848:ARG:NH2	2:B:996:ARG:HD2	2.15	0.52
6:F:109:VAL:CG1	6:F:110:ASP:N	2.71	0.52
8:H:84:ALA:HA	8:H:87:ARG:HD2	1.92	0.52
10:J:52:THR:HG22	10:J:52:THR:O	2.08	0.52
12:L:38:LEU:HD11	12:L:49:LYS:HE2	1.92	0.52
1:A:164:ARG:HG2	1:A:165:GLY:H	1.74	0.52
1:A:852:TYR:CD2	1:A:1060:PRO:CB	2.91	0.52
1:A:962:ARG:O	1:A:965:GLN:N	2.38	0.52
1:A:967:ALA:CA	1:A:1044:TRP:HZ3	2.22	0.52
1:A:1036:ARG:NH1	1:A:1036:ARG:CG	2.72	0.52
1:A:1037:LEU:HD12	1:A:1042:PHE:HA	1.91	0.52
4:D:59:ILE:HG23	7:G:47:CYS:SG	2.49	0.52
5:E:55:ARG:HD2	5:E:84:ASP:HA	1.92	0.52
6:F:78:GLN:O	6:F:78:GLN:HG2	2.10	0.52
7:G:27:LYS:O	7:G:31:LEU:HG	2.10	0.52
8:H:115:TYR:CE2	8:H:124:ARG:HG3	2.43	0.52
1:A:388:LEU:O	1:A:392:VAL:HG23	2.09	0.52
1:A:596:THR:O	1:A:598:LEU:N	2.42	0.52
1:A:864:ILE:HG21	1:A:1370:LEU:CD2	2.39	0.52
1:A:929:LEU:H	1:A:929:LEU:CD2	2.22	0.52
1:A:1219:THR:CG2	1:A:1220:PHE:N	2.72	0.52
1:A:1364:ASN:O	1:A:1365:TYR:C	2.53	0.52
2:B:203:PHE:CD2	2:B:461:LEU:HD21	2.45	0.52
2:B:274:PRO:HB2	2:B:359:GLU:HB3	1.91	0.52
2:B:637:LEU:HD12	2:B:693:ILE:CD1	2.40	0.52
2:B:655:LYS:HE3	2:B:659:ALA:HB2	1.92	0.52
5:E:136:ASN:O	5:E:140:LEU:HG	2.09	0.52
6:F:117:PRO:HA	6:F:120:ILE:HB	1.90	0.52
7:G:139:ILE:HD13	7:G:140:LYS:HG3	1.92	0.52
1:A:67:CYS:O	1:A:68:GLN:C	2.52	0.52
1:A:1289:ARG:HD3	1:A:1290:LYS:O	2.10	0.52
1:A:1329:THR:HG23	1:A:1331:SER:H	1.74	0.52
2:B:310:MET:O	2:B:311:LEU:C	2.52	0.52
2:B:427:ASP:O	2:B:430:ARG:HB2	2.10	0.52
2:B:702:LEU:O	2:B:739:THR:N	2.42	0.52
2:B:791:THR:O	2:B:792:MET:HB3	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:860:MET:HG2	2:B:861:ASP:N	2.24	0.52
7:G:111:THR:CG2	7:G:114:LEU:HD22	2.38	0.52
8:H:30:SER:HB3	8:H:36:CYS:HB3	1.91	0.52
9:I:7:CYS:C	9:I:8:ARG:O	2.51	0.52
1:A:28:ARG:HG3	1:A:83:HIS:ND1	2.24	0.52
1:A:567:LYS:CB	1:A:568:PRO:CD	2.74	0.52
1:A:1041:ALA:O	1:A:1045:VAL:HG23	2.09	0.52
1:A:1121:GLU:CG	1:A:1122:PRO:HD2	2.39	0.52
1:A:1272:THR:C	1:A:1273:LEU:HD12	2.34	0.52
1:A:1365:TYR:HB2	5:E:203:GLU:OE1	2.10	0.52
1:A:1378:GLN:O	5:E:177:ARG:HD2	2.10	0.52
2:B:345:LYS:HA	2:B:348:ARG:HG2	1.91	0.52
2:B:766:ARG:NH2	2:B:1020:ARG:HB3	2.25	0.52
2:B:958:GLN:C	2:B:960:GLY:H	2.18	0.52
3:C:178:PHE:HD2	3:C:178:PHE:C	2.18	0.52
5:E:153:HIS:ND1	5:E:153:HIS:N	2.58	0.52
5:E:177:ARG:CD	5:E:215:MET:HE2	2.39	0.52
11:K:61:TYR:CD1	11:K:61:TYR:C	2.88	0.52
1:A:578:LEU:C	1:A:580:VAL:H	2.18	0.52
1:A:645:LEU:O	1:A:649:ILE:HG13	2.09	0.52
1:A:814:PHE:HE1	2:B:519:TRP:CA	2.23	0.52
1:A:1313:LEU:C	1:A:1315:GLU:N	2.68	0.52
1:A:1446:ASP:O	1:A:1447:GLU:C	2.52	0.52
2:B:640:VAL:HG22	2:B:651:LEU:HD23	1.92	0.52
2:B:774:GLY:C	2:B:776:GLN:H	2.18	0.52
5:E:168:TYR:HB3	5:E:170:LEU:HG	1.90	0.52
1:A:130:ASP:H	1:A:134:ARG:HH21	1.58	0.52
1:A:649:ILE:O	1:A:650:GLN:C	2.53	0.52
1:A:933:TYR:HD2	1:A:933:TYR:C	2.17	0.52
1:A:1377:THR:HG22	5:E:176:PRO:HB3	1.92	0.52
2:B:483:LEU:HD12	2:B:484:ASN:N	2.25	0.52
2:B:541:LEU:HB2	2:B:747:MET:CE	2.38	0.52
2:B:582:VAL:HG22	2:B:626:ILE:CG2	2.39	0.52
2:B:1084:GLN:C	2:B:1085:ILE:HD12	2.34	0.52
3:C:166:GLU:HA	11:K:6:ARG:HB3	1.92	0.52
3:C:178:PHE:C	3:C:178:PHE:CD2	2.88	0.52
8:H:82:PRO:C	8:H:84:ALA:N	2.68	0.52
11:K:33:ILE:HB	11:K:35:PHE:HE1	1.76	0.52
1:A:63:ARG:CG	1:A:74:MET:HE1	2.39	0.51
1:A:172:PRO:HG3	1:A:185:TRP:CZ2	2.46	0.51
1:A:401:GLY:H	1:A:435:HIS:HD2	1.58	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:653:VAL:O	1:A:654:ASN:C	2.53	0.51
1:A:693:VAL:O	1:A:693:VAL:HG12	2.10	0.51
1:A:741:ASN:C	1:A:741:ASN:ND2	2.67	0.51
2:B:479:VAL:O	2:B:480:SER:HB3	2.09	0.51
2:B:781:PHE:HD1	2:B:782:LEU:HG	1.75	0.51
3:C:92:CYS:SG	3:C:94:LYS:HB2	2.50	0.51
6:F:146:TRP:HB3	6:F:151:LEU:HD11	1.92	0.51
11:K:74:ARG:C	11:K:75:ILE:HD13	2.36	0.51
1:A:656:TRP:CH2	1:A:660:ASN:CG	2.88	0.51
1:A:933:TYR:CD2	1:A:933:TYR:C	2.88	0.51
1:A:1218:GLN:O	1:A:1221:LYS:HE3	2.10	0.51
1:A:1345:ARG:O	1:A:1348:LEU:HB3	2.10	0.51
2:B:428:ILE:HG22	2:B:432:MET:HE2	1.91	0.51
2:B:527:THR:OG1	2:B:528:PRO:HD2	2.09	0.51
2:B:770:GLN:HB2	2:B:985:GLY:N	2.24	0.51
5:E:96:PHE:CD1	5:E:100:ILE:HD11	2.45	0.51
6:F:134:ILE:HG22	6:F:136:ARG:HG3	1.92	0.51
6:F:138:LEU:O	6:F:139:PRO:C	2.52	0.51
1:A:364:VAL:O	1:A:364:VAL:HG13	2.09	0.51
1:A:416:ARG:C	1:A:417:TYR:HD2	2.18	0.51
1:A:419:LYS:HG3	1:A:420:ARG:HG3	1.92	0.51
1:A:577:ILE:HA	1:A:580:VAL:CG2	2.39	0.51
1:A:913:LEU:HD12	1:A:914:GLU:H	1.73	0.51
1:A:1170:ILE:HG22	1:A:1174:PHE:CE1	2.46	0.51
2:B:54:PHE:CE2	2:B:59:LEU:HD13	2.45	0.51
2:B:101:MET:HE3	2:B:169:ARG:HH12	1.75	0.51
2:B:287:ARG:HG2	2:B:292:ILE:HA	1.92	0.51
2:B:459:TYR:CE1	2:B:469:GLN:HB2	2.46	0.51
2:B:770:GLN:HG2	2:B:983:ARG:C	2.36	0.51
6:F:100:GLN:O	6:F:103:MET:HB2	2.10	0.51
7:G:165:GLU:HG3	7:G:168:LEU:HD12	1.90	0.51
10:J:7:CYS:CB	10:J:49:MET:HE3	2.41	0.51
12:L:61:THR:HG21	12:L:63:ARG:NE	2.25	0.51
1:A:350:ARG:C	1:A:351:THR:CG2	2.84	0.51
1:A:492:PRO:HG3	1:A:501:LEU:CD1	2.36	0.51
1:A:1037:LEU:HD12	1:A:1042:PHE:CA	2.40	0.51
2:B:311:LEU:O	2:B:312:GLU:C	2.54	0.51
2:B:378:LEU:O	2:B:382:ILE:HG13	2.10	0.51
2:B:408:LEU:CD1	2:B:545:ILE:HG21	2.41	0.51
2:B:522:VAL:CG1	2:B:537:LYS:HB3	2.40	0.51
2:B:640:VAL:O	2:B:640:VAL:HG12	2.09	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:886:LYS:HE2	2:B:940:PRO:CG	2.40	0.51
1:A:18:GLN:HG2	1:A:1418:LEU:HD13	1.91	0.51
1:A:108:MET:HE3	1:A:210:ILE:CD1	2.40	0.51
1:A:151:ASP:HA	1:A:162:VAL:O	2.11	0.51
1:A:496:GLU:O	1:A:499:ALA:HB3	2.11	0.51
1:A:917:SER:C	1:A:919:ILE:H	2.19	0.51
1:A:1324:PRO:CB	5:E:142:VAL:HG11	2.40	0.51
2:B:203:PHE:HD2	2:B:461:LEU:HD21	1.75	0.51
2:B:408:LEU:HD13	2:B:545:ILE:HG21	1.93	0.51
3:C:186:LEU:HB3	3:C:188:HIS:CD2	2.45	0.51
7:G:1:MET:HE2	7:G:1:MET:C	2.35	0.51
7:G:122:ASN:HB2	7:G:131:GLN:HG2	1.92	0.51
1:A:87:ALA:HB3	1:A:276:LEU:HD23	1.92	0.51
1:A:340:LEU:HD21	2:B:1200:ALA:N	2.26	0.51
1:A:853:ASP:HB3	6:F:138:LEU:HD21	1.93	0.51
1:A:981:LEU:HD23	1:A:1039:LYS:HA	1.93	0.51
1:A:1008:GLN:O	1:A:1011:GLN:HB3	2.11	0.51
1:A:1348:LEU:HD23	1:A:1372:VAL:HG13	1.92	0.51
1:A:1421:CYS:C	1:A:1423:GLY:H	2.19	0.51
2:B:638:PHE:HE1	2:B:743:ILE:HA	1.74	0.51
2:B:1193:GLN:C	2:B:1194:ILE:HG22	2.34	0.51
4:D:26:THR:C	4:D:28:GLN:H	2.18	0.51
11:K:12:LEU:HA	11:K:37:LYS:HG3	1.91	0.51
1:A:44:THR:O	1:A:45:GLN:HB2	2.10	0.51
1:A:68:GLN:NE2	1:A:68:GLN:O	2.43	0.51
1:A:441:PRO:O	1:A:492:PRO:HG2	2.11	0.51
1:A:1011:GLN:NE2	1:A:1015:VAL:HG21	2.25	0.51
2:B:114:PRO:O	2:B:118:ARG:HG3	2.09	0.51
2:B:514:LEU:HD12	2:B:518:HIS:HD2	1.76	0.51
2:B:800:GLN:CB	10:J:52:THR:HG22	2.38	0.51
2:B:822:ASN:HD22	2:B:822:ASN:N	2.08	0.51
3:C:77:ILE:N	3:C:129:ILE:HD11	2.25	0.51
3:C:115:SER:HB2	3:C:142:VAL:HB	1.93	0.51
4:D:18:VAL:O	4:D:19:GLU:CB	2.57	0.51
8:H:31:THR:O	8:H:31:THR:HG22	2.10	0.51
12:L:44:ASP:O	12:L:45:ALA:HB3	2.10	0.51
1:A:338:GLY:HA2	2:B:1129:ARG:HH22	1.75	0.51
1:A:855:THR:HA	1:A:866:PHE:O	2.11	0.51
2:B:575:PRO:HG2	2:B:576:ASP:H	1.76	0.51
2:B:795:ILE:HD12	2:B:795:ILE:N	2.26	0.51
2:B:906:SER:O	2:B:907:GLY:C	2.54	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:1022:THR:HG23	2:B:1022:THR:O	2.10	0.51
3:C:53:THR:HG21	3:C:152:GLU:HB3	1.92	0.51
3:C:57:VAL:HG23	3:C:58:LEU:N	2.26	0.51
6:F:95:GLY:O	6:F:96:THR:C	2.54	0.51
7:G:111:THR:HG22	7:G:114:LEU:CD1	2.39	0.51
8:H:63:LEU:CB	8:H:90:ALA:HB3	2.41	0.51
8:H:95:TYR:HE2	8:H:97:MET:SD	2.34	0.51
1:A:204:THR:O	1:A:205:GLU:C	2.54	0.51
1:A:288:ALA:HA	1:A:291:GLU:CD	2.36	0.51
1:A:518:LYS:CB	1:A:519:PRO:HD2	2.40	0.51
1:A:586:ILE:HD11	1:A:633:VAL:HG22	1.93	0.51
1:A:666:ILE:HD12	1:A:667:GLY:N	2.24	0.51
1:A:852:TYR:CE1	6:F:136:ARG:HG2	2.45	0.51
1:A:866:PHE:O	1:A:867:ILE:HD12	2.09	0.51
1:A:1220:PHE:O	1:A:1221:LYS:HB2	2.11	0.51
1:A:1424:VAL:HG11	2:B:1139:ILE:HD13	1.92	0.51
1:A:1437:GLY:O	1:A:1441:PHE:CE2	2.64	0.51
2:B:32:ALA:HB3	2:B:658:ILE:HD11	1.92	0.51
2:B:698:GLU:HA	2:B:701:ILE:HG12	1.93	0.51
2:B:744:HIS:CG	2:B:745:PRO:HD2	2.45	0.51
2:B:1116:ARG:HG3	2:B:1198:TYR:CG	2.45	0.51
3:C:49:VAL:O	12:L:66:GLN:HA	2.10	0.51
3:C:184:ASN:ND2	3:C:189:THR:H	2.04	0.51
4:D:4:SER:O	4:D:5:THR:HB	2.11	0.51
8:H:59:ILE:O	8:H:60:ALA:CB	2.59	0.51
8:H:63:LEU:HD23	8:H:90:ALA:C	2.36	0.51
1:A:403:LYS:C	1:A:415:LEU:HB2	2.36	0.51
1:A:563:PRO:HB3	1:A:572:TRP:CE2	2.46	0.51
1:A:868:TYR:CE1	1:A:1064:VAL:HG22	2.45	0.51
1:A:1259:MET:HE2	1:A:1263:ILE:CG1	2.40	0.51
1:A:1384:VAL:HA	1:A:1389:PHE:CE2	2.46	0.51
2:B:957:ASN:ND2	2:B:961:LEU:HB2	2.21	0.51
4:D:50:LEU:HD23	4:D:50:LEU:H	1.75	0.51
4:D:160:VAL:HG12	4:D:161:GLY:N	2.26	0.51
6:F:84:TYR:CA	6:F:152:ILE:HD12	2.41	0.51
8:H:59:ILE:HG22	8:H:60:ALA:H	1.75	0.51
10:J:8:PHE:CE2	10:J:49:MET:SD	3.04	0.51
11:K:68:PHE:N	11:K:68:PHE:HD2	2.08	0.51
1:A:331:GLY:O	1:A:332:LYS:O	2.29	0.50
1:A:353:ILE:HD12	1:A:470:LEU:HD21	1.92	0.50
1:A:490:HIS:CD2	1:A:490:HIS:N	2.79	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:863:VAL:C	1:A:864:ILE:HD12	2.36	0.50
1:A:993:LEU:HB2	1:A:1046:LEU:HD22	1.92	0.50
1:A:1130:GLN:O	1:A:1134:ILE:HG13	2.11	0.50
1:A:1293:SER:OG	1:A:1294:PRO:HD2	2.11	0.50
1:A:1444:MET:HE2	6:F:133:VAL:O	2.10	0.50
2:B:37:PHE:CE1	2:B:41:LYS:HG3	2.45	0.50
2:B:345:LYS:HE3	2:B:349:ILE:HD11	1.93	0.50
2:B:972:LYS:HD2	2:B:1098:MET:HE1	1.93	0.50
3:C:18:VAL:HG12	3:C:18:VAL:O	2.10	0.50
6:F:147:SER:OG	6:F:150:GLU:HG3	2.12	0.50
10:J:12:LYS:O	10:J:13:VAL:C	2.54	0.50
11:K:75:ILE:HD13	11:K:75:ILE:N	2.26	0.50
12:L:38:LEU:HD12	12:L:39:SER:N	2.26	0.50
1:A:332:LYS:H	1:A:337:ARG:CB	2.23	0.50
1:A:532:ARG:CD	1:A:749:ALA:HB2	2.41	0.50
2:B:240:ILE:HD12	2:B:241:ARG:N	2.25	0.50
2:B:785:TYR:HE2	10:J:60:PHE:CZ	2.29	0.50
2:B:1058:LEU:HD23	2:B:1061:GLU:OE2	2.10	0.50
3:C:66:ARG:NH2	10:J:3:VAL:O	2.45	0.50
6:F:101:ILE:HD11	6:F:107:VAL:HG13	1.92	0.50
11:K:68:PHE:O	11:K:70:ARG:HG2	2.11	0.50
1:A:69:THR:HG21	2:B:1174:LYS:HZ2	1.73	0.50
1:A:184:SER:HB3	1:A:199:LEU:HD23	1.92	0.50
1:A:427:GLN:HG3	1:A:430:TRP:CZ2	2.45	0.50
1:A:881:GLN:HE22	1:A:959:ASN:HA	1.75	0.50
2:B:91:SER:OG	2:B:133:LYS:HB2	2.12	0.50
2:B:221:ASN:OD1	2:B:242:SER:HA	2.10	0.50
2:B:936:ASP:CG	2:B:937:ALA:H	2.19	0.50
2:B:1037:LEU:H	2:B:1037:LEU:HD12	1.77	0.50
8:H:63:LEU:CG	8:H:90:ALA:HB3	2.41	0.50
1:A:46:THR:O	1:A:47:ARG:C	2.55	0.50
1:A:849:MET:CE	1:A:1061:GLY:HA2	2.42	0.50
2:B:60:GLN:NE2	2:B:95:ILE:HB	2.25	0.50
3:C:76:ASP:O	3:C:79:GLN:HG2	2.11	0.50
3:C:261:ALA:HA	3:C:264:GLN:OE1	2.12	0.50
4:D:47:LEU:HD13	4:D:48:ILE:N	2.26	0.50
5:E:164:LEU:HD22	5:E:211:TYR:CD2	2.45	0.50
1:A:442:VAL:O	1:A:457:ALA:HA	2.10	0.50
1:A:814:PHE:HD2	1:A:814:PHE:C	2.19	0.50
1:A:966:ASN:O	1:A:970:THR:CB	2.60	0.50
1:A:1329:THR:O	1:A:1329:THR:HG22	2.10	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:323:VAL:O	2:B:323:VAL:HG12	2.10	0.50
2:B:796:LEU:HD21	2:B:821:GLN:HG3	1.92	0.50
2:B:912:ILE:HD11	2:B:966:VAL:HG23	1.92	0.50
2:B:971:THR:OG1	3:C:61:GLU:HG3	2.12	0.50
3:C:8:VAL:C	3:C:9:LYS:HG3	2.36	0.50
4:D:35:LEU:HD11	4:D:173:HIS:CD2	2.46	0.50
7:G:81:PRO:O	7:G:82:PHE:CD2	2.65	0.50
8:H:58:THR:HG22	8:H:59:ILE:N	2.26	0.50
8:H:118:PHE:HD1	8:H:121:LEU:O	1.95	0.50
11:K:56:VAL:O	11:K:56:VAL:HG12	2.11	0.50
1:A:302:THR:HA	1:A:305:ASP:O	2.12	0.50
1:A:353:ILE:HG21	1:A:487:MET:CG	2.40	0.50
1:A:354:SER:HA	1:A:482:PHE:CE2	2.47	0.50
1:A:464:PRO:O	1:A:465:TYR:O	2.30	0.50
1:A:573:SER:OG	1:A:576:GLN:HB2	2.11	0.50
1:A:858:ASN:ND2	1:A:861:GLY:H	2.09	0.50
1:A:1120:LEU:O	1:A:1323:ASP:HB2	2.12	0.50
2:B:628:THR:HG23	2:B:628:THR:O	2.10	0.50
2:B:637:LEU:HB3	2:B:740:HIS:HB3	1.93	0.50
2:B:780:VAL:HG21	10:J:56:LEU:HD11	1.94	0.50
2:B:1077:THR:OG1	2:B:1079:LYS:HB3	2.11	0.50
2:B:1183:LYS:HD2	2:B:1183:LYS:O	2.11	0.50
2:B:1212:ILE:O	2:B:1212:ILE:HG22	2.11	0.50
4:D:128:VAL:O	4:D:131:GLU:HB3	2.12	0.50
5:E:80:VAL:HG12	5:E:82:PHE:HE1	1.77	0.50
7:G:106:MET:HG2	7:G:107:LYS:N	2.25	0.50
1:A:185:TRP:HE3	1:A:185:TRP:H	1.58	0.50
1:A:489:LEU:C	1:A:490:HIS:CD2	2.90	0.50
1:A:973:ILE:HG21	1:A:1036:ARG:O	2.12	0.50
1:A:1329:THR:H	1:A:1335:ILE:HD11	1.77	0.50
2:B:370:PHE:HD2	2:B:373:ARG:HD2	1.77	0.50
2:B:429:PHE:CD2	2:B:433:GLN:NE2	2.79	0.50
2:B:429:PHE:HD2	2:B:433:GLN:NE2	2.10	0.50
2:B:877:PRO:CA	2:B:934:LYS:HZ3	2.25	0.50
2:B:1107:ALA:O	2:B:1108:ARG:HB2	2.11	0.50
2:B:1147:LEU:O	2:B:1148:LYS:C	2.53	0.50
5:E:117:THR:HG22	5:E:119:SER:H	1.76	0.50
12:L:27:LEU:HD13	12:L:37:LYS:HG2	1.92	0.50
1:A:129:LYS:O	1:A:130:ASP:HB2	2.12	0.50
1:A:379:VAL:HG12	1:A:380:VAL:H	1.76	0.50
1:A:1011:GLN:NE2	1:A:1015:VAL:CG2	2.75	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1323:ASP:OD2	1:A:1325:THR:HG22	2.12	0.50
2:B:60:GLN:HE22	2:B:95:ILE:HB	1.76	0.50
2:B:844:SER:O	2:B:847:ASP:HB2	2.12	0.50
3:C:237:SER:C	3:C:238:ILE:HG13	2.37	0.50
4:D:220:LEU:HD23	4:D:221:TYR:N	2.23	0.50
5:E:23:VAL:O	5:E:28:TYR:HB2	2.11	0.50
7:G:53:ASN:HD22	7:G:53:ASN:N	2.10	0.50
8:H:106:GLU:HG2	8:H:112:ILE:HD11	1.92	0.50
8:H:139:ASN:O	8:H:140:ALA:CB	2.59	0.50
1:A:79:GLY:HA3	1:A:243:PRO:CB	2.37	0.50
1:A:476:SER:N	1:A:477:PRO:CD	2.75	0.50
1:A:947:PHE:N	1:A:947:PHE:CD2	2.80	0.50
1:A:1198:ASP:HB3	1:A:1201:ALA:CB	2.41	0.50
1:A:1445:ILE:HD11	7:G:70:PHE:CE2	2.47	0.50
2:B:44:VAL:O	2:B:45:SER:C	2.54	0.50
2:B:1078:GLY:O	2:B:1079:LYS:C	2.55	0.50
2:B:1160:VAL:CG1	2:B:1161:HIS:H	2.25	0.50
3:C:127:ARG:O	3:C:129:ILE:N	2.45	0.50
4:D:54:GLU:CD	4:D:164:ILE:HD11	2.37	0.50
4:D:125:SER:O	4:D:128:VAL:HB	2.12	0.50
11:K:35:PHE:N	11:K:35:PHE:CD1	2.78	0.50
1:A:24:PRO:O	1:A:28:ARG:HB2	2.12	0.49
1:A:40:THR:CG2	1:A:54:ASN:HD21	2.24	0.49
1:A:73:GLY:O	1:A:74:MET:C	2.55	0.49
1:A:100:LYS:HE2	1:A:104:GLU:OE2	2.12	0.49
1:A:903:ASN:HD22	1:A:905:ASP:N	1.87	0.49
1:A:1331:SER:OG	1:A:1333:ILE:HG22	2.12	0.49
2:B:785:TYR:C	2:B:787:VAL:H	2.20	0.49
2:B:1096:ARG:HB2	2:B:1096:ARG:NH1	2.27	0.49
3:C:20:PHE:HE1	3:C:230:MET:HE3	1.77	0.49
3:C:70:ILE:O	3:C:72:LEU:CD1	2.56	0.49
3:C:77:ILE:HA	3:C:129:ILE:HD11	1.93	0.49
3:C:93:ASP:OD2	3:C:122:SER:HB2	2.11	0.49
7:G:9:LEU:HG	7:G:11:ILE:CG1	2.41	0.49
1:A:490:HIS:HE1	2:B:1130:PHE:HB2	1.77	0.49
1:A:1198:ASP:HB3	1:A:1201:ALA:HB3	1.94	0.49
2:B:58:THR:O	2:B:62:ILE:HG13	2.12	0.49
2:B:1065:GLN:HE21	2:B:1067:ARG:H	1.58	0.49
4:D:40:HIS:CD2	7:G:73:LYS:HE3	2.47	0.49
5:E:100:ILE:HG23	5:E:105:PHE:HD1	1.74	0.49
5:E:106:GLN:HA	5:E:130:ALA:HB2	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:E:182:ASP:OD1	5:E:183:PRO:HD2	2.12	0.49
1:A:9:ALA:HB3	2:B:1193:GLN:HB2	1.93	0.49
1:A:58:LEU:HD13	1:A:243:PRO:HA	1.94	0.49
1:A:91:PHE:N	1:A:297:GLN:HE22	2.09	0.49
1:A:374:LEU:HD11	2:B:1105:ALA:HB1	1.94	0.49
1:A:451:HIS:HD2	1:A:453:MET:HE2	1.76	0.49
1:A:1322:ILE:HG13	1:A:1323:ASP:N	2.26	0.49
2:B:545:ILE:HG22	2:B:546:SER:O	2.12	0.49
3:C:183:TRP:O	3:C:184:ASN:C	2.55	0.49
4:D:6:SER:HB3	7:G:8:SER:OG	2.11	0.49
4:D:155:ARG:NH2	4:D:221:TYR:CD1	2.81	0.49
5:E:3:GLN:HE21	5:E:52:ARG:HH12	1.58	0.49
7:G:104:GLY:C	7:G:106:MET:H	2.20	0.49
10:J:57:ILE:CG2	10:J:58:GLU:N	2.75	0.49
1:A:144:THR:O	1:A:146:MET:HE2	2.12	0.49
1:A:253:ASN:HB3	2:B:935:ARG:NH1	2.28	0.49
1:A:370:ILE:HG22	1:A:374:LEU:HD13	1.92	0.49
1:A:1030:ARG:HD3	1:A:1035:TYR:OH	2.13	0.49
1:A:1197:LEU:HD12	1:A:1209:MET:SD	2.52	0.49
2:B:168:GLY:HA2	2:B:454:THR:OG1	2.12	0.49
2:B:269:ILE:HG21	2:B:282:ILE:HD13	1.95	0.49
2:B:780:VAL:HB	2:B:817:LEU:HD22	1.93	0.49
2:B:1016:ALA:HB1	2:B:1020:ARG:HE	1.76	0.49
3:C:165:LYS:O	11:K:6:ARG:HD2	2.13	0.49
4:D:126:ILE:C	4:D:128:VAL:N	2.70	0.49
6:F:139:PRO:O	6:F:141:GLY:N	2.45	0.49
7:G:8:SER:HA	7:G:72:VAL:O	2.12	0.49
7:G:116:PRO:C	7:G:118:ASP:H	2.19	0.49
10:J:1:MET:H1	10:J:56:LEU:N	2.10	0.49
11:K:57:LEU:H	11:K:77:THR:HA	1.77	0.49
12:L:38:LEU:O	12:L:39:SER:HB2	2.12	0.49
1:A:425:GLN:CD	1:A:425:GLN:N	2.71	0.49
1:A:670:ILE:HD13	1:A:670:ILE:H	1.78	0.49
1:A:868:TYR:CE1	1:A:1064:VAL:CG1	2.92	0.49
1:A:1157:ASP:C	1:A:1159:ARG:H	2.20	0.49
1:A:1370:LEU:O	1:A:1371:LEU:C	2.55	0.49
1:A:1444:MET:HE1	6:F:135:ARG:CB	2.43	0.49
2:B:222:ILE:HD11	2:B:627:PHE:HZ	1.77	0.49
2:B:510:LYS:O	2:B:511:PRO:C	2.55	0.49
2:B:822:ASN:ND2	10:J:52:THR:HG21	2.27	0.49
2:B:860:MET:HE2	2:B:965:LYS:HE2	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:900:ALA:CB	12:L:61:THR:OG1	2.61	0.49
2:B:1138:MET:HE2	2:B:1146:PHE:CE2	2.47	0.49
3:C:86:CYS:SG	3:C:87:PHE:N	2.86	0.49
5:E:4:GLU:HB3	5:E:7:ARG:HE	1.76	0.49
12:L:61:THR:HG21	12:L:63:ARG:HE	1.77	0.49
1:A:16:GLU:CD	2:B:1220:ARG:HA	2.37	0.49
1:A:506:ALA:C	1:A:508:PRO:HD2	2.38	0.49
1:A:741:ASN:HD21	1:A:743:VAL:CG2	2.12	0.49
1:A:761:MET:HA	1:A:804:TYR:HB2	1.95	0.49
1:A:814:PHE:C	1:A:814:PHE:CD2	2.90	0.49
1:A:1424:VAL:O	1:A:1425:SER:C	2.56	0.49
2:B:176:SER:O	2:B:182:SER:HB3	2.13	0.49
2:B:604:ARG:HG3	2:B:611:PRO:HA	1.94	0.49
2:B:642:ASP:CA	2:B:649:LYS:HG3	2.42	0.49
9:I:17:ARG:NH2	9:I:30:ARG:NH2	2.61	0.49
1:A:335:ARG:O	1:A:339:ASN:HB2	2.13	0.49
1:A:384:ASN:O	1:A:385:ILE:C	2.54	0.49
1:A:967:ALA:HA	1:A:1044:TRP:CZ3	2.47	0.49
1:A:1100:ARG:NH1	1:A:1111:MET:HE2	2.27	0.49
1:A:1146:VAL:HG12	1:A:1201:ALA:O	2.12	0.49
1:A:1313:LEU:HG	1:A:1317:MET:HE3	1.94	0.49
1:A:1325:THR:O	5:E:147:HIS:CD2	2.66	0.49
2:B:1170:THR:HB	2:B:1182:CYS:HB2	1.95	0.49
3:C:45:ALA:HB1	3:C:72:LEU:HD13	1.94	0.49
3:C:109:SER:O	3:C:110:THR:C	2.55	0.49
5:E:93:MET:O	5:E:97:VAL:HG23	2.13	0.49
1:A:14:VAL:N	1:A:1432:GLN:HE22	2.01	0.49
1:A:34:LYS:CE	1:A:57:ARG:NH1	2.71	0.49
1:A:316:GLN:HG2	1:A:317:LYS:HG2	1.93	0.49
1:A:1407:GLU:CD	1:A:1407:GLU:H	2.21	0.49
2:B:167:ILE:HG22	2:B:453:ILE:HG13	1.94	0.49
2:B:226:PHE:CE2	2:B:398:ARG:NH2	2.80	0.49
2:B:298:LEU:HD22	2:B:298:LEU:N	2.27	0.49
2:B:361:LEU:N	2:B:362:PRO:HD3	2.26	0.49
3:C:15:LYS:O	3:C:240:VAL:HG22	2.13	0.49
3:C:112:ASN:HB3	3:C:114:TYR:HE1	1.78	0.49
4:D:71:LYS:HA	4:D:74:GLN:HG3	1.93	0.49
7:G:13:LEU:HD23	7:G:14:HIS:H	1.75	0.49
8:H:115:TYR:CZ	8:H:124:ARG:HG3	2.48	0.49
11:K:42:LEU:O	11:K:46:ILE:HG13	2.13	0.49
1:A:41:MET:CB	1:A:49:LYS:HA	2.43	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:224:PHE:CD2	1:A:231:PRO:CD	2.95	0.49
1:A:313:GLN:O	1:A:314:ALA:C	2.56	0.49
1:A:362:ASP:HB3	1:A:507:VAL:CG1	2.39	0.49
1:A:572:TRP:HH2	8:H:79:TRP:CZ3	2.31	0.49
1:A:883:LEU:HD21	1:A:956:LEU:HD11	1.95	0.49
1:A:886:ILE:HD11	1:A:943:LEU:C	2.37	0.49
1:A:1116:LEU:HB2	1:A:1311:VAL:HG22	1.93	0.49
1:A:1332:PHE:N	1:A:1332:PHE:CD2	2.80	0.49
2:B:578:THR:HG23	2:B:622:LYS:C	2.38	0.49
2:B:705:MET:N	2:B:710:LEU:HD12	2.25	0.49
2:B:980:PHE:HE1	2:B:990:ILE:HG13	1.78	0.49
2:B:1203:LEU:HA	2:B:1206:GLU:OE1	2.13	0.49
3:C:8:VAL:HG12	3:C:9:LYS:H	1.78	0.49
4:D:52:LEU:O	4:D:54:GLU:N	2.44	0.49
6:F:109:VAL:CG1	6:F:110:ASP:H	2.25	0.49
12:L:55:ILE:O	12:L:56:LEU:CB	2.60	0.49
1:A:507:VAL:CG2	1:A:521:MET:HE1	2.42	0.49
1:A:563:PRO:HB3	1:A:572:TRP:CZ2	2.47	0.49
1:A:567:LYS:CB	8:H:95:TYR:HA	2.43	0.49
1:A:741:ASN:HD22	1:A:743:VAL:N	2.11	0.49
1:A:901:LEU:HB2	1:A:926:GLN:HG2	1.94	0.49
1:A:997:LEU:HD22	1:A:1053:PHE:CD2	2.48	0.49
1:A:1437:GLY:O	1:A:1441:PHE:HE2	1.96	0.49
2:B:288:ALA:HB1	2:B:331:LEU:HG	1.94	0.49
2:B:290:GLY:O	2:B:292:ILE:HG13	2.13	0.49
2:B:581:PHE:O	2:B:626:ILE:HB	2.12	0.49
2:B:1106:ARG:HH11	2:B:1126:GLY:HA2	1.78	0.49
4:D:179:GLN:O	4:D:183:LEU:HD12	2.13	0.49
5:E:185:ALA:HA	5:E:190:LEU:HG	1.95	0.49
5:E:198:ILE:HB	5:E:210:SER:HB3	1.94	0.49
1:A:14:VAL:HB	1:A:1430:LEU:CD1	2.43	0.48
1:A:28:ARG:HG3	1:A:83:HIS:CE1	2.48	0.48
1:A:58:LEU:CB	1:A:80:HIS:O	2.61	0.48
1:A:252:PHE:O	1:A:253:ASN:HB3	2.13	0.48
1:A:341:MET:HE2	1:A:843:LYS:NZ	2.28	0.48
1:A:426:LEU:O	1:A:427:GLN:HG2	2.13	0.48
1:A:453:MET:HG2	1:A:520:CYS:SG	2.53	0.48
1:A:491:VAL:HG13	1:A:492:PRO:O	2.13	0.48
1:A:863:VAL:HG12	1:A:864:ILE:N	2.28	0.48
1:A:998:LEU:O	1:A:999:VAL:HG23	2.13	0.48
2:B:188:ASP:OD2	2:B:188:ASP:N	2.44	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:446:LEU:HD23	2:B:448:ILE:HD11	1.94	0.48
2:B:515:HIS:CD2	2:B:517:THR:OG1	2.66	0.48
2:B:876:LYS:HG3	2:B:893:LEU:HD12	1.95	0.48
3:C:167:HIS:CD2	12:L:70:ARG:HG2	2.49	0.48
4:D:126:ILE:C	4:D:128:VAL:H	2.20	0.48
12:L:47:ARG:HB2	12:L:47:ARG:NH1	2.27	0.48
1:A:81:PHE:CE2	1:A:243:PRO:HD3	2.48	0.48
1:A:125:ALA:O	1:A:128:ILE:HB	2.13	0.48
1:A:874:ASP:O	1:A:876:ALA:N	2.46	0.48
1:A:1152:ILE:HD11	9:I:44:TYR:HD2	1.77	0.48
1:A:1329:THR:CG2	1:A:1331:SER:H	2.26	0.48
2:B:190:TYR:HD2	10:J:63:TYR:CE2	2.31	0.48
2:B:983:ARG:HH11	2:B:1091:TYR:HB3	1.78	0.48
4:D:51:ASN:HB3	4:D:178:ALA:O	2.13	0.48
6:F:76:LYS:O	6:F:79:ARG:HD3	2.13	0.48
6:F:80:ALA:HB3	6:F:144:GLU:OE2	2.13	0.48
7:G:30:LEU:CD2	7:G:31:LEU:HD23	2.43	0.48
10:J:17:LYS:O	10:J:20:SER:N	2.45	0.48
1:A:254:GLU:O	1:A:255:SER:C	2.56	0.48
1:A:380:VAL:HB	1:A:430:TRP:H	1.78	0.48
1:A:419:LYS:C	1:A:421:ALA:H	2.21	0.48
2:B:202:TYR:CD2	2:B:202:TYR:N	2.65	0.48
2:B:1007:VAL:HG23	2:B:1008:PRO:CD	2.43	0.48
2:B:1079:LYS:N	3:C:27:LEU:HD21	2.29	0.48
3:C:68:GLY:O	3:C:169:LYS:HB2	2.13	0.48
3:C:235:VAL:HG12	10:J:13:VAL:CG1	2.43	0.48
5:E:152:LYS:HG3	5:E:154:ILE:CD1	2.43	0.48
7:G:13:LEU:HD21	7:G:17:PHE:HB2	1.95	0.48
11:K:31:VAL:HG12	11:K:32:VAL:N	2.28	0.48
1:A:567:LYS:HB3	8:H:95:TYR:HA	1.95	0.48
1:A:588:LEU:HB3	1:A:607:ILE:HB	1.95	0.48
1:A:981:LEU:CD2	1:A:1039:LYS:HA	2.42	0.48
1:A:1018:PHE:O	1:A:1019:CYS:C	2.57	0.48
1:A:1265:ASN:C	1:A:1267:MET:H	2.21	0.48
1:A:1393:ASN:O	1:A:1394:THR:O	2.30	0.48
2:B:108:VAL:HG23	2:B:109:THR:N	2.29	0.48
2:B:629:ASP:HB3	2:B:632:ARG:HE	1.78	0.48
2:B:973:ILE:HG23	2:B:974:PRO:HD2	1.95	0.48
2:B:1206:GLU:O	2:B:1209:ALA:HB3	2.13	0.48
3:C:167:HIS:O	3:C:169:LYS:N	2.47	0.48
4:D:56:ARG:HD3	4:D:149:THR:CA	2.41	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:K:34:THR:O	11:K:34:THR:HG22	2.13	0.48
1:A:81:PHE:HA	1:A:243:PRO:HG3	1.96	0.48
1:A:255:SER:OG	2:B:918:ILE:HD13	2.14	0.48
1:A:463:ILE:CB	1:A:464:PRO:CD	2.87	0.48
1:A:728:LYS:O	1:A:732:LEU:HB2	2.14	0.48
1:A:744:LYS:HG2	1:A:748:MET:HE2	1.96	0.48
2:B:120:ARG:CD	2:B:955:THR:HG21	2.44	0.48
2:B:582:VAL:HB	2:B:587:HIS:CD2	2.48	0.48
2:B:736:THR:O	2:B:736:THR:HG22	2.13	0.48
3:C:57:VAL:HG23	3:C:58:LEU:HD23	1.96	0.48
4:D:73:SER:CB	7:G:21:ARG:HH21	2.26	0.48
4:D:173:HIS:C	4:D:175:PHE:N	2.71	0.48
5:E:79:TRP:HB2	5:E:105:PHE:CE1	2.48	0.48
10:J:6:ARG:N	10:J:14:VAL:HB	2.28	0.48
12:L:38:LEU:HD21	12:L:48:CYS:HA	1.95	0.48
1:A:58:LEU:HD12	1:A:244:PRO:CD	2.44	0.48
1:A:257:ARG:HG2	1:A:258:GLY:H	1.78	0.48
1:A:722:LEU:HD23	1:A:799:PHE:CD1	2.49	0.48
1:A:773:LYS:H	1:A:773:LYS:HD2	1.78	0.48
1:A:946:VAL:C	1:A:947:PHE:HD2	2.21	0.48
1:A:1021:LEU:O	1:A:1024:SER:HB3	2.13	0.48
1:A:1151:GLU:HG2	9:I:45:ARG:HB2	1.96	0.48
2:B:642:ASP:HA	2:B:649:LYS:HG3	1.95	0.48
3:C:124:LEU:HB3	3:C:126:GLY:H	1.77	0.48
3:C:179:GLU:HG2	3:C:180:TYR:N	2.28	0.48
3:C:183:TRP:O	3:C:185:LYS:N	2.47	0.48
5:E:100:ILE:CG2	5:E:105:PHE:HB2	2.43	0.48
5:E:151:PRO:HD2	5:E:153:HIS:HE1	1.77	0.48
6:F:72:LYS:O	6:F:143:PHE:CD2	2.67	0.48
9:I:33:SER:O	9:I:35:VAL:HG23	2.13	0.48
12:L:59:ALA:O	12:L:60:ARG:O	2.31	0.48
1:A:446:ARG:HD2	1:A:480:ALA:HB2	1.94	0.48
1:A:722:LEU:N	1:A:722:LEU:HD12	2.29	0.48
1:A:1105:LEU:HD22	1:A:1384:VAL:HG21	1.96	0.48
1:A:1290:LYS:O	1:A:1291:VAL:HG23	2.14	0.48
2:B:516:ASN:ND2	2:B:516:ASN:N	2.62	0.48
2:B:637:LEU:HB2	2:B:693:ILE:HD12	1.94	0.48
2:B:979:LYS:C	2:B:980:PHE:CD1	2.91	0.48
3:C:57:VAL:CG1	10:J:60:PHE:HB3	2.37	0.48
8:H:58:THR:HG22	8:H:59:ILE:H	1.79	0.48
12:L:51:CYS:HB3	12:L:53:HIS:HB2	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:145:LYS:HE3	1:A:145:LYS:HA	1.96	0.48
1:A:245:PRO:O	2:B:1114:LEU:CD1	2.61	0.48
1:A:492:PRO:HA	1:A:497:THR:HG21	1.96	0.48
1:A:967:ALA:HA	1:A:1044:TRP:HZ3	1.77	0.48
1:A:993:LEU:O	1:A:994:GLN:C	2.56	0.48
1:A:1256:GLU:C	1:A:1258:HIS:H	2.22	0.48
1:A:1430:LEU:O	1:A:1432:GLN:HG3	2.12	0.48
2:B:247:GLY:HA2	2:B:418:LYS:NZ	2.28	0.48
2:B:640:VAL:O	2:B:641:GLU:C	2.57	0.48
2:B:757:PRO:O	2:B:758:PHE:HB2	2.13	0.48
3:C:240:VAL:O	3:C:241:ASP:C	2.56	0.48
6:F:100:GLN:HA	6:F:103:MET:HB2	1.96	0.48
8:H:48:PRO:O	8:H:49:VAL:HG23	2.14	0.48
1:A:350:ARG:O	1:A:351:THR:HG22	2.13	0.48
1:A:388:LEU:HB3	1:A:432:VAL:HG21	1.95	0.48
1:A:756:ILE:O	1:A:759:ALA:HB3	2.14	0.48
1:A:831:THR:O	1:A:834:THR:HB	2.14	0.48
1:A:1227:ILE:CA	1:A:1228:TRP:CE3	2.96	0.48
2:B:179:CYS:SG	2:B:181:LEU:HB2	2.54	0.48
2:B:496:ARG:O	2:B:539:LEU:HD12	2.14	0.48
2:B:796:LEU:O	2:B:799:PRO:HD3	2.14	0.48
2:B:1114:LEU:HD11	2:B:1202:LEU:HD11	1.94	0.48
3:C:70:ILE:HG22	3:C:71:PRO:HD2	1.95	0.48
3:C:97:VAL:HG21	3:C:129:ILE:HG22	1.96	0.48
4:D:32:GLU:HB3	4:D:33:PHE:CD1	2.49	0.48
5:E:61:GLN:HB2	5:E:79:TRP:CE3	2.49	0.48
5:E:153:HIS:CD2	5:E:184:VAL:HG11	2.49	0.48
7:G:61:ILE:HG13	7:G:68:ALA:HB2	1.94	0.48
8:H:125:LEU:HG	8:H:126:GLU:N	2.28	0.48
1:A:89:PRO:O	1:A:204:THR:HG21	2.14	0.48
1:A:105:CYS:O	1:A:114:LEU:HG	2.12	0.48
1:A:499:ALA:O	1:A:500:GLU:C	2.56	0.48
1:A:523:ILE:HG23	1:A:527:THR:HB	1.95	0.48
1:A:974:ASP:OD2	1:A:977:LYS:HB2	2.14	0.48
1:A:1227:ILE:CA	1:A:1228:TRP:HE3	2.27	0.48
2:B:181:LEU:O	2:B:182:SER:C	2.57	0.48
2:B:351:TYR:CD1	2:B:355:ILE:HD11	2.47	0.48
2:B:409:ALA:O	2:B:413:LEU:HD12	2.14	0.48
2:B:879:ARG:NH2	2:B:885:MET:HE2	2.29	0.48
2:B:1003:ALA:CB	3:C:179:GLU:HB2	2.43	0.48
3:C:3:GLU:HB3	11:K:104:ASN:ND2	2.29	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:4:SER:O	4:D:5:THR:CB	2.62	0.48
5:E:119:SER:HA	5:E:122:LYS:CG	2.43	0.48
5:E:151:PRO:HD2	5:E:153:HIS:CE1	2.49	0.48
7:G:22:MET:HE2	7:G:70:PHE:CE1	2.48	0.48
1:A:535:THR:CG2	1:A:616:VAL:HA	2.44	0.47
1:A:850:VAL:HG23	1:A:1064:VAL:CG2	2.44	0.47
1:A:1001:ARG:O	1:A:1002:GLY:O	2.32	0.47
1:A:1141:THR:CG2	1:A:1205:LYS:HD3	2.43	0.47
1:A:1260:LEU:HA	1:A:1263:ILE:CD1	2.44	0.47
1:A:1398:MET:HE1	1:A:1424:VAL:HG23	1.95	0.47
2:B:500:THR:O	2:B:535:LEU:HD22	2.14	0.47
2:B:758:PHE:CZ	2:B:1044:ALA:HA	2.49	0.47
2:B:862:GLN:HA	2:B:962:LYS:O	2.14	0.47
2:B:1192:TYR:CD1	2:B:1218:THR:HG21	2.48	0.47
4:D:49:ALA:HB3	4:D:174:PRO:O	2.13	0.47
8:H:130:ARG:HB3	8:H:133:ASN:HB2	1.96	0.47
9:I:17:ARG:HB3	9:I:26:LEU:HB3	1.96	0.47
12:L:32:ALA:HB3	12:L:55:ILE:HG13	1.92	0.47
1:A:515:GLN:HE22	1:A:1074:GLU:HB3	1.80	0.47
1:A:947:PHE:N	1:A:947:PHE:HD2	2.11	0.47
1:A:985:ASP:O	1:A:986:ILE:C	2.57	0.47
1:A:1261:LYS:O	1:A:1264:GLU:HB3	2.14	0.47
2:B:27:ALA:O	2:B:29:ASP:N	2.47	0.47
2:B:113:TYR:HD2	2:B:114:PRO:HD2	1.77	0.47
2:B:485:ARG:NH1	2:B:782:LEU:HD21	2.29	0.47
2:B:797:TYR:HB3	2:B:798:TYR:CD2	2.43	0.47
2:B:1046:PRO:O	2:B:1048:THR:HG23	2.14	0.47
3:C:82:TYR:CD2	3:C:84:ARG:NH2	2.82	0.47
4:D:175:PHE:O	4:D:176:GLU:C	2.57	0.47
5:E:84:ASP:O	5:E:86:PRO:HD3	2.14	0.47
7:G:56:ILE:HG23	7:G:57:GLN:N	2.26	0.47
7:G:112:LYS:HA	7:G:115:MET:HE3	1.96	0.47
8:H:37:LYS:HD2	8:H:126:GLU:OE2	2.13	0.47
11:K:22:ASP:O	11:K:31:VAL:HG13	2.13	0.47
12:L:63:ARG:O	12:L:64:LEU:C	2.57	0.47
1:A:27:VAL:C	1:A:29:ALA:N	2.71	0.47
1:A:532:ARG:C	1:A:534:LEU:H	2.22	0.47
1:A:1226:VAL:HG22	1:A:1240:CYS:CB	2.43	0.47
2:B:276:ILE:HD11	2:B:334:ILE:HG23	1.96	0.47
2:B:277:LYS:HG2	2:B:336:ARG:O	2.14	0.47
2:B:352:ALA:O	2:B:353:LYS:C	2.58	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:706:GLN:HA	2:B:707:PRO:HD2	1.72	0.47
2:B:743:ILE:H	2:B:743:ILE:HG12	1.41	0.47
2:B:1101:ASP:O	2:B:1122:ARG:CZ	2.63	0.47
4:D:71:LYS:HA	4:D:74:GLN:CG	2.44	0.47
5:E:131:THR:HG21	5:E:191:LYS:NZ	2.29	0.47
5:E:143:ASN:HB3	5:E:146:HIS:HB2	1.95	0.47
1:A:1289:ARG:HH22	1:A:1326:ARG:NH1	2.12	0.47
1:A:1364:ASN:O	1:A:1366:ARG:N	2.47	0.47
1:A:1391:ARG:O	1:A:1392:SER:O	2.33	0.47
2:B:222:ILE:C	2:B:240:ILE:HD13	2.39	0.47
2:B:365:THR:HG23	2:B:367:LEU:N	2.29	0.47
2:B:582:VAL:HG22	2:B:626:ILE:HG21	1.96	0.47
2:B:612:GLU:HG2	2:B:613:VAL:N	2.28	0.47
2:B:918:ILE:HD12	2:B:935:ARG:CZ	2.44	0.47
2:B:1205:GLN:O	2:B:1208:MET:HB2	2.15	0.47
3:C:41:ILE:HD12	3:C:243:VAL:HG13	1.96	0.47
4:D:154:PHE:HZ	4:D:214:LEU:CD1	2.28	0.47
5:E:124:VAL:CG1	5:E:132:ILE:HB	2.32	0.47
1:A:127:ALA:O	1:A:129:LYS:HG3	2.15	0.47
1:A:464:PRO:HG2	1:A:465:TYR:HD1	1.79	0.47
1:A:1037:LEU:HD12	1:A:1042:PHE:N	2.30	0.47
1:A:1173:HIS:HB3	1:A:1227:ILE:HD13	1.97	0.47
2:B:32:ALA:O	2:B:35:SER:HB2	2.15	0.47
2:B:43:LEU:HD11	2:B:811:TYR:O	2.14	0.47
3:C:125:MET:HB2	3:C:127:ARG:HE	1.78	0.47
7:G:151:ILE:HD12	7:G:151:ILE:N	2.30	0.47
12:L:38:LEU:CD1	12:L:49:LYS:HE2	2.44	0.47
1:A:336:ILE:HD11	2:B:1203:LEU:HD12	1.92	0.47
1:A:390:GLN:HE21	1:A:394:ASN:HD21	1.63	0.47
1:A:451:HIS:ND1	1:A:451:HIS:N	2.62	0.47
1:A:482:PHE:CD1	2:B:836:GLU:HB2	2.50	0.47
1:A:501:LEU:HD21	2:B:1146:PHE:CD1	2.49	0.47
1:A:675:THR:HG22	1:A:675:THR:O	2.15	0.47
1:A:845:LEU:O	1:A:846:GLU:C	2.58	0.47
1:A:852:TYR:HE2	1:A:1060:PRO:HB2	1.72	0.47
1:A:864:ILE:N	1:A:864:ILE:CD1	2.78	0.47
1:A:1400:CYS:HA	1:A:1408:ILE:HD12	1.97	0.47
2:B:190:TYR:CD1	10:J:62:ARG:HD3	2.49	0.47
2:B:498:THR:HB	2:B:537:LYS:O	2.14	0.47
2:B:834:ASN:O	2:B:1013:ASN:HB2	2.15	0.47
3:C:241:ASP:HB3	11:K:109:TRP:CZ3	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:G:8:SER:HB2	7:G:71:ASN:OD1	2.13	0.47
9:I:17:ARG:HG3	9:I:18:GLU:N	2.30	0.47
10:J:41:LEU:HD23	10:J:41:LEU:N	2.29	0.47
12:L:30:ILE:HG22	12:L:31:CYS:O	2.15	0.47
1:A:19:PHE:CZ	1:A:1397:LEU:CD2	2.92	0.47
1:A:377:PRO:HB2	1:A:431:LYS:HZ2	1.80	0.47
1:A:434:ARG:CZ	1:A:437:MET:HG3	2.44	0.47
1:A:444:PHE:HE2	1:A:470:LEU:HD13	1.80	0.47
1:A:567:LYS:HZ2	8:H:46:LEU:HB2	1.79	0.47
1:A:591:PHE:HD2	1:A:595:THR:CB	2.17	0.47
1:A:600:PRO:CA	8:H:25:ARG:NH1	2.75	0.47
1:A:826:ASP:O	1:A:830:LYS:HB2	2.14	0.47
1:A:1111:MET:SD	1:A:1114:PRO:HB3	2.54	0.47
1:A:1348:LEU:O	1:A:1352:VAL:HG23	2.14	0.47
2:B:25:ILE:HD11	2:B:651:LEU:CD1	2.41	0.47
2:B:335:GLY:O	2:B:336:ARG:HB2	2.14	0.47
2:B:523:CYS:SG	2:B:524:PRO:HD2	2.55	0.47
2:B:579:ARG:HA	2:B:589:VAL:HG22	1.97	0.47
2:B:868:MET:HE3	2:B:868:MET:HA	1.96	0.47
2:B:1023:VAL:HG13	2:B:1026:LEU:HD12	1.95	0.47
3:C:213:PRO:O	3:C:214:ASN:CB	2.63	0.47
4:D:40:HIS:HD2	7:G:73:LYS:HE3	1.80	0.47
4:D:198:LEU:HB2	4:D:199:ASN:H	1.53	0.47
7:G:27:LYS:HE2	7:G:54:ILE:HD12	1.97	0.47
7:G:111:THR:H	7:G:114:LEU:HD22	1.79	0.47
1:A:26:GLU:O	1:A:30:ILE:HG13	2.14	0.47
1:A:164:ARG:CG	1:A:165:GLY:N	2.78	0.47
1:A:496:GLU:O	1:A:500:GLU:HG3	2.14	0.47
1:A:532:ARG:HH12	1:A:745:GLN:NE2	2.13	0.47
1:A:647:GLY:O	1:A:648:ASN:C	2.57	0.47
1:A:857:ARG:HA	1:A:864:ILE:HD13	1.97	0.47
1:A:1308:THR:HG23	1:A:1310:GLY:N	2.26	0.47
2:B:123:THR:C	2:B:125:SER:H	2.23	0.47
2:B:204:ILE:O	2:B:204:ILE:HG22	2.13	0.47
2:B:205:ILE:O	2:B:206:ASN:C	2.57	0.47
2:B:421:PHE:CD1	2:B:424:LEU:HD23	2.50	0.47
2:B:705:MET:HA	2:B:705:MET:HE3	1.96	0.47
2:B:782:LEU:HB3	2:B:784:ASN:OD1	2.15	0.47
2:B:822:ASN:HD22	2:B:822:ASN:H	1.62	0.47
2:B:1040:ASN:O	2:B:1041:GLU:C	2.57	0.47
2:B:1076:HIS:ND1	2:B:1076:HIS:N	2.63	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:39:ALA:CA	3:C:164:ALA:HB3	2.40	0.47
3:C:44:LEU:HD13	3:C:129:ILE:HG23	1.96	0.47
3:C:46:ILE:HG22	3:C:47:ASP:N	2.29	0.47
3:C:70:ILE:HG23	3:C:142:VAL:HG11	1.96	0.47
3:C:112:ASN:HD21	3:C:146:LYS:HG2	1.79	0.47
3:C:141:GLY:O	3:C:142:VAL:HB	2.14	0.47
3:C:148:ARG:NH1	10:J:64:ASN:HA	2.30	0.47
3:C:181:ASP:CG	3:C:186:LEU:HD13	2.40	0.47
5:E:163:GLU:OE2	5:E:167:ARG:HG3	2.13	0.47
9:I:8:ARG:O	9:I:9:ASP:HB2	2.15	0.47
1:A:339:ASN:O	1:A:343:LYS:HE2	2.15	0.47
1:A:396:PRO:O	1:A:397:ASN:CG	2.58	0.47
1:A:501:LEU:HD21	2:B:1146:PHE:CE1	2.49	0.47
1:A:678:GLU:O	1:A:681:GLU:HG2	2.15	0.47
1:A:808:LEU:HD21	1:A:816:HIS:HD2	1.80	0.47
1:A:929:LEU:HD22	1:A:929:LEU:N	2.29	0.47
1:A:1074:GLU:N	1:A:1075:PRO:HD2	2.29	0.47
1:A:1139:GLU:O	1:A:1139:GLU:HG2	2.14	0.47
2:B:850:LEU:HD12	2:B:851:PHE:N	2.30	0.47
3:C:128:ASN:OD1	3:C:131:HIS:HD2	1.98	0.47
10:J:57:ILE:CG2	10:J:58:GLU:H	2.28	0.47
1:A:356:ASP:OD2	2:B:833:TYR:CE2	2.67	0.47
1:A:690:VAL:HG13	1:A:690:VAL:O	2.14	0.47
1:A:811:GLN:H	1:A:811:GLN:CD	2.23	0.47
1:A:1219:THR:HG21	1:A:1271:ILE:HD11	1.96	0.47
2:B:802:PRO:HG3	2:B:1091:TYR:CZ	2.50	0.47
2:B:821:GLN:HE22	2:B:851:PHE:H	1.63	0.47
2:B:877:PRO:C	2:B:878:GLN:HG3	2.40	0.47
4:D:15:LEU:O	4:D:17:LYS:N	2.48	0.47
4:D:53:SER:O	4:D:54:GLU:C	2.57	0.47
10:J:58:GLU:HA	10:J:61:LEU:HD12	1.97	0.47
12:L:47:ARG:NH2	12:L:54:ARG:HG2	2.30	0.47
1:A:166:GLY:O	1:A:167:CYS:O	2.33	0.46
1:A:227:VAL:HG12	1:A:228:PHE:CD1	2.51	0.46
1:A:659:HIS:ND1	2:B:1074:ASN:ND2	2.63	0.46
1:A:809:THR:HB	1:A:810:PRO:HD2	1.97	0.46
1:A:990:VAL:HG12	1:A:991:LYS:N	2.31	0.46
2:B:615:MET:HB3	2:B:626:ILE:HA	1.96	0.46
2:B:711:GLU:CB	2:B:712:PRO:HD3	2.43	0.46
2:B:899:ILE:HD12	2:B:911:ILE:HG23	1.96	0.46
2:B:916:THR:HB	2:B:935:ARG:HG3	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:999:MET:HG2	2:B:1007:VAL:HG23	1.97	0.46
2:B:1017:ILE:H	2:B:1018:PRO:CD	2.27	0.46
2:B:1065:GLN:NE2	2:B:1066:SER:H	2.12	0.46
2:B:1160:VAL:CG1	2:B:1161:HIS:N	2.73	0.46
5:E:136:ASN:OD1	5:E:137:GLU:N	2.48	0.46
5:E:202:SER:HG	5:E:204:THR:HG22	1.76	0.46
6:F:79:ARG:HG3	6:F:144:GLU:OE1	2.15	0.46
7:G:94:CYS:SG	7:G:130:TYR:CE1	3.08	0.46
8:H:63:LEU:HG	8:H:90:ALA:HB3	1.97	0.46
8:H:130:ARG:CZ	8:H:130:ARG:HB2	2.44	0.46
1:A:54:ASN:HB3	1:A:247:ARG:HH12	1.80	0.46
1:A:316:GLN:O	1:A:317:LYS:C	2.58	0.46
1:A:357:PRO:HD2	2:B:833:TYR:CE2	2.51	0.46
1:A:451:HIS:O	1:A:452:LYS:C	2.57	0.46
1:A:913:LEU:HB2	1:A:1032:LEU:HD22	1.97	0.46
1:A:1436:ILE:O	1:A:1438:THR:N	2.47	0.46
1:A:1450:LEU:HD13	6:F:131:PRO:HG2	1.97	0.46
2:B:533:CYS:C	2:B:535:LEU:N	2.72	0.46
2:B:641:GLU:HB2	2:B:643:ASP:CG	2.40	0.46
2:B:739:THR:HB	2:B:740:HIS:ND1	2.29	0.46
2:B:1065:GLN:CD	2:B:1066:SER:H	2.23	0.46
3:C:175:ALA:HB2	10:J:10:CYS:CB	2.46	0.46
3:C:245:VAL:O	3:C:246:ARG:C	2.58	0.46
4:D:159:THR:O	4:D:162:ALA:HB3	2.15	0.46
5:E:83:CYS:HB3	5:E:85:GLU:HG3	1.98	0.46
5:E:153:HIS:HB3	5:E:196:VAL:CG2	2.45	0.46
5:E:156:LEU:HD12	5:E:195:VAL:HB	1.96	0.46
6:F:127:GLU:O	6:F:128:LYS:C	2.57	0.46
7:G:115:MET:SD	7:G:119:LEU:HD23	2.56	0.46
1:A:986:ILE:HD13	1:A:1031:VAL:HG21	1.97	0.46
1:A:1102:LYS:O	1:A:1103:GLU:C	2.57	0.46
1:A:1215:ARG:O	1:A:1216:ILE:C	2.58	0.46
1:A:1300:LYS:C	1:A:1302:PRO:HD3	2.40	0.46
1:A:1313:LEU:HB3	1:A:1338:VAL:HG21	1.96	0.46
2:B:132:VAL:HG23	2:B:165:VAL:HB	1.96	0.46
2:B:800:GLN:HB3	10:J:52:THR:HG21	1.96	0.46
3:C:238:ILE:HG23	3:C:239:PRO:HD2	1.97	0.46
11:K:32:VAL:HG23	11:K:74:ARG:CG	2.43	0.46
1:A:583:PRO:HB2	1:A:637:LYS:HE3	1.98	0.46
1:A:682:THR:O	1:A:682:THR:HG22	2.15	0.46
1:A:814:PHE:HE1	2:B:519:TRP:CB	2.26	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:868:TYR:HE2	1:A:1366:ARG:NE	2.14	0.46
2:B:63:ILE:HG23	2:B:63:ILE:O	2.14	0.46
2:B:129:PHE:CE2	2:B:166:PHE:HD1	2.32	0.46
2:B:226:PHE:HA	2:B:395:GLN:CG	2.45	0.46
2:B:405:ARG:NH1	2:B:632:ARG:CG	2.79	0.46
2:B:565:PRO:C	2:B:567:GLU:N	2.73	0.46
2:B:579:ARG:HB2	2:B:586:TRP:NE1	2.30	0.46
2:B:596:LEU:O	2:B:600:LEU:HG	2.15	0.46
2:B:856:PHE:N	2:B:856:PHE:HD1	2.14	0.46
3:C:173:ALA:O	3:C:174:ALA:HB3	2.14	0.46
3:C:191:TYR:HB3	3:C:201:TRP:CD1	2.51	0.46
5:E:31:THR:OG1	5:E:34:GLU:HB2	2.16	0.46
5:E:80:VAL:HG12	5:E:82:PHE:CE1	2.50	0.46
6:F:101:ILE:HA	6:F:105:ALA:HB3	1.96	0.46
8:H:143:LEU:HD12	8:H:143:LEU:N	2.30	0.46
10:J:1:MET:N	10:J:56:LEU:N	2.63	0.46
11:K:5:ASP:O	11:K:6:ARG:C	2.59	0.46
1:A:66:LYS:O	1:A:67:CYS:CB	2.63	0.46
1:A:245:PRO:O	2:B:1114:LEU:HD11	2.15	0.46
1:A:533:LYS:HB3	1:A:533:LYS:HE2	1.69	0.46
1:A:566:ILE:HG22	1:A:566:ILE:O	2.15	0.46
1:A:795:GLU:CD	1:A:795:GLU:N	2.71	0.46
1:A:1308:THR:CG2	1:A:1310:GLY:O	2.61	0.46
1:A:1422:ARG:HH12	2:B:1224:PHE:HD2	1.62	0.46
2:B:56:ASP:HB3	2:B:57:TYR:CD1	2.51	0.46
2:B:244:LEU:HD21	2:B:366:GLN:NE2	2.31	0.46
2:B:651:LEU:C	2:B:653:VAL:H	2.24	0.46
2:B:858:SER:HA	2:B:966:VAL:O	2.15	0.46
3:C:39:ALA:HA	3:C:164:ALA:CB	2.41	0.46
3:C:123:ASN:HD21	3:C:125:MET:HG2	1.81	0.46
7:G:139:ILE:HG23	7:G:140:LYS:N	2.31	0.46
8:H:107:VAL:O	8:H:108:SER:HB2	2.16	0.46
1:A:382:PRO:HD3	1:A:428:TYR:CZ	2.50	0.46
1:A:503:GLN:O	1:A:504:LEU:HD12	2.15	0.46
1:A:541:ILE:HG22	1:A:542:GLU:H	1.81	0.46
1:A:898:ARG:HA	1:A:933:TYR:CE1	2.50	0.46
1:A:964:ILE:O	1:A:964:ILE:CG2	2.64	0.46
1:A:1097:GLY:O	1:A:1098:VAL:C	2.59	0.46
1:A:1116:LEU:HB2	1:A:1329:THR:OG1	2.15	0.46
2:B:191:LYS:HE2	2:B:191:LYS:HB3	1.78	0.46
2:B:241:ARG:HG2	2:B:253:THR:CG2	2.46	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:248:SER:O	2:B:249:ARG:C	2.59	0.46
2:B:435:THR:C	2:B:437:GLU:H	2.24	0.46
2:B:611:PRO:CB	2:B:685:LEU:HD11	2.44	0.46
2:B:744:HIS:CD2	2:B:745:PRO:HD2	2.51	0.46
2:B:800:GLN:HB2	2:B:821:GLN:HA	1.98	0.46
2:B:834:ASN:O	2:B:1013:ASN:CB	2.63	0.46
2:B:1004:GLU:HB3	2:B:1006:ILE:HD11	1.97	0.46
3:C:114:TYR:C	3:C:116:LYS:H	2.24	0.46
5:E:3:GLN:NE2	5:E:52:ARG:HH12	2.13	0.46
5:E:121:MET:C	5:E:123:LEU:H	2.23	0.46
5:E:177:ARG:HH11	5:E:215:MET:HE2	1.79	0.46
5:E:191:LYS:HB3	5:E:191:LYS:HE2	1.69	0.46
8:H:130:ARG:HA	8:H:133:ASN:HB2	1.97	0.46
10:J:17:LYS:O	10:J:18:TRP:C	2.56	0.46
1:A:58:LEU:HD12	1:A:244:PRO:HG3	1.98	0.46
1:A:92:HIS:HD2	1:A:236:LEU:HD21	1.81	0.46
1:A:535:THR:CG2	1:A:617:VAL:HG23	2.46	0.46
1:A:645:LEU:HD11	1:A:649:ILE:HD11	1.97	0.46
1:A:692:ASP:O	1:A:694:THR:N	2.49	0.46
1:A:881:GLN:NE2	1:A:958:VAL:O	2.49	0.46
1:A:1159:ARG:O	1:A:1160:SER:HB3	2.16	0.46
1:A:1280:GLU:O	1:A:1281:ARG:C	2.58	0.46
1:A:1282:VAL:HA	1:A:1307:GLU:O	2.16	0.46
2:B:706:GLN:C	2:B:708:GLU:H	2.24	0.46
2:B:861:ASP:OD1	2:B:862:GLN:N	2.48	0.46
3:C:124:LEU:C	3:C:126:GLY:N	2.73	0.46
5:E:198:ILE:CD1	5:E:212:ARG:HH11	2.27	0.46
1:A:69:THR:HG21	2:B:1174:LYS:HZ3	1.81	0.46
1:A:95:PHE:CE2	1:A:1410:PHE:CD2	3.03	0.46
1:A:521:MET:O	1:A:646:PHE:CE2	2.68	0.46
1:A:898:ARG:O	1:A:1029:ARG:NH1	2.49	0.46
1:A:901:LEU:HG	1:A:926:GLN:HE21	1.80	0.46
1:A:1437:GLY:C	1:A:1439:GLY:H	2.23	0.46
2:B:103:ASN:HB2	2:B:169:ARG:NH2	2.31	0.46
2:B:114:PRO:O	2:B:115:GLN:C	2.57	0.46
2:B:294:ASP:C	2:B:296:GLU:N	2.74	0.46
2:B:887:HIS:CD2	2:B:888:GLY:N	2.84	0.46
3:C:167:HIS:HA	11:K:6:ARG:NH1	2.30	0.46
5:E:121:MET:HA	5:E:124:VAL:HG23	1.97	0.46
5:E:131:THR:C	5:E:132:ILE:HG13	2.41	0.46
6:F:70:LYS:HD3	6:F:70:LYS:HA	1.42	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:F:89:GLU:HB3	6:F:134:ILE:CD1	2.45	0.46
8:H:38:LEU:HD12	8:H:39:THR:N	2.31	0.46
11:K:65:HIS:CD2	11:K:66:PRO:N	2.84	0.46
1:A:56:PRO:O	1:A:57:ARG:CG	2.60	0.46
1:A:222:LEU:O	1:A:222:LEU:HD13	2.16	0.46
1:A:381:THR:HG23	6:F:104:ASN:OD1	2.16	0.46
1:A:451:HIS:CD2	1:A:1074:GLU:HG3	2.51	0.46
1:A:923:LEU:HD23	1:A:923:LEU:HA	1.54	0.46
1:A:952:ALA:C	1:A:954:TRP:HD1	2.23	0.46
1:A:1167:GLU:HA	1:A:1170:ILE:CD1	2.45	0.46
1:A:1215:ARG:O	1:A:1218:GLN:N	2.49	0.46
1:A:1265:ASN:C	1:A:1267:MET:N	2.73	0.46
2:B:698:GLU:HG2	2:B:701:ILE:HD11	1.98	0.46
2:B:859:TYR:CZ	2:B:941:LEU:HD12	2.50	0.46
2:B:898:LEU:HD13	2:B:952:VAL:HG11	1.98	0.46
2:B:906:SER:HA	2:B:946:ASN:HB3	1.97	0.46
2:B:1013:ASN:HD21	2:B:1015:HIS:CD2	2.31	0.46
4:D:14:ARG:O	4:D:16:LYS:N	2.43	0.46
5:E:168:TYR:HD2	5:E:168:TYR:HA	1.69	0.46
7:G:21:ARG:O	7:G:22:MET:C	2.59	0.46
7:G:95:SER:O	7:G:121:PHE:CE1	2.68	0.46
8:H:38:LEU:HD13	8:H:125:LEU:HD12	1.97	0.46
11:K:55:LYS:HB2	11:K:81:TYR:HE1	1.81	0.46
1:A:41:MET:CE	1:A:257:ARG:HG3	2.46	0.46
1:A:340:LEU:HD22	2:B:1200:ALA:HB2	1.97	0.46
1:A:456:MET:CE	1:A:507:VAL:HG23	2.46	0.46
1:A:805:LEU:HD13	2:B:1052:VAL:HG21	1.98	0.46
1:A:850:VAL:HG12	1:A:850:VAL:O	2.16	0.46
1:A:1228:TRP:CE3	1:A:1228:TRP:N	2.84	0.46
2:B:241:ARG:HG2	2:B:253:THR:HG21	1.97	0.46
2:B:308:TRP:HH2	9:I:45:ARG:HG2	1.81	0.46
2:B:446:LEU:O	2:B:447:ALA:CB	2.64	0.46
2:B:1095:LEU:O	2:B:1096:ARG:O	2.34	0.46
3:C:131:HIS:HA	3:C:132:PRO:HD2	1.76	0.46
3:C:167:HIS:ND1	3:C:167:HIS:C	2.74	0.46
3:C:251:LEU:O	3:C:255:VAL:HG23	2.15	0.46
4:D:208:GLU:CG	4:D:212:LYS:HE3	2.43	0.46
5:E:61:GLN:HB2	5:E:79:TRP:HE3	1.81	0.46
5:E:132:ILE:HD12	5:E:132:ILE:N	2.31	0.46
5:E:178:ILE:O	5:E:178:ILE:HG23	2.16	0.46
11:K:58:PHE:HE2	11:K:74:ARG:HB3	1.80	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:445:ASN:HA	1:A:478:TYR:HE2	1.81	0.45
1:A:1267:MET:O	1:A:1271:ILE:HB	2.15	0.45
2:B:802:PRO:HG3	2:B:1091:TYR:CE1	2.51	0.45
2:B:1072:MET:HG3	2:B:1085:ILE:HD13	1.99	0.45
3:C:180:TYR:O	3:C:181:ASP:C	2.59	0.45
3:C:248:ILE:HD13	11:K:101:LEU:HD22	1.99	0.45
7:G:122:ASN:ND2	7:G:125:SER:HB3	2.30	0.45
1:A:135:PHE:CD1	1:A:222:LEU:HD21	2.51	0.45
1:A:203:SER:O	1:A:206:GLU:HB3	2.17	0.45
1:A:528:LEU:CD1	1:A:749:ALA:HB1	2.46	0.45
1:A:529:CYS:SG	1:A:662:PHE:CZ	3.10	0.45
1:A:545:GLN:O	1:A:546:VAL:C	2.59	0.45
1:A:670:ILE:HG23	1:A:805:LEU:HD21	1.99	0.45
1:A:829:VAL:C	1:A:831:THR:N	2.74	0.45
1:A:1349:TYR:HE2	1:A:1353:TYR:CD1	2.34	0.45
2:B:424:LEU:O	2:B:428:ILE:HG13	2.16	0.45
2:B:515:HIS:CD2	2:B:516:ASN:N	2.83	0.45
2:B:758:PHE:N	2:B:759:PRO:CD	2.80	0.45
2:B:844:SER:O	2:B:848:ARG:HG3	2.15	0.45
2:B:1058:LEU:O	2:B:1061:GLU:HB2	2.15	0.45
7:G:44:TYR:CE1	7:G:157:ILE:HB	2.51	0.45
7:G:119:LEU:HA	7:G:132:SER:OG	2.16	0.45
8:H:32:THR:HG22	8:H:33:GLN:OE1	2.16	0.45
11:K:30:ALA:HB2	11:K:76:GLN:HG3	1.97	0.45
1:A:14:VAL:H	1:A:1432:GLN:NE2	2.00	0.45
1:A:19:PHE:HZ	1:A:1397:LEU:CD2	2.22	0.45
1:A:261:ASP:OD2	1:A:261:ASP:N	2.48	0.45
1:A:391:LEU:O	1:A:392:VAL:C	2.60	0.45
1:A:445:ASN:HA	1:A:478:TYR:CE2	2.51	0.45
1:A:925:LEU:O	1:A:929:LEU:HD23	2.17	0.45
1:A:1229:SER:HB2	1:A:1230:GLU:H	1.48	0.45
2:B:27:ALA:O	2:B:28:GLU:C	2.59	0.45
2:B:102:VAL:O	2:B:109:THR:HG23	2.17	0.45
2:B:496:ARG:NH1	2:B:541:LEU:HA	2.31	0.45
2:B:599:THR:HG22	2:B:600:LEU:HD23	1.99	0.45
2:B:797:TYR:HD2	2:B:798:TYR:HE2	1.64	0.45
2:B:1016:ALA:CB	2:B:1020:ARG:HE	2.29	0.45
3:C:88:CYS:O	3:C:89:GLU:C	2.58	0.45
4:D:172:LEU:HD11	4:D:202:ILE:HG21	1.97	0.45
5:E:135:PHE:HB3	5:E:140:LEU:HD11	1.99	0.45
5:E:139:ALA:O	5:E:140:LEU:HD23	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:F:127:GLU:O	6:F:129:LYS:N	2.49	0.45
12:L:36:SER:O	12:L:37:LYS:C	2.60	0.45
12:L:38:LEU:CG	12:L:39:SER:H	2.29	0.45
1:A:135:PHE:HD1	1:A:222:LEU:HD21	1.80	0.45
1:A:303:TYR:O	1:A:325:ILE:HG13	2.16	0.45
1:A:689:LYS:O	1:A:689:LYS:HG2	2.16	0.45
1:A:833:GLU:O	1:A:837:ILE:HG13	2.16	0.45
2:B:243:ALA:O	2:B:244:LEU:C	2.59	0.45
2:B:259:TYR:N	2:B:259:TYR:CD1	2.85	0.45
2:B:542:MET:HB3	2:B:636:PRO:HD3	1.98	0.45
2:B:912:ILE:HD11	2:B:966:VAL:HG21	1.97	0.45
2:B:1070:GLU:O	2:B:1084:GLN:HB3	2.15	0.45
2:B:1076:HIS:CD2	11:K:40:HIS:CE1	2.90	0.45
6:F:125:LEU:O	6:F:128:LYS:N	2.48	0.45
6:F:130:ILE:HB	6:F:148:VAL:HG21	1.99	0.45
11:K:59:ALA:HA	11:K:74:ARG:O	2.15	0.45
12:L:40:LEU:HD22	12:L:44:ASP:OD1	2.17	0.45
1:A:58:LEU:HD12	1:A:244:PRO:CG	2.47	0.45
1:A:239:LEU:CD1	1:A:240:PRO:HD2	2.36	0.45
1:A:380:VAL:HG23	1:A:430:TRP:C	2.41	0.45
1:A:627:GLY:O	1:A:628:GLY:C	2.60	0.45
1:A:929:LEU:CD2	1:A:929:LEU:N	2.79	0.45
1:A:1129:GLU:O	1:A:1133:LEU:HG	2.16	0.45
1:A:1135:ARG:HG2	1:A:1136:SER:N	2.32	0.45
1:A:1138:ILE:HA	1:A:1275:GLY:HA2	1.99	0.45
1:A:1365:TYR:CD1	5:E:203:GLU:HG2	2.52	0.45
2:B:602:THR:HA	2:B:605:ARG:HB2	1.98	0.45
2:B:705:MET:HE1	2:B:742:GLU:HG2	1.98	0.45
2:B:779:GLY:O	2:B:795:ILE:HA	2.17	0.45
3:C:121:VAL:O	3:C:122:SER:C	2.59	0.45
3:C:130:GLY:O	3:C:132:PRO:HD3	2.17	0.45
3:C:175:ALA:HB2	10:J:10:CYS:HB2	1.97	0.45
5:E:96:PHE:CZ	5:E:100:ILE:HD11	2.51	0.45
7:G:13:LEU:HD13	7:G:22:MET:HE3	1.98	0.45
8:H:15:VAL:HG13	8:H:24:CYS:HB3	1.98	0.45
1:A:444:PHE:CE2	1:A:470:LEU:HD13	2.51	0.45
1:A:590:ARG:O	1:A:591:PHE:CB	2.57	0.45
1:A:808:LEU:HD23	1:A:812:GLU:C	2.42	0.45
1:A:1004:ASN:HD22	1:A:1007:ILE:HG13	1.82	0.45
1:A:1061:GLY:O	1:A:1062:GLU:C	2.59	0.45
1:A:1280:GLU:HB3	1:A:1281:ARG:H	1.61	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1395:GLY:HA3	1:A:1426:GLU:OE2	2.16	0.45
2:B:312:GLU:O	2:B:313:MET:C	2.60	0.45
2:B:839:MET:HG3	2:B:1010:LEU:HD21	1.98	0.45
2:B:887:HIS:CD2	2:B:888:GLY:H	2.34	0.45
3:C:58:LEU:HD13	3:C:62:PHE:CD2	2.51	0.45
7:G:1:MET:SD	7:G:79:PHE:CE1	3.10	0.45
8:H:4:THR:HG21	8:H:7:ASP:OD2	2.16	0.45
8:H:35:GLN:H	8:H:35:GLN:HG3	1.57	0.45
8:H:47:PHE:CD2	8:H:95:TYR:HD1	2.34	0.45
10:J:17:LYS:HB3	10:J:39:LEU:HD23	1.98	0.45
10:J:56:LEU:O	10:J:60:PHE:CD2	2.70	0.45
1:A:68:GLN:O	1:A:68:GLN:CD	2.59	0.45
1:A:152:VAL:CG1	1:A:153:PRO:HD2	2.47	0.45
1:A:252:PHE:O	1:A:253:ASN:CB	2.65	0.45
1:A:456:MET:HE3	1:A:507:VAL:HG23	1.98	0.45
1:A:763:ALA:O	1:A:803:SER:HB3	2.17	0.45
1:A:868:TYR:O	1:A:869:GLY:C	2.59	0.45
2:B:234:ILE:CG1	2:B:257:LYS:HB3	2.45	0.45
2:B:666:TYR:C	2:B:668:ASP:H	2.25	0.45
2:B:811:TYR:CD1	2:B:811:TYR:N	2.83	0.45
2:B:841:MET:CE	2:B:1010:LEU:HG	2.46	0.45
2:B:918:ILE:HG21	2:B:935:ARG:NH2	2.32	0.45
2:B:942:ARG:HD3	2:B:942:ARG:HA	1.67	0.45
2:B:956:THR:HA	2:B:961:LEU:O	2.17	0.45
4:D:146:GLN:O	4:D:149:THR:HG22	2.17	0.45
5:E:18:THR:O	5:E:19:VAL:C	2.59	0.45
7:G:14:HIS:HD2	7:G:16:SER:H	1.63	0.45
7:G:132:SER:O	7:G:133:SER:C	2.58	0.45
12:L:39:SER:O	12:L:40:LEU:CG	2.57	0.45
1:A:447:GLN:HE21	2:B:1134:GLU:CD	2.24	0.45
1:A:578:LEU:C	1:A:580:VAL:N	2.74	0.45
1:A:598:LEU:HD13	8:H:115:TYR:HE2	1.81	0.45
1:A:897:TYR:HD2	1:A:936:LEU:CD1	2.30	0.45
1:A:1189:SER:C	1:A:1191:TRP:H	2.25	0.45
1:A:1279:ILE:HG22	1:A:1282:VAL:CG2	2.47	0.45
2:B:102:VAL:HG23	2:B:112:LEU:HD22	1.97	0.45
2:B:1193:GLN:C	2:B:1194:ILE:CG2	2.90	0.45
4:D:29:LEU:HB3	4:D:33:PHE:HB2	1.98	0.45
4:D:39:ASN:ND2	4:D:40:HIS:N	2.64	0.45
4:D:52:LEU:H	4:D:182:SER:HB3	1.82	0.45
11:K:47:ARG:NH1	11:K:51:LEU:HD22	2.32	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:55:ASP:HB2	1:A:59:GLY:O	2.17	0.45
1:A:527:THR:O	1:A:528:LEU:C	2.60	0.45
1:A:1173:HIS:CB	1:A:1227:ILE:HD13	2.47	0.45
1:A:1188:GLN:OE1	1:A:1241:ARG:HD3	2.17	0.45
2:B:66:ASP:OD1	2:B:422:LYS:HE2	2.17	0.45
2:B:199:MET:HE2	2:B:492:LEU:HD23	1.98	0.45
2:B:291:ILE:HG21	2:B:300:HIS:CD2	2.52	0.45
2:B:398:ARG:NH1	2:B:398:ARG:CB	2.80	0.45
2:B:950:ASP:HB2	2:B:969:ARG:HB3	1.99	0.45
3:C:67:LEU:HA	3:C:70:ILE:HD11	1.97	0.45
3:C:241:ASP:O	3:C:245:VAL:HG23	2.17	0.45
4:D:192:LYS:HD2	4:D:199:ASN:HA	1.98	0.45
5:E:79:TRP:HE1	5:E:81:GLU:HB2	1.82	0.45
6:F:85:MET:SD	6:F:90:ARG:HB2	2.57	0.45
7:G:79:PHE:C	7:G:79:PHE:HD1	2.22	0.45
7:G:79:PHE:HD1	7:G:80:LYS:N	2.14	0.45
7:G:106:MET:HG3	7:G:157:ILE:O	2.16	0.45
11:K:56:VAL:O	11:K:56:VAL:CG1	2.64	0.45
1:A:414:ASP:C	1:A:416:ARG:H	2.25	0.45
1:A:709:THR:CG2	1:A:710:LEU:N	2.79	0.45
1:A:899:VAL:CG1	1:A:908:LEU:HD21	2.43	0.45
1:A:1313:LEU:O	1:A:1314:SER:C	2.59	0.45
2:B:416:LEU:HD11	2:B:460:ALA:HB3	1.98	0.45
2:B:710:LEU:HA	2:B:733:HIS:HB3	1.98	0.45
2:B:948:ILE:HD12	2:B:969:ARG:HH11	1.81	0.45
2:B:1013:ASN:ND2	2:B:1015:HIS:HD2	2.13	0.45
2:B:1109:GLY:H	2:B:1111:MET:HE2	1.82	0.45
4:D:129:LEU:O	4:D:132:GLN:HG3	2.17	0.45
4:D:151:PHE:HD1	4:D:151:PHE:N	2.15	0.45
7:G:163:ILE:O	7:G:163:ILE:HG13	2.17	0.45
10:J:32:GLU:H	10:J:32:GLU:CD	2.24	0.45
1:A:58:LEU:HD13	1:A:244:PRO:HD3	1.97	0.44
1:A:181:LEU:HD23	1:A:181:LEU:HA	1.87	0.44
1:A:474:VAL:HG23	1:A:521:MET:CE	2.47	0.44
1:A:886:ILE:HD11	1:A:943:LEU:O	2.16	0.44
1:A:1069:ALA:C	1:A:1071:SER:N	2.73	0.44
1:A:1132:LYS:HA	1:A:1135:ARG:HB3	1.99	0.44
1:A:1311:VAL:HG11	1:A:1329:THR:HG21	1.99	0.44
1:A:1445:ILE:CD1	7:G:59:GLY:H	2.29	0.44
2:B:179:CYS:O	2:B:181:LEU:N	2.50	0.44
2:B:218:SER:O	2:B:219:ALA:C	2.60	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:219:ALA:HA	2:B:405:ARG:HG2	1.98	0.44
2:B:360:PHE:CZ	2:B:361:LEU:HD13	2.52	0.44
2:B:483:LEU:HG	2:B:485:ARG:HD3	1.98	0.44
2:B:842:ASN:C	2:B:842:ASN:ND2	2.75	0.44
2:B:1037:LEU:HD12	2:B:1037:LEU:N	2.32	0.44
2:B:1073:TYR:OH	3:C:179:GLU:HG3	2.17	0.44
2:B:1109:GLY:HA3	2:B:1110:PRO:HD3	1.80	0.44
4:D:35:LEU:HD12	4:D:35:LEU:N	2.25	0.44
4:D:185:CYS:HB3	4:D:211:LEU:HD22	1.99	0.44
5:E:187:TYR:CD2	5:E:188:LEU:HD23	2.50	0.44
6:F:101:ILE:CG2	6:F:117:PRO:HB3	2.46	0.44
7:G:129:SER:OG	7:G:130:TYR:N	2.49	0.44
11:K:23:PRO:HA	11:K:31:VAL:HG22	1.98	0.44
1:A:93:VAL:HG11	1:A:305:ASP:HB3	1.99	0.44
1:A:434:ARG:HH21	1:A:437:MET:H	1.65	0.44
1:A:886:ILE:HG13	1:A:943:LEU:HB2	1.98	0.44
1:A:1196:GLU:HG3	1:A:1197:LEU:N	2.31	0.44
1:A:1396:ALA:HB2	1:A:1417:GLU:OE1	2.17	0.44
2:B:101:MET:HB2	2:B:169:ARG:HH12	1.82	0.44
2:B:241:ARG:HH21	2:B:251:ILE:HD13	1.82	0.44
2:B:1098:MET:HE3	2:B:1101:ASP:OD2	2.17	0.44
4:D:52:LEU:C	4:D:53:SER:HG	2.22	0.44
8:H:5:LEU:O	8:H:6:PHE:HB2	2.17	0.44
11:K:63:VAL:O	11:K:63:VAL:CG2	2.61	0.44
1:A:27:VAL:O	1:A:29:ALA:N	2.50	0.44
1:A:434:ARG:HH21	1:A:437:MET:HG3	1.80	0.44
1:A:664:THR:O	1:A:664:THR:CG2	2.64	0.44
1:A:868:TYR:O	1:A:870:GLU:N	2.50	0.44
1:A:1153:TYR:CZ	1:A:1163:ILE:HD11	2.53	0.44
1:A:1290:LYS:HA	1:A:1299:VAL:O	2.17	0.44
2:B:244:LEU:HD12	2:B:250:PHE:HB2	1.99	0.44
2:B:254:LEU:HD22	2:B:361:LEU:CD1	2.47	0.44
2:B:519:TRP:CD1	2:B:519:TRP:C	2.95	0.44
2:B:582:VAL:HG22	2:B:626:ILE:HB	1.99	0.44
2:B:750:GLY:H	2:B:753:ALA:HB3	1.80	0.44
2:B:766:ARG:HH21	2:B:1020:ARG:HB3	1.81	0.44
2:B:1076:HIS:CD2	11:K:40:HIS:CD2	3.05	0.44
3:C:129:ILE:HG22	3:C:129:ILE:O	2.16	0.44
4:D:173:HIS:C	4:D:175:PHE:H	2.25	0.44
5:E:67:GLU:O	5:E:70:SER:HB3	2.17	0.44
5:E:145:THR:HG21	5:E:187:TYR:CE2	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:H:82:PRO:HB2	8:H:83:GLN:H	1.67	0.44
11:K:5:ASP:CB	11:K:7:PHE:CZ	2.96	0.44
11:K:114:LEU:O	11:K:114:LEU:HD23	2.17	0.44
1:A:409:SER:O	1:A:410:GLY:C	2.59	0.44
1:A:671:ALA:HB3	1:A:676:MET:HG3	1.99	0.44
1:A:1356:ILE:HG21	1:A:1363:VAL:HG23	1.98	0.44
1:A:1435:PRO:HA	1:A:1439:GLY:O	2.18	0.44
2:B:866:TYR:HD2	2:B:870:ILE:HD12	1.82	0.44
2:B:980:PHE:CD1	2:B:980:PHE:N	2.86	0.44
2:B:1003:ALA:HB1	3:C:179:GLU:HB2	1.99	0.44
2:B:1102:LYS:C	2:B:1122:ARG:HH11	2.25	0.44
2:B:1185:CYS:SG	4:D:17:LYS:HE3	2.57	0.44
2:B:1204:PHE:CZ	2:B:1216:LEU:HD21	2.52	0.44
3:C:258:ILE:HG13	11:K:19:LEU:HD11	1.98	0.44
4:D:53:SER:C	4:D:55:ALA:N	2.74	0.44
4:D:56:ARG:HB2	4:D:148:LEU:HB3	1.99	0.44
4:D:185:CYS:SG	4:D:186:ASP:N	2.90	0.44
4:D:203:SER:OG	4:D:206:GLU:HB2	2.17	0.44
5:E:98:ILE:HA	5:E:101:GLN:HB3	1.99	0.44
5:E:150:VAL:HG12	5:E:151:PRO:HD2	1.98	0.44
12:L:61:THR:HG21	12:L:63:ARG:NH2	2.32	0.44
1:A:88:LYS:HE2	1:A:205:GLU:OE2	2.17	0.44
1:A:100:LYS:O	1:A:104:GLU:HG3	2.18	0.44
1:A:523:ILE:HG22	1:A:528:LEU:HB3	1.99	0.44
1:A:598:LEU:HA	8:H:122:LEU:CD1	2.46	0.44
1:A:670:ILE:HG22	1:A:805:LEU:HD21	1.98	0.44
1:A:861:GLY:O	1:A:862:ASN:C	2.60	0.44
1:A:1072:ILE:C	1:A:1075:PRO:HD2	2.43	0.44
2:B:54:PHE:HB2	2:B:410:GLY:HA2	1.99	0.44
2:B:362:PRO:O	2:B:366:GLN:HG3	2.17	0.44
2:B:552:MET:HA	2:B:552:MET:HE3	1.99	0.44
3:C:192:TRP:O	3:C:192:TRP:CE3	2.71	0.44
4:D:26:THR:C	4:D:28:GLN:N	2.76	0.44
4:D:202:ILE:HD11	4:D:207:LEU:N	2.32	0.44
12:L:40:LEU:HD22	12:L:44:ASP:CG	2.42	0.44
1:A:839:ARG:O	1:A:840:ARG:C	2.60	0.44
1:A:1232:ASN:O	1:A:1233:ASP:C	2.60	0.44
2:B:515:HIS:CG	2:B:516:ASN:H	2.34	0.44
2:B:1076:HIS:CD2	11:K:40:HIS:NE2	2.78	0.44
4:D:66:ARG:HG3	7:G:51:TYR:CD1	2.53	0.44
5:E:187:TYR:C	5:E:188:LEU:HD23	2.43	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:F:81:THR:O	6:F:82:THR:C	2.60	0.44
7:G:99:PHE:HE1	7:G:101:VAL:HG23	1.82	0.44
8:H:94:ASP:O	8:H:95:TYR:HB2	2.17	0.44
12:L:55:ILE:CG1	12:L:56:LEU:H	2.31	0.44
1:A:205:GLU:CD	1:A:205:GLU:N	2.74	0.44
1:A:329:LEU:O	1:A:330:LYS:O	2.36	0.44
1:A:567:LYS:CE	8:H:46:LEU:HB2	2.48	0.44
1:A:628:GLY:O	1:A:629:LEU:C	2.59	0.44
2:B:879:ARG:HE	2:B:879:ARG:CA	2.31	0.44
2:B:1034:VAL:HA	2:B:1037:LEU:HD12	1.99	0.44
5:E:89:GLY:C	5:E:91:LYS:N	2.76	0.44
5:E:117:THR:HB	5:E:120:ALA:CB	2.46	0.44
5:E:184:VAL:O	5:E:187:TYR:HB3	2.18	0.44
8:H:81:PRO:O	8:H:82:PRO:O	2.36	0.44
11:K:61:TYR:O	11:K:61:TYR:CG	2.69	0.44
1:A:310:GLY:O	1:A:312:PRO:HD2	2.18	0.44
1:A:325:ILE:C	1:A:327:ALA:H	2.25	0.44
1:A:595:THR:HG23	1:A:603:ASN:HB2	2.00	0.44
1:A:897:TYR:CD2	1:A:936:LEU:CD1	3.01	0.44
1:A:1163:ILE:HG23	1:A:1164:PRO:HD2	1.99	0.44
1:A:1167:GLU:HA	1:A:1170:ILE:HD12	1.99	0.44
1:A:1239:ARG:C	1:A:1240:CYS:SG	3.00	0.44
2:B:759:PRO:C	2:B:761:HIS:N	2.76	0.44
2:B:885:MET:HA	2:B:936:ASP:HB2	2.00	0.44
2:B:911:ILE:HD11	2:B:941:LEU:HB2	2.00	0.44
2:B:1080:LYS:HG3	3:C:180:TYR:OH	2.18	0.44
6:F:69:LEU:C	6:F:71:GLU:H	2.25	0.44
7:G:165:GLU:H	7:G:165:GLU:HG2	1.37	0.44
9:I:7:CYS:HB2	9:I:34:TYR:CD2	2.53	0.44
1:A:35:ILE:H	1:A:35:ILE:CD1	2.20	0.44
1:A:195:ASP:O	1:A:196:GLU:O	2.36	0.44
1:A:360:GLU:O	1:A:362:ASP:N	2.51	0.44
1:A:494:SER:O	1:A:495:GLU:C	2.61	0.44
1:A:633:VAL:O	1:A:636:GLU:N	2.51	0.44
1:A:809:THR:HG21	2:B:730:ARG:HG3	2.00	0.44
1:A:858:ASN:HD22	1:A:858:ASN:N	2.11	0.44
1:A:1215:ARG:HA	1:A:1215:ARG:HD2	1.80	0.44
2:B:121:ASN:HA	2:B:207:GLY:CA	2.48	0.44
2:B:179:CYS:C	2:B:181:LEU:H	2.25	0.44
2:B:430:ARG:HH11	2:B:430:ARG:CB	2.30	0.44
2:B:1192:TYR:CG	2:B:1218:THR:HG21	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:59:ALA:O	3:C:60:ASP:C	2.59	0.44
3:C:126:GLY:C	3:C:127:ARG:HG2	2.43	0.44
4:D:14:ARG:C	4:D:16:LYS:H	2.24	0.44
4:D:33:PHE:CE2	7:G:3:PHE:CD1	3.06	0.44
4:D:178:ALA:O	4:D:181:GLY:N	2.50	0.44
5:E:63:ASN:CB	5:E:64:PRO:HD2	2.47	0.44
5:E:90:VAL:O	5:E:90:VAL:CG2	2.66	0.44
7:G:14:HIS:CD2	7:G:15:PRO:CD	3.01	0.44
7:G:143:ILE:HG12	7:G:170:ALA:CA	2.43	0.44
8:H:54:SER:C	8:H:55:LEU:HD12	2.42	0.44
11:K:35:PHE:N	11:K:35:PHE:HD1	2.15	0.44
12:L:26:THR:CG2	12:L:27:LEU:H	2.23	0.44
1:A:14:VAL:HA	2:B:1217:TYR:O	2.18	0.43
1:A:211:PHE:HA	1:A:214:ILE:CD1	2.47	0.43
1:A:663:SER:HB2	2:B:1085:ILE:HG13	1.99	0.43
1:A:690:VAL:HG21	1:A:722:LEU:HD11	1.98	0.43
1:A:809:THR:OG1	1:A:812:GLU:HG3	2.18	0.43
1:A:841:LEU:CD2	1:A:845:LEU:HD11	2.48	0.43
1:A:1162:VAL:O	1:A:1162:VAL:HG12	2.18	0.43
1:A:1170:ILE:HG22	1:A:1174:PHE:HE1	1.81	0.43
1:A:1397:LEU:HA	1:A:1400:CYS:SG	2.58	0.43
1:A:1402:PHE:O	1:A:1404:GLU:N	2.51	0.43
2:B:225:VAL:HG22	2:B:396:ASP:OD2	2.18	0.43
2:B:563:MET:HE1	2:B:580:VAL:HG11	1.99	0.43
2:B:878:GLN:OE1	2:B:879:ARG:HG2	2.18	0.43
2:B:913:GLY:HA2	2:B:938:SER:OG	2.17	0.43
3:C:242:GLN:O	3:C:246:ARG:HG2	2.17	0.43
4:D:66:ARG:HD2	4:D:133:THR:HB	1.99	0.43
7:G:160:ILE:HD12	7:G:160:ILE:N	2.32	0.43
12:L:33:GLU:H	12:L:33:GLU:CD	2.26	0.43
1:A:92:HIS:CE1	1:A:1410:PHE:HE2	2.33	0.43
1:A:390:GLN:HE21	1:A:394:ASN:ND2	2.16	0.43
1:A:523:ILE:HD13	1:A:622:VAL:HG23	1.99	0.43
1:A:946:VAL:HG13	5:E:201:LYS:HB3	1.98	0.43
1:A:1009:ASN:O	1:A:1010:ALA:C	2.60	0.43
1:A:1106:ASN:O	1:A:1107:VAL:HB	2.18	0.43
1:A:1207:LEU:HD21	1:A:1274:ARG:NE	2.33	0.43
1:A:1223:ASP:HA	1:A:1243:VAL:HG22	1.97	0.43
2:B:205:ILE:O	2:B:207:GLY:N	2.51	0.43
2:B:276:ILE:O	2:B:276:ILE:HG22	2.18	0.43
2:B:360:PHE:CD2	2:B:361:LEU:HB2	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:521:LEU:HD11	2:B:695:ALA:HB2	2.00	0.43
2:B:591:ARG:O	2:B:592:ASN:C	2.61	0.43
2:B:852:ARG:HH22	12:L:70:ARG:C	2.26	0.43
2:B:1073:TYR:HE2	3:C:180:TYR:CE2	2.36	0.43
3:C:53:THR:HB	3:C:154:LYS:HB2	1.99	0.43
4:D:48:ILE:O	4:D:48:ILE:CG2	2.66	0.43
5:E:24:LYS:HG2	5:E:25:ASP:N	2.34	0.43
5:E:98:ILE:HG22	5:E:102:GLU:OE2	2.19	0.43
5:E:117:THR:C	5:E:119:SER:H	2.26	0.43
5:E:147:HIS:HD2	5:E:148:GLU:N	2.16	0.43
8:H:5:LEU:HD22	8:H:133:ASN:O	2.18	0.43
8:H:95:TYR:HE2	8:H:97:MET:CG	2.31	0.43
10:J:57:ILE:O	10:J:58:GLU:C	2.60	0.43
11:K:7:PHE:C	11:K:9:LEU:N	2.76	0.43
11:K:23:PRO:CA	11:K:31:VAL:HG13	2.46	0.43
1:A:91:PHE:HE2	1:A:99:ILE:HG21	1.82	0.43
1:A:114:LEU:O	1:A:115:LEU:HG	2.18	0.43
1:A:298:PHE:CE2	1:A:314:ALA:HB2	2.53	0.43
1:A:532:ARG:C	1:A:534:LEU:N	2.76	0.43
1:A:631:HIS:O	1:A:632:VAL:C	2.62	0.43
1:A:774:ARG:HG3	1:A:797:LYS:HE3	2.00	0.43
1:A:1001:ARG:O	1:A:1002:GLY:C	2.61	0.43
1:A:1040:GLN:O	1:A:1041:ALA:C	2.62	0.43
1:A:1332:PHE:H	1:A:1332:PHE:HD2	1.66	0.43
2:B:125:SER:HB2	2:B:169:ARG:HB3	2.00	0.43
2:B:219:ALA:H	2:B:404:LYS:HA	1.84	0.43
2:B:321:GLY:O	2:B:322:PHE:C	2.61	0.43
2:B:605:ARG:HG2	2:B:688:GLY:CA	2.48	0.43
2:B:1106:ARG:HD3	2:B:1126:GLY:C	2.43	0.43
3:C:130:GLY:O	3:C:132:PRO:CD	2.67	0.43
7:G:41:LYS:HD3	7:G:42:PHE:CZ	2.54	0.43
8:H:37:LYS:O	8:H:38:LEU:HB2	2.19	0.43
1:A:218:ASP:O	1:A:219:PHE:C	2.61	0.43
1:A:336:ILE:C	1:A:338:GLY:H	2.27	0.43
1:A:337:ARG:HG2	2:B:1132:GLU:OE2	2.19	0.43
1:A:472:LEU:HD11	2:B:835:GLN:NE2	2.33	0.43
1:A:588:LEU:HG	1:A:589:GLN:N	2.32	0.43
1:A:595:THR:HG23	1:A:603:ASN:CB	2.49	0.43
1:A:645:LEU:O	1:A:646:PHE:C	2.59	0.43
1:A:852:TYR:HD2	1:A:852:TYR:HA	1.70	0.43
1:A:1215:ARG:HA	1:A:1218:GLN:HG2	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1215:ARG:HH22	1:A:1272:THR:CB	2.31	0.43
1:A:1344:GLY:C	1:A:1346:ALA:N	2.77	0.43
2:B:179:CYS:C	2:B:181:LEU:N	2.74	0.43
2:B:510:LYS:HG3	2:B:511:PRO:CD	2.27	0.43
2:B:530:GLY:O	2:B:533:CYS:HB2	2.19	0.43
2:B:639:ILE:CG2	2:B:641:GLU:HG2	2.48	0.43
2:B:878:GLN:C	2:B:879:ARG:HG2	2.43	0.43
2:B:1191:ILE:CG2	2:B:1192:TYR:N	2.82	0.43
2:B:1191:ILE:HG22	2:B:1192:TYR:N	2.34	0.43
3:C:191:TYR:HD2	3:C:201:TRP:NE1	2.16	0.43
5:E:142:VAL:HG12	5:E:143:ASN:N	2.33	0.43
6:F:124:GLU:HG2	6:F:129:LYS:O	2.19	0.43
7:G:89:GLY:HA3	7:G:103:VAL:CG2	2.48	0.43
7:G:96:GLN:H	7:G:96:GLN:HG3	1.69	0.43
11:K:91:CYS:O	11:K:94:ILE:HB	2.19	0.43
1:A:7:SER:HB3	2:B:1193:GLN:NE2	2.33	0.43
1:A:37:PHE:N	1:A:37:PHE:CD1	2.87	0.43
1:A:117:GLU:H	1:A:117:GLU:CD	2.26	0.43
1:A:284:ALA:C	1:A:286:HIS:H	2.27	0.43
1:A:284:ALA:C	1:A:286:HIS:N	2.77	0.43
1:A:995:GLU:O	1:A:995:GLU:CD	2.61	0.43
1:A:1114:PRO:O	1:A:1311:VAL:HG23	2.19	0.43
2:B:209:GLU:O	2:B:482:VAL:HG12	2.17	0.43
2:B:486:TYR:HD1	2:B:1096:ARG:HH22	1.65	0.43
2:B:514:LEU:HD12	2:B:518:HIS:CD2	2.53	0.43
2:B:516:ASN:H	2:B:516:ASN:ND2	2.16	0.43
2:B:852:ARG:HG2	2:B:973:ILE:HG23	2.00	0.43
2:B:1130:PHE:CE2	2:B:1150:ARG:HG2	2.54	0.43
3:C:9:LYS:O	3:C:20:PHE:HB2	2.18	0.43
3:C:62:PHE:O	3:C:66:ARG:HG2	2.18	0.43
3:C:88:CYS:SG	3:C:91:HIS:C	3.01	0.43
3:C:259:LEU:O	3:C:262:LEU:HB2	2.19	0.43
4:D:20:GLU:O	4:D:20:GLU:HG2	2.18	0.43
4:D:71:LYS:O	4:D:74:GLN:HB2	2.19	0.43
4:D:163:VAL:O	4:D:166:LEU:HB3	2.19	0.43
5:E:82:PHE:CD1	5:E:82:PHE:N	2.87	0.43
5:E:138:ALA:HA	5:E:141:VAL:CG2	2.45	0.43
7:G:1:MET:SD	7:G:79:PHE:CD1	3.12	0.43
11:K:43:GLY:O	11:K:47:ARG:HB2	2.17	0.43
11:K:84:LYS:O	11:K:87:LEU:HB3	2.17	0.43
1:A:49:LYS:NZ	1:A:61:ILE:HG13	2.28	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:671:ALA:N	1:A:676:MET:HE3	2.18	0.43
1:A:735:VAL:O	1:A:735:VAL:HG12	2.19	0.43
1:A:1037:LEU:HD13	1:A:1041:ALA:CB	2.49	0.43
2:B:641:GLU:HB2	2:B:643:ASP:OD1	2.19	0.43
2:B:693:ILE:HD11	2:B:740:HIS:CD2	2.52	0.43
2:B:745:PRO:HB2	2:B:1047:PHE:CD1	2.53	0.43
2:B:798:TYR:HE1	10:J:4:PRO:HA	1.84	0.43
2:B:866:TYR:HB2	2:B:870:ILE:HB	2.01	0.43
2:B:1023:VAL:HG12	2:B:1027:ILE:CD1	2.46	0.43
3:C:201:TRP:HA	3:C:201:TRP:CE3	2.53	0.43
6:F:83:PRO:HG2	6:F:84:TYR:CD1	2.54	0.43
7:G:106:MET:HG2	7:G:107:LYS:H	1.83	0.43
7:G:143:ILE:HG23	7:G:169:GLY:C	2.43	0.43
11:K:45:LEU:HG	11:K:94:ILE:HD13	2.00	0.43
1:A:36:ARG:C	1:A:270:LEU:HD21	2.44	0.43
1:A:98:LYS:HE2	1:A:224:PHE:HE1	1.83	0.43
1:A:548:ASN:HD21	11:K:47:ARG:NH2	2.15	0.43
1:A:679:ILE:CG2	1:A:729:ALA:HB1	2.47	0.43
1:A:692:ASP:C	1:A:694:THR:N	2.77	0.43
1:A:852:TYR:CD1	6:F:136:ARG:HB3	2.53	0.43
1:A:1152:ILE:CG2	1:A:1260:LEU:HD23	2.45	0.43
2:B:100:PRO:HB2	2:B:180:TYR:HE1	1.84	0.43
2:B:102:VAL:CG2	2:B:112:LEU:HD22	2.49	0.43
2:B:235:SER:C	2:B:236:HIS:CD2	2.97	0.43
2:B:240:ILE:HD12	2:B:241:ARG:H	1.83	0.43
2:B:583:ASN:HD21	2:B:628:THR:HG22	1.84	0.43
2:B:660:LYS:O	2:B:663:ALA:HB3	2.18	0.43
2:B:796:LEU:HD23	2:B:799:PRO:HA	1.99	0.43
2:B:900:ALA:O	2:B:903:VAL:HG23	2.18	0.43
2:B:1026:LEU:H	2:B:1026:LEU:HG	1.56	0.43
3:C:100:THR:HB	3:C:119:VAL:HG23	2.01	0.43
4:D:119:ARG:HB2	4:D:221:TYR:CE1	2.53	0.43
4:D:192:LYS:HD2	4:D:199:ASN:N	2.33	0.43
5:E:63:ASN:CB	5:E:64:PRO:CD	2.96	0.43
5:E:202:SER:C	5:E:204:THR:H	2.27	0.43
6:F:85:MET:O	6:F:155:LEU:HD21	2.19	0.43
6:F:125:LEU:C	6:F:127:GLU:N	2.76	0.43
7:G:84:GLY:HA2	7:G:146:LYS:HE2	2.00	0.43
7:G:121:PHE:HB2	7:G:130:TYR:CE2	2.54	0.43
1:A:37:PHE:HB2	1:A:52:GLY:HA3	2.00	0.43
1:A:59:GLY:HA3	1:A:244:PRO:HG2	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:185:TRP:HZ3	1:A:200:ARG:HB3	1.84	0.43
1:A:417:TYR:O	1:A:418:SER:C	2.62	0.43
1:A:500:GLU:CD	1:A:1438:THR:HG21	2.44	0.43
1:A:535:THR:HG21	1:A:617:VAL:HG23	2.00	0.43
1:A:578:LEU:O	1:A:580:VAL:N	2.52	0.43
1:A:876:ALA:C	1:A:878:ILE:H	2.27	0.43
1:A:1038:THR:O	1:A:1039:LYS:C	2.59	0.43
2:B:459:TYR:C	2:B:459:TYR:HD2	2.25	0.43
2:B:651:LEU:O	2:B:654:ARG:NE	2.50	0.43
2:B:1212:ILE:O	2:B:1214:PRO:HD3	2.17	0.43
3:C:49:VAL:HG22	3:C:157:CYS:HG	1.84	0.43
5:E:11:ARG:C	5:E:13:TRP:N	2.77	0.43
5:E:29:PHE:C	5:E:29:PHE:CD2	2.97	0.43
5:E:74:ASP:N	5:E:74:ASP:OD1	2.51	0.43
8:H:6:PHE:CD2	8:H:7:ASP:N	2.87	0.43
11:K:31:VAL:CG1	11:K:32:VAL:N	2.82	0.43
1:A:135:PHE:CD1	1:A:222:LEU:CD2	2.98	0.43
1:A:338:GLY:HA2	2:B:1129:ARG:NH2	2.34	0.43
1:A:794:PRO:C	1:A:796:SER:H	2.27	0.43
1:A:851:HIS:CB	6:F:139:PRO:HG3	2.48	0.43
1:A:1219:THR:HG23	1:A:1220:PHE:N	2.33	0.43
1:A:1433:MET:HE3	1:A:1433:MET:HB2	1.93	0.43
1:A:1437:GLY:C	1:A:1439:GLY:N	2.76	0.43
2:B:115:GLN:O	2:B:118:ARG:HB2	2.19	0.43
2:B:333:PHE:CD2	2:B:333:PHE:O	2.72	0.43
2:B:498:THR:O	2:B:536:VAL:HA	2.18	0.43
2:B:557:PHE:O	2:B:561:TRP:HD1	2.02	0.43
2:B:764:SER:N	2:B:765:PRO:CD	2.81	0.43
2:B:1033:LYS:O	2:B:1036:ALA:HB3	2.19	0.43
7:G:138:THR:O	7:G:140:LYS:N	2.51	0.43
1:A:22:PHE:HB3	2:B:1211:ASN:ND2	2.34	0.43
1:A:744:LYS:O	1:A:748:MET:CB	2.65	0.43
1:A:1067:LEU:O	1:A:1068:ALA:C	2.62	0.43
2:B:281:PRO:HG2	2:B:284:ILE:HG13	2.00	0.43
2:B:998:ASP:OD1	2:B:998:ASP:N	2.51	0.43
2:B:1167:GLY:HA3	2:B:1215:ARG:HA	2.00	0.43
3:C:46:ILE:HG13	3:C:72:LEU:HD21	2.01	0.43
5:E:190:LEU:N	5:E:190:LEU:HD23	2.33	0.43
7:G:91:VAL:HG12	7:G:92:VAL:N	2.34	0.43
12:L:55:ILE:HG12	12:L:56:LEU:N	2.34	0.43
1:A:15:LYS:HG3	2:B:1219:ASP:HA	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:34:LYS:NZ	1:A:57:ARG:NH1	2.67	0.42
1:A:42:ASP:HB3	1:A:45:GLN:CA	2.49	0.42
1:A:53:LEU:CD2	1:A:54:ASN:H	2.19	0.42
1:A:407:ARG:C	1:A:409:SER:H	2.27	0.42
1:A:445:ASN:C	1:A:487:MET:HE3	2.44	0.42
1:A:853:ASP:OD1	1:A:855:THR:CB	2.67	0.42
1:A:1015:VAL:O	1:A:1016:THR:C	2.61	0.42
2:B:59:LEU:HD21	2:B:128:LEU:HD11	2.01	0.42
2:B:166:PHE:CD2	2:B:166:PHE:C	2.96	0.42
2:B:215:GLN:HE21	2:B:476:ARG:HH11	1.67	0.42
2:B:294:ASP:OD1	9:I:12:ASN:HA	2.19	0.42
2:B:601:ARG:C	2:B:603:LEU:H	2.26	0.42
2:B:638:PHE:HB2	2:B:741:CYS:HB3	2.00	0.42
2:B:724:ASP:HB3	2:B:727:LYS:CG	2.48	0.42
2:B:857:ARG:HG2	2:B:858:SER:H	1.83	0.42
2:B:1106:ARG:CZ	2:B:1110:PRO:HD2	2.48	0.42
4:D:145:MET:O	4:D:149:THR:HG22	2.18	0.42
4:D:188:ALA:CA	4:D:191:ALA:HB3	2.33	0.42
7:G:51:TYR:CD2	7:G:51:TYR:C	2.97	0.42
11:K:94:ILE:O	11:K:98:LEU:HG	2.18	0.42
11:K:107:THR:O	11:K:111:LEU:HG	2.19	0.42
1:A:62:ASP:OD1	1:A:62:ASP:C	2.61	0.42
1:A:224:PHE:HE2	1:A:231:PRO:HA	1.83	0.42
1:A:311:GLN:O	1:A:312:PRO:C	2.61	0.42
1:A:343:LYS:HG2	2:B:1151:LEU:HD12	2.02	0.42
1:A:526:ASP:HB2	2:B:835:GLN:OE1	2.19	0.42
1:A:661:GLY:CA	2:B:1081:LEU:HD22	2.49	0.42
1:A:662:PHE:O	2:B:828:ALA:HA	2.18	0.42
1:A:821:ARG:O	1:A:825:ILE:HG13	2.19	0.42
1:A:837:ILE:O	1:A:838:GLN:C	2.59	0.42
1:A:845:LEU:O	1:A:1065:GLY:HA3	2.19	0.42
2:B:736:THR:HA	2:B:738:PHE:CE1	2.54	0.42
2:B:807:ARG:HB3	2:B:807:ARG:HH11	1.84	0.42
2:B:1043:ASP:O	2:B:1050:ILE:HD11	2.19	0.42
2:B:1173:ALA:HB1	2:B:1175:LEU:HD21	2.01	0.42
3:C:45:ALA:HA	3:C:72:LEU:HD13	2.00	0.42
3:C:70:ILE:HG22	3:C:71:PRO:CD	2.48	0.42
4:D:56:ARG:CD	4:D:149:THR:HA	2.43	0.42
4:D:198:LEU:N	4:D:198:LEU:HD23	2.34	0.42
5:E:78:LEU:HD23	5:E:78:LEU:C	2.44	0.42
5:E:92:THR:O	5:E:95:THR:HB	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:E:185:ALA:CA	5:E:190:LEU:HG	2.49	0.42
6:F:87:LYS:HG3	6:F:88:TYR:CE1	2.52	0.42
8:H:84:ALA:O	8:H:85:GLY:C	2.61	0.42
8:H:102:TYR:CE2	8:H:115:TYR:C	2.97	0.42
12:L:31:CYS:O	12:L:35:SER:HA	2.19	0.42
1:A:619:LYS:O	1:A:623:GLY:HA3	2.19	0.42
1:A:783:THR:HG21	1:A:796:SER:O	2.19	0.42
1:A:1025:ARG:O	1:A:1035:TYR:HE1	2.02	0.42
1:A:1100:ARG:HA	1:A:1103:GLU:OE1	2.20	0.42
1:A:1109:LYS:HA	1:A:1109:LYS:HD2	1.71	0.42
1:A:1161:THR:H	1:A:1170:ILE:HD13	1.84	0.42
1:A:1325:THR:O	5:E:148:GLU:HB3	2.20	0.42
1:A:1362:TYR:CE1	1:A:1363:VAL:O	2.72	0.42
1:A:1402:PHE:CD1	1:A:1403:GLU:HG2	2.54	0.42
2:B:70:ILE:HG13	2:B:429:PHE:HZ	1.84	0.42
2:B:281:PRO:O	2:B:283:VAL:N	2.53	0.42
2:B:459:TYR:O	2:B:463:THR:OG1	2.38	0.42
2:B:502:ILE:C	2:B:502:ILE:HD12	2.44	0.42
2:B:522:VAL:HG12	2:B:523:CYS:H	1.78	0.42
2:B:610:ASN:HA	2:B:611:PRO:HD3	1.92	0.42
2:B:610:ASN:C	2:B:612:GLU:H	2.26	0.42
2:B:972:LYS:CD	2:B:1098:MET:HE1	2.49	0.42
3:C:167:HIS:HA	11:K:6:ARG:HH12	1.84	0.42
5:E:167:ARG:O	5:E:168:TYR:CG	2.72	0.42
11:K:30:ALA:HB2	11:K:76:GLN:CG	2.49	0.42
11:K:53:ASP:C	11:K:55:LYS:N	2.77	0.42
1:A:41:MET:O	1:A:42:ASP:O	2.37	0.42
1:A:109:HIS:NE2	1:A:169:ASN:ND2	2.67	0.42
1:A:385:ILE:O	1:A:386:ASP:C	2.62	0.42
1:A:437:MET:C	1:A:438:ASP:O	2.63	0.42
1:A:809:THR:HG23	1:A:812:GLU:OE1	2.19	0.42
1:A:1365:TYR:O	1:A:1366:ARG:C	2.62	0.42
1:A:1397:LEU:N	1:A:1397:LEU:HD23	2.35	0.42
2:B:566:LEU:HD22	2:B:586:TRP:O	2.20	0.42
2:B:641:GLU:HB2	2:B:643:ASP:OD2	2.19	0.42
2:B:707:PRO:HB3	2:B:741:CYS:SG	2.58	0.42
3:C:69:LEU:O	10:J:6:ARG:NH1	2.52	0.42
4:D:173:HIS:O	4:D:175:PHE:N	2.53	0.42
1:A:92:HIS:CE1	1:A:1410:PHE:CE2	3.06	0.42
1:A:399:HIS:C	1:A:399:HIS:HD1	2.26	0.42
1:A:568:PRO:HB2	3:C:221:TYR:CE1	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:571:LEU:HD22	8:H:46:LEU:HD11	2.01	0.42
1:A:848:ILE:HA	1:A:857:ARG:O	2.20	0.42
1:A:904:THR:O	1:A:904:THR:HG22	2.19	0.42
1:A:1095:THR:HB	1:A:1100:ARG:HB2	2.01	0.42
1:A:1152:ILE:HG13	9:I:44:TYR:CD2	2.54	0.42
1:A:1450:LEU:HD21	7:G:19:GLY:O	2.19	0.42
2:B:378:LEU:O	2:B:378:LEU:HD12	2.19	0.42
2:B:498:THR:O	2:B:536:VAL:HG13	2.19	0.42
2:B:627:PHE:CE1	2:B:696:GLU:HG2	2.55	0.42
2:B:737:THR:O	2:B:738:PHE:C	2.62	0.42
2:B:797:TYR:C	2:B:799:PRO:HD3	2.44	0.42
2:B:809:MET:O	2:B:810:GLU:C	2.62	0.42
2:B:830:TYR:HB3	2:B:831:SER:H	1.44	0.42
2:B:1043:ASP:C	2:B:1045:SER:H	2.26	0.42
5:E:135:PHE:HD2	5:E:140:LEU:HD21	1.83	0.42
6:F:88:TYR:CD1	6:F:88:TYR:N	2.86	0.42
7:G:34:VAL:O	7:G:35:GLU:C	2.63	0.42
10:J:9:SER:HB2	10:J:45:CYS:CB	2.50	0.42
12:L:55:ILE:HG12	12:L:56:LEU:H	1.85	0.42
1:A:41:MET:HA	1:A:49:LYS:HA	2.01	0.42
1:A:59:GLY:O	1:A:244:PRO:HG2	2.19	0.42
1:A:548:ASN:ND2	11:K:47:ARG:HH21	2.15	0.42
1:A:661:GLY:O	1:A:662:PHE:HB2	2.19	0.42
1:A:784:LEU:C	1:A:786:HIS:H	2.27	0.42
1:A:907:THR:HG21	1:A:920:LEU:HD21	2.01	0.42
1:A:919:ILE:HD11	1:A:925:LEU:CD1	2.49	0.42
1:A:1211:GLN:O	1:A:1214:GLU:HB2	2.19	0.42
2:B:448:ILE:O	2:B:449:ASN:C	2.62	0.42
2:B:817:LEU:N	2:B:818:PRO:CD	2.81	0.42
2:B:916:THR:HA	2:B:917:PRO:HD3	1.82	0.42
3:C:242:GLN:C	3:C:244:VAL:N	2.76	0.42
5:E:116:ILE:HG22	5:E:120:ALA:HB3	2.01	0.42
5:E:162:ARG:NH1	5:E:162:ARG:HB3	2.34	0.42
8:H:89:LEU:C	8:H:91:ASP:N	2.76	0.42
8:H:130:ARG:N	8:H:130:ARG:HD3	2.35	0.42
1:A:151:ASP:OD1	1:A:163:SER:HB3	2.19	0.42
1:A:344:ARG:HA	2:B:1129:ARG:HA	2.02	0.42
1:A:424:ILE:HG22	1:A:425:GLN:O	2.19	0.42
1:A:881:GLN:CD	1:A:959:ASN:HA	2.45	0.42
1:A:914:GLU:HB2	1:A:979:SER:O	2.19	0.42
1:A:946:VAL:HG22	5:E:201:LYS:CD	2.44	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:999:VAL:C	1:A:1000:LEU:HG	2.45	0.42
1:A:1291:VAL:CG2	1:A:1292:PRO:HD2	2.49	0.42
2:B:129:PHE:CD2	2:B:166:PHE:HB2	2.54	0.42
2:B:286:PHE:CE1	2:B:378:LEU:HG	2.54	0.42
2:B:745:PRO:O	2:B:747:MET:N	2.53	0.42
2:B:1155:SER:O	2:B:1156:ASP:C	2.62	0.42
3:C:47:ASP:OD2	3:C:170:TRP:HH2	2.02	0.42
3:C:167:HIS:CD2	12:L:70:ARG:CG	3.03	0.42
3:C:187:LYS:HA	3:C:187:LYS:HD3	1.77	0.42
3:C:265:MET:HE3	3:C:265:MET:HB2	1.89	0.42
4:D:146:GLN:HA	4:D:149:THR:HG22	2.02	0.42
5:E:17:ARG:O	5:E:21:GLU:HG3	2.19	0.42
6:F:82:THR:CG2	6:F:83:PRO:HD2	2.50	0.42
7:G:79:PHE:CE2	7:G:105:PRO:HD2	2.55	0.42
1:A:89:PRO:HD2	1:A:205:GLU:HG3	2.02	0.42
1:A:524:VAL:O	1:A:525:GLN:C	2.62	0.42
1:A:555:ASP:OD2	1:A:644:LYS:HE2	2.20	0.42
1:A:779:PHE:CZ	2:B:517:THR:HA	2.55	0.42
1:A:851:HIS:HB3	6:F:139:PRO:HG3	2.02	0.42
1:A:1163:ILE:CG2	1:A:1164:PRO:HD2	2.50	0.42
1:A:1284:MET:CA	1:A:1306:LEU:HD23	2.50	0.42
1:A:1362:TYR:HD1	1:A:1362:TYR:C	2.28	0.42
1:A:1407:GLU:CD	1:A:1407:GLU:N	2.78	0.42
2:B:277:LYS:HD3	2:B:277:LYS:N	2.35	0.42
2:B:514:LEU:HD12	2:B:514:LEU:HA	1.91	0.42
2:B:792:MET:O	2:B:793:ALA:HB2	2.20	0.42
2:B:1060:ARG:C	2:B:1062:HIS:H	2.27	0.42
2:B:1130:PHE:HD2	2:B:1151:LEU:HD11	1.85	0.42
3:C:77:ILE:HD12	3:C:77:ILE:O	2.19	0.42
1:A:522:GLY:HA2	1:A:630:ILE:CD1	2.50	0.42
1:A:607:ILE:HG12	1:A:612:ILE:HA	2.02	0.42
1:A:1368:MET:HB2	1:A:1368:MET:HE3	1.52	0.42
2:B:173:MET:HE2	2:B:203:PHE:HE1	1.85	0.42
2:B:335:GLY:HA2	2:B:348:ARG:HB2	2.01	0.42
2:B:405:ARG:HA	2:B:631:GLY:O	2.19	0.42
2:B:461:LEU:HD12	2:B:461:LEU:HA	1.85	0.42
2:B:650:GLU:HG3	2:B:651:LEU:H	1.84	0.42
2:B:1174:LYS:O	2:B:1176:ASN:N	2.44	0.42
3:C:136:ASP:OD2	3:C:137:LYS:N	2.53	0.42
4:D:216:ASN:C	4:D:218:GLU:N	2.73	0.42
6:F:146:TRP:HB3	6:F:151:LEU:CD1	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:J:31:ASP:OD1	10:J:34:THR:HB	2.20	0.42
11:K:47:ARG:NH1	11:K:47:ARG:HB3	2.35	0.42
1:A:61:ILE:HG22	1:A:62:ASP:H	1.84	0.42
1:A:101:LYS:HG2	1:A:139:TRP:CE2	2.55	0.42
1:A:225:ASN:HD22	1:A:227:VAL:H	1.68	0.42
1:A:442:VAL:HG12	1:A:490:HIS:O	2.20	0.42
1:A:565:ILE:O	1:A:570:PRO:HA	2.20	0.42
1:A:670:ILE:HG22	1:A:805:LEU:HD11	2.01	0.42
1:A:802:ASN:ND2	2:B:728:ARG:HB2	2.35	0.42
1:A:963:ILE:HD11	1:A:1048:ASN:CB	2.48	0.42
1:A:1035:TYR:O	1:A:1036:ARG:CB	2.58	0.42
1:A:1156:PRO:HA	1:A:1190:PRO:CB	2.35	0.42
1:A:1426:GLU:H	1:A:1426:GLU:HG2	1.51	0.42
2:B:467:GLY:HA3	2:B:475:SER:HB3	2.01	0.42
2:B:515:HIS:CG	2:B:516:ASN:N	2.88	0.42
2:B:515:HIS:O	2:B:518:HIS:HB2	2.20	0.42
2:B:619:ILE:C	2:B:621:GLU:H	2.28	0.42
2:B:785:TYR:CE2	10:J:60:PHE:HE1	2.37	0.42
2:B:840:ILE:HG21	2:B:1011:ILE:HD12	2.02	0.42
2:B:899:ILE:CG2	2:B:949:VAL:HG21	2.42	0.42
2:B:1074:ASN:OD1	2:B:1075:GLY:N	2.53	0.42
2:B:1098:MET:HE3	2:B:1098:MET:HB2	1.90	0.42
5:E:54:GLN:O	5:E:57:MET:HB3	2.20	0.42
5:E:100:ILE:HG23	5:E:105:PHE:HB2	2.01	0.42
6:F:125:LEU:HA	6:F:130:ILE:HD11	2.01	0.42
7:G:79:PHE:HD2	7:G:105:PRO:HD2	1.85	0.42
1:A:61:ILE:HG22	1:A:62:ASP:N	2.35	0.41
1:A:335:ARG:HA	1:A:339:ASN:ND2	2.27	0.41
1:A:403:LYS:HB3	1:A:404:TYR:CD1	2.54	0.41
1:A:425:GLN:NE2	1:A:425:GLN:N	2.54	0.41
1:A:503:GLN:OE1	1:A:503:GLN:HA	2.20	0.41
1:A:899:VAL:CB	1:A:929:LEU:HD12	2.36	0.41
1:A:942:PHE:CD2	1:A:942:PHE:C	2.98	0.41
1:A:948:VAL:O	1:A:949:ASP:C	2.63	0.41
1:A:1099:PRO:HG2	1:A:1100:ARG:H	1.85	0.41
1:A:1242:VAL:O	1:A:1243:VAL:CG1	2.67	0.41
1:A:1264:GLU:HG3	1:A:1265:ASN:OD1	2.20	0.41
1:A:1445:ILE:HD12	1:A:1445:ILE:C	2.45	0.41
2:B:263:GLY:O	2:B:264:SER:C	2.62	0.41
2:B:278:GLN:HB3	2:B:279:ASP:H	1.69	0.41
2:B:860:MET:HG3	2:B:965:LYS:HG2	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:900:ALA:HA	2:B:901:PRO:HD3	1.86	0.41
2:B:1133:MET:HB3	2:B:1133:MET:HE2	1.86	0.41
5:E:152:LYS:HD2	5:E:152:LYS:HA	1.86	0.41
7:G:146:LYS:HB2	7:G:168:LEU:HD11	2.02	0.41
7:G:160:ILE:CG2	7:G:161:GLY:N	2.81	0.41
7:G:163:ILE:HA	7:G:168:LEU:HD13	2.02	0.41
11:K:55:LYS:HB2	11:K:81:TYR:CE1	2.55	0.41
1:A:162:VAL:HG12	1:A:163:SER:H	1.85	0.41
1:A:284:ALA:HB1	1:A:289:ILE:HD11	2.03	0.41
1:A:399:HIS:O	1:A:435:HIS:HD2	2.04	0.41
1:A:474:VAL:HG23	1:A:477:PRO:HG2	2.02	0.41
1:A:475:THR:HG23	1:A:476:SER:N	2.35	0.41
1:A:541:ILE:HD12	1:A:577:ILE:CD1	2.49	0.41
1:A:580:VAL:O	1:A:580:VAL:HG12	2.19	0.41
1:A:654:ASN:O	1:A:657:LEU:HB3	2.20	0.41
1:A:767:GLN:HE21	1:A:774:ARG:HB2	1.81	0.41
1:A:1277:GLU:O	1:A:1278:ASN:HB2	2.19	0.41
1:A:1444:MET:HE2	1:A:1444:MET:HB2	1.80	0.41
2:B:68:THR:HG22	2:B:91:SER:CA	2.50	0.41
2:B:778:MET:SD	2:B:794:ASN:HB3	2.59	0.41
2:B:781:PHE:CD1	2:B:782:LEU:HG	2.54	0.41
2:B:1051:THR:HB	2:B:1054:GLY:H	1.84	0.41
3:C:16:ASP:HA	3:C:240:VAL:HG13	2.02	0.41
3:C:77:ILE:CA	3:C:129:ILE:HD11	2.49	0.41
3:C:93:ASP:O	3:C:127:ARG:NH2	2.52	0.41
3:C:124:LEU:HD23	3:C:126:GLY:HA2	2.02	0.41
4:D:146:GLN:O	4:D:147:TYR:C	2.63	0.41
4:D:213:GLU:O	4:D:216:ASN:HB2	2.20	0.41
8:H:43:ASN:OD1	8:H:43:ASN:C	2.63	0.41
9:I:15:TYR:N	9:I:15:TYR:HD1	2.16	0.41
11:K:99:GLY:O	11:K:102:LYS:HB3	2.20	0.41
1:A:389:THR:N	1:A:426:LEU:HD11	2.35	0.41
1:A:501:LEU:HA	1:A:505:CYS:SG	2.60	0.41
1:A:928:LEU:O	1:A:929:LEU:C	2.62	0.41
1:A:1011:GLN:O	1:A:1012:ARG:C	2.63	0.41
1:A:1127:ASP:CG	1:A:1130:GLN:HB2	2.46	0.41
1:A:1139:GLU:O	1:A:1140:HIS:C	2.63	0.41
1:A:1297:GLU:CD	1:A:1297:GLU:H	2.28	0.41
1:A:1396:ALA:HA	1:A:1399:ARG:CZ	2.49	0.41
2:B:487:THR:O	2:B:490:SER:N	2.53	0.41
2:B:549:THR:CG2	2:B:550:ASP:H	2.32	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:846:ILE:HG23	2:B:974:PRO:HG2	2.02	0.41
2:B:1038:SER:O	10:J:33:GLY:HA3	2.20	0.41
4:D:123:LEU:HD22	4:D:149:THR:HG21	2.02	0.41
4:D:160:VAL:O	4:D:161:GLY:C	2.62	0.41
1:A:258:GLY:O	1:A:259:GLU:O	2.39	0.41
1:A:528:LEU:CG	1:A:529:CYS:N	2.81	0.41
1:A:562:THR:CG2	1:A:563:PRO:HD2	2.50	0.41
1:A:690:VAL:HG23	1:A:721:PHE:HD1	1.85	0.41
1:A:829:VAL:HG11	2:B:508:LEU:HD13	2.03	0.41
1:A:1342:GLU:C	1:A:1344:GLY:N	2.76	0.41
2:B:21:GLU:O	2:B:21:GLU:HG3	2.21	0.41
2:B:519:TRP:HE1	2:B:635:ARG:NH2	2.18	0.41
2:B:754:SER:HB2	2:B:812:LEU:CD1	2.49	0.41
2:B:1107:ALA:O	2:B:1108:ARG:HB3	2.18	0.41
2:B:1159:ARG:HD2	2:B:1159:ARG:C	2.46	0.41
4:D:48:ILE:HG21	7:G:4:ILE:HB	2.01	0.41
4:D:198:LEU:O	4:D:199:ASN:HB2	2.20	0.41
5:E:45:LYS:HB3	5:E:46:TYR:CE1	2.54	0.41
6:F:152:ILE:HG22	6:F:153:VAL:H	1.85	0.41
9:I:17:ARG:CG	9:I:18:GLU:H	2.31	0.41
11:K:11:LEU:HA	11:K:11:LEU:HD22	1.65	0.41
1:A:35:ILE:HG13	1:A:52:GLY:O	2.21	0.41
1:A:69:THR:O	1:A:70:CYS:C	2.64	0.41
1:A:325:ILE:C	1:A:327:ALA:N	2.75	0.41
1:A:767:GLN:HE22	1:A:797:LYS:CB	2.34	0.41
1:A:784:LEU:O	1:A:786:HIS:N	2.53	0.41
1:A:1006:ILE:HD12	5:E:167:ARG:HB2	2.01	0.41
1:A:1011:GLN:HE22	1:A:1015:VAL:CG2	2.32	0.41
1:A:1343:ALA:O	1:A:1346:ALA:HB3	2.21	0.41
1:A:1362:TYR:CD1	1:A:1363:VAL:N	2.87	0.41
2:B:596:LEU:HD21	2:B:624:LEU:HD22	2.03	0.41
2:B:1122:ARG:C	2:B:1124:ARG:N	2.77	0.41
4:D:194:LEU:N	4:D:194:LEU:CD2	2.82	0.41
5:E:213:ILE:CG1	5:E:214:CYS:N	2.81	0.41
6:F:96:THR:HG22	6:F:100:GLN:OE1	2.20	0.41
9:I:33:SER:O	9:I:34:TYR:C	2.62	0.41
1:A:512:VAL:O	1:A:512:VAL:HG12	2.20	0.41
1:A:531:ILE:CG1	1:A:578:LEU:HD21	2.51	0.41
1:A:1008:GLN:O	1:A:1009:ASN:C	2.64	0.41
1:A:1120:LEU:HD12	1:A:1120:LEU:N	2.35	0.41
1:A:1166:ASP:O	1:A:1170:ILE:HG13	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1259:MET:HG3	1:A:1263:ILE:HD11	2.03	0.41
1:A:1263:ILE:HG13	1:A:1263:ILE:H	1.43	0.41
2:B:119:LEU:HD23	2:B:953:LEU:CD1	2.49	0.41
2:B:120:ARG:HG2	2:B:955:THR:HG21	2.01	0.41
2:B:121:ASN:N	2:B:121:ASN:HD22	2.19	0.41
2:B:547:VAL:HG12	2:B:612:GLU:OE2	2.20	0.41
2:B:637:LEU:HD13	2:B:740:HIS:HB2	2.01	0.41
2:B:650:GLU:CG	2:B:651:LEU:N	2.82	0.41
2:B:850:LEU:HA	10:J:8:PHE:HE1	1.84	0.41
3:C:53:THR:CG2	3:C:152:GLU:HB3	2.51	0.41
3:C:123:ASN:ND2	3:C:125:MET:HG2	2.36	0.41
4:D:180:LEU:HD11	4:D:198:LEU:HD11	2.01	0.41
5:E:30:ILE:HG23	5:E:34:GLU:HB3	2.01	0.41
5:E:119:SER:C	5:E:122:LYS:HB2	2.45	0.41
10:J:18:TRP:CH2	10:J:22:LEU:HD11	2.55	0.41
12:L:67:PHE:CD1	12:L:67:PHE:N	2.89	0.41
1:A:75:ASN:O	1:A:76:GLU:HB2	2.20	0.41
1:A:125:ALA:HA	1:A:128:ILE:CD1	2.51	0.41
1:A:284:ALA:CB	1:A:289:ILE:HD11	2.50	0.41
1:A:285:PRO:O	1:A:288:ALA:N	2.53	0.41
1:A:919:ILE:HD11	1:A:925:LEU:HD11	2.02	0.41
1:A:1068:ALA:HB1	1:A:1367:HIS:HA	2.03	0.41
1:A:1230:GLU:C	1:A:1232:ASN:N	2.79	0.41
2:B:243:ALA:HA	2:B:250:PHE:O	2.19	0.41
2:B:976:ILE:HD13	2:B:976:ILE:HA	1.85	0.41
3:C:73:GLN:O	3:C:129:ILE:HA	2.20	0.41
3:C:136:ASP:O	3:C:137:LYS:C	2.64	0.41
3:C:148:ARG:H	3:C:151:GLN:CG	2.26	0.41
4:D:202:ILE:HG12	4:D:207:LEU:HB2	2.02	0.41
6:F:85:MET:HA	6:F:151:LEU:HD23	2.02	0.41
1:A:243:PRO:O	1:A:246:VAL:HG23	2.21	0.41
1:A:335:ARG:HD3	1:A:335:ARG:HA	1.82	0.41
1:A:339:ASN:HB3	2:B:1117:GLN:HE22	1.86	0.41
1:A:445:ASN:HB2	1:A:455:MET:HA	2.03	0.41
1:A:954:TRP:HA	1:A:955:PRO:HD3	1.97	0.41
1:A:1025:ARG:HA	1:A:1030:ARG:CD	2.51	0.41
1:A:1101:LEU:HB2	1:A:1355:VAL:HG11	2.02	0.41
1:A:1293:SER:OG	1:A:1295:THR:HG22	2.21	0.41
2:B:51:PHE:O	2:B:52:ASN:C	2.64	0.41
2:B:235:SER:CB	2:B:236:HIS:HD2	2.33	0.41
2:B:430:ARG:CG	2:B:430:ARG:NH1	2.63	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:1007:VAL:CG2	2:B:1008:PRO:N	2.83	0.41
2:B:1034:VAL:HA	2:B:1037:LEU:HD11	2.03	0.41
2:B:1132:GLU:O	2:B:1135:ARG:HB3	2.21	0.41
3:C:58:LEU:HD13	3:C:62:PHE:CE2	2.56	0.41
3:C:166:GLU:C	11:K:6:ARG:HH11	2.29	0.41
5:E:12:LEU:HD13	5:E:12:LEU:HA	1.82	0.41
5:E:13:TRP:O	5:E:16:PHE:HB3	2.21	0.41
5:E:136:ASN:OD1	5:E:138:ALA:N	2.54	0.41
6:F:86:THR:O	6:F:87:LYS:C	2.63	0.41
1:A:120:GLU:HG3	1:A:123:ARG:NH1	2.36	0.41
1:A:208:LEU:O	1:A:209:ASN:C	2.64	0.41
1:A:385:ILE:CG1	1:A:428:TYR:HE2	2.33	0.41
1:A:446:ARG:HB3	1:A:478:TYR:HB3	2.01	0.41
1:A:677:ARG:HD2	1:A:677:ARG:C	2.46	0.41
1:A:710:LEU:H	1:A:710:LEU:HD12	1.85	0.41
1:A:730:GLY:O	1:A:731:ARG:C	2.62	0.41
1:A:745:GLN:HA	1:A:748:MET:HE3	2.02	0.41
1:A:809:THR:CG2	2:B:730:ARG:HG3	2.51	0.41
1:A:830:LYS:HE2	1:A:1081:LEU:HB2	2.03	0.41
1:A:1153:TYR:CB	1:A:1192:LEU:HD23	2.47	0.41
1:A:1213:GLY:O	1:A:1217:LYS:HB2	2.21	0.41
1:A:1216:ILE:H	1:A:1216:ILE:HG13	1.71	0.41
1:A:1236:LEU:C	1:A:1237:ILE:HG13	2.45	0.41
1:A:1351:GLU:O	1:A:1352:VAL:C	2.64	0.41
2:B:56:ASP:HB3	2:B:57:TYR:CE1	2.55	0.41
2:B:167:ILE:HD12	2:B:167:ILE:N	2.36	0.41
2:B:202:TYR:C	2:B:203:PHE:HD1	2.29	0.41
2:B:205:ILE:HG22	2:B:206:ASN:N	2.35	0.41
2:B:512:ARG:O	2:B:513:GLN:O	2.39	0.41
2:B:582:VAL:C	2:B:584:GLY:H	2.29	0.41
2:B:822:ASN:N	2:B:822:ASN:ND2	2.69	0.41
2:B:841:MET:O	2:B:993:THR:HA	2.21	0.41
2:B:874:PHE:HD1	2:B:962:LYS:HD3	1.85	0.41
2:B:979:LYS:C	2:B:980:PHE:CG	2.99	0.41
2:B:983:ARG:NH1	2:B:1091:TYR:HB3	2.35	0.41
2:B:1166:CYS:O	2:B:1166:CYS:SG	2.78	0.41
3:C:70:ILE:HA	3:C:71:PRO:HD3	1.86	0.41
3:C:258:ILE:HA	3:C:258:ILE:HD12	1.85	0.41
5:E:20:LYS:O	5:E:21:GLU:C	2.64	0.41
5:E:147:HIS:O	5:E:150:VAL:HG23	2.20	0.41
6:F:76:LYS:HA	6:F:79:ARG:CD	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:F:103:MET:CE	7:G:66:GLY:H	2.31	0.41
7:G:94:CYS:SG	7:G:130:TYR:HE1	2.43	0.41
8:H:102:TYR:HE2	8:H:116:TYR:N	2.19	0.41
8:H:106:GLU:HG2	8:H:112:ILE:CD1	2.50	0.41
9:I:7:CYS:O	9:I:11:ASN:HA	2.21	0.41
10:J:54:VAL:HG12	10:J:56:LEU:HD23	2.03	0.41
11:K:44:ASN:O	11:K:48:ALA:HB2	2.20	0.41
11:K:102:LYS:O	11:K:103:THR:C	2.63	0.41
1:A:33:ALA:HA	1:A:57:ARG:NH1	2.36	0.41
1:A:54:ASN:HD22	1:A:54:ASN:N	2.17	0.41
1:A:204:THR:O	1:A:206:GLU:N	2.54	0.41
1:A:224:PHE:HE2	1:A:234:MET:HE2	1.85	0.41
1:A:242:PRO:HA	1:A:243:PRO:HD3	1.85	0.41
1:A:269:ILE:O	1:A:270:LEU:C	2.62	0.41
1:A:277:GLU:C	1:A:279:LEU:H	2.29	0.41
1:A:311:GLN:O	1:A:313:GLN:N	2.53	0.41
1:A:403:LYS:C	1:A:404:TYR:CD1	2.99	0.41
1:A:766:GLY:HA2	1:A:799:PHE:CE1	2.55	0.41
1:A:836:TYR:CE2	1:A:840:ARG:CD	3.04	0.41
1:A:857:ARG:HA	1:A:864:ILE:CD1	2.51	0.41
2:B:39:ARG:NH2	2:B:665:GLU:HG2	2.35	0.41
2:B:92:PHE:CE1	2:B:421:PHE:HZ	2.38	0.41
2:B:313:MET:HE3	2:B:386:LEU:HD13	2.02	0.41
2:B:407:ASP:HB3	2:B:412:LEU:HD21	2.03	0.41
2:B:420:LEU:HD23	2:B:420:LEU:HA	1.94	0.41
2:B:839:MET:HE3	2:B:1010:LEU:HD21	2.03	0.41
2:B:941:LEU:HG	2:B:942:ARG:H	1.86	0.41
2:B:1224:PHE:CZ	5:E:171:LYS:HE3	2.56	0.41
3:C:168:ALA:O	3:C:169:LYS:C	2.64	0.41
3:C:221:TYR:CZ	3:C:222:LYS:HE3	2.56	0.41
3:C:259:LEU:HD12	3:C:259:LEU:HA	1.87	0.41
4:D:8:PHE:CE1	4:D:37:GLN:HB2	2.55	0.41
4:D:13:ARG:O	4:D:15:LEU:N	2.54	0.41
5:E:178:ILE:O	5:E:179:GLN:O	2.39	0.41
5:E:188:LEU:HB2	5:E:190:LEU:HD21	2.03	0.41
6:F:74:ILE:HD12	6:F:144:GLU:CG	2.50	0.41
7:G:12:THR:HG22	7:G:13:LEU:N	2.36	0.41
10:J:41:LEU:HD23	10:J:41:LEU:H	1.86	0.41
1:A:25:GLU:OE2	1:A:25:GLU:N	2.54	0.40
1:A:35:ILE:HD12	1:A:35:ILE:N	2.22	0.40
1:A:49:LYS:CE	1:A:61:ILE:HG13	2.51	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:265:LYS:CD	1:A:322:VAL:HG21	2.37	0.40
1:A:779:PHE:CE1	1:A:785:PRO:HD3	2.55	0.40
1:A:900:ASP:HB3	1:A:906:HIS:HB3	2.03	0.40
1:A:1445:ILE:HD11	7:G:59:GLY:N	2.37	0.40
2:B:254:LEU:HD12	2:B:272:THR:O	2.21	0.40
2:B:418:LYS:C	2:B:420:LEU:N	2.79	0.40
2:B:639:ILE:HD11	2:B:691:GLU:HB2	2.03	0.40
2:B:640:VAL:O	2:B:640:VAL:CG1	2.69	0.40
2:B:820:GLY:H	2:B:1091:TYR:HH	1.68	0.40
2:B:821:GLN:NE2	2:B:851:PHE:H	2.18	0.40
2:B:854:LEU:HD23	2:B:854:LEU:HA	1.84	0.40
2:B:996:ARG:H	2:B:996:ARG:HG3	1.39	0.40
2:B:997:GLU:C	2:B:999:MET:H	2.28	0.40
2:B:1074:ASN:OD1	2:B:1074:ASN:C	2.64	0.40
2:B:1096:ARG:O	2:B:1097:HIS:C	2.63	0.40
3:C:115:SER:HB3	3:C:142:VAL:HB	2.03	0.40
3:C:206:ASN:O	3:C:207:CYS:C	2.63	0.40
4:D:13:ARG:C	4:D:15:LEU:N	2.79	0.40
5:E:13:TRP:HA	5:E:42:PHE:CE2	2.56	0.40
6:F:69:LEU:HB3	6:F:71:GLU:CD	2.46	0.40
10:J:48:ARG:NH1	10:J:48:ARG:CG	2.78	0.40
1:A:195:ASP:OD1	1:A:195:ASP:N	2.54	0.40
1:A:323:LYS:N	1:A:323:LYS:HD2	2.36	0.40
1:A:391:LEU:O	1:A:394:ASN:N	2.55	0.40
1:A:441:PRO:O	1:A:492:PRO:CG	2.70	0.40
1:A:505:CYS:O	1:A:506:ALA:C	2.64	0.40
1:A:741:ASN:HD21	1:A:743:VAL:H	1.70	0.40
1:A:909:ASP:C	1:A:911:SER:N	2.79	0.40
1:A:1004:ASN:HD21	1:A:1006:ILE:HB	1.86	0.40
2:B:189:LEU:HD13	2:B:196:PRO:HA	2.03	0.40
2:B:522:VAL:CG1	2:B:523:CYS:H	2.31	0.40
2:B:526:GLU:HG2	2:B:538:ASN:HB2	2.03	0.40
2:B:563:MET:O	2:B:565:PRO:HD3	2.21	0.40
2:B:1108:ARG:HG2	2:B:1109:GLY:H	1.85	0.40
2:B:1183:LYS:HD2	2:B:1183:LYS:C	2.47	0.40
3:C:8:VAL:CG1	3:C:9:LYS:N	2.85	0.40
3:C:263:THR:O	3:C:266:ASP:HB2	2.21	0.40
5:E:127:ILE:N	5:E:128:PRO:HD3	2.36	0.40
5:E:171:LYS:H	5:E:171:LYS:HG2	1.51	0.40
6:F:131:PRO:O	6:F:132:LEU:HD23	2.21	0.40
7:G:21:ARG:HD3	7:G:21:ARG:HA	1.89	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:G:80:LYS:HG2	7:G:82:PHE:CE2	2.56	0.40
10:J:9:SER:HB2	10:J:45:CYS:SG	2.61	0.40
10:J:24:LEU:HD23	10:J:24:LEU:HA	1.87	0.40
1:A:305:ASP:OD1	1:A:306:ASN:N	2.55	0.40
1:A:379:VAL:HG12	1:A:380:VAL:N	2.36	0.40
1:A:389:THR:O	1:A:390:GLN:C	2.63	0.40
1:A:490:HIS:HB3	2:B:1150:ARG:CZ	2.50	0.40
1:A:531:ILE:HG12	1:A:578:LEU:HD21	2.03	0.40
1:A:698:GLN:C	1:A:700:ASN:H	2.29	0.40
1:A:1266:THR:O	1:A:1266:THR:HG22	2.21	0.40
1:A:1279:ILE:HG22	1:A:1282:VAL:HG21	2.03	0.40
2:B:38:PHE:CD2	2:B:43:LEU:CD2	3.05	0.40
2:B:216:GLU:CA	2:B:406:LEU:HD23	2.45	0.40
2:B:221:ASN:HA	2:B:241:ARG:O	2.21	0.40
2:B:508:LEU:HA	2:B:512:ARG:HE	1.86	0.40
2:B:529:GLU:OE1	2:B:769:TYR:HE2	2.04	0.40
2:B:589:VAL:HG12	2:B:590:HIS:N	2.35	0.40
2:B:773:MET:O	2:B:776:GLN:HB2	2.22	0.40
2:B:833:TYR:C	2:B:835:GLN:N	2.73	0.40
2:B:1076:HIS:HD2	11:K:40:HIS:CD2	2.39	0.40
2:B:1114:LEU:HD12	2:B:1114:LEU:HA	1.78	0.40
3:C:13:ALA:HA	3:C:18:VAL:HA	2.03	0.40
3:C:136:ASP:HB2	3:C:141:GLY:CA	2.51	0.40
4:D:205:ASP:C	4:D:208:GLU:HB3	2.46	0.40
5:E:52:ARG:HA	5:E:53:PRO:HD3	1.79	0.40
5:E:124:VAL:O	5:E:132:ILE:CD1	2.69	0.40
7:G:91:VAL:CG1	7:G:92:VAL:N	2.85	0.40
10:J:5:VAL:HA	10:J:15:GLY:N	2.34	0.40
10:J:8:PHE:CD2	10:J:49:MET:SD	3.15	0.40
11:K:38:GLU:HB3	11:K:71:PHE:HE2	1.86	0.40
1:A:98:LYS:HE2	1:A:224:PHE:CE1	2.56	0.40
1:A:154:SER:C	1:A:156:ASP:H	2.29	0.40
1:A:640:GLN:O	1:A:641:VAL:C	2.64	0.40
1:A:670:ILE:O	1:A:670:ILE:HG12	2.21	0.40
1:A:1069:ALA:O	1:A:1070:GLN:C	2.65	0.40
1:A:1398:MET:CE	1:A:1424:VAL:HG23	2.51	0.40
2:B:42:GLY:C	2:B:43:LEU:HD23	2.46	0.40
2:B:259:TYR:HB2	2:B:268:THR:HG23	2.03	0.40
2:B:276:ILE:O	2:B:276:ILE:CG2	2.68	0.40
2:B:347:LYS:HG2	2:B:347:LYS:H	1.71	0.40
2:B:780:VAL:HG11	10:J:56:LEU:HD12	2.02	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:796:LEU:HD11	2:B:821:GLN:HE21	1.86	0.40
2:B:1030:LEU:HD12	2:B:1030:LEU:HA	1.89	0.40
2:B:1109:GLY:H	2:B:1111:MET:CE	2.35	0.40
3:C:260:LEU:HG	3:C:264:GLN:NE2	2.37	0.40
5:E:136:ASN:OD1	5:E:136:ASN:C	2.65	0.40
5:E:171:LYS:O	5:E:172:GLU:C	2.64	0.40
6:F:82:THR:HA	6:F:83:PRO:HD3	1.99	0.40
1:A:127:ALA:O	1:A:129:LYS:N	2.55	0.40
1:A:497:THR:HG23	2:B:1146:PHE:HD1	1.86	0.40
1:A:1205:LYS:O	1:A:1206:ASP:C	2.65	0.40
2:B:213:ILE:H	2:B:479:VAL:HG12	1.86	0.40
2:B:421:PHE:O	2:B:424:LEU:HB3	2.22	0.40
2:B:636:PRO:HG2	2:B:743:ILE:HD11	2.03	0.40
2:B:990:ILE:HG22	2:B:992:ILE:N	2.34	0.40
2:B:1138:MET:HB3	2:B:1147:LEU:HD21	2.00	0.40
4:D:124:GLU:O	4:D:128:VAL:HG23	2.21	0.40
5:E:33:GLU:C	5:E:35:VAL:N	2.79	0.40
8:H:38:LEU:HD13	8:H:125:LEU:CD1	2.52	0.40
11:K:5:ASP:HB2	11:K:7:PHE:CE2	2.54	0.40
12:L:53:HIS:HB3	12:L:55:ILE:HD11	2.00	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	1406/1733 (81%)	974 (69%)	284 (20%)	148 (10%)	0	6
2	B	1081/1224 (88%)	754 (70%)	232 (22%)	95 (9%)	0	8
3	C	264/318 (83%)	187 (71%)	48 (18%)	29 (11%)	0	6
4	D	174/219 (80%)	116 (67%)	43 (25%)	15 (9%)	0	9

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
5	E	212/215 (99%)	156 (74%)	41 (19%)	15 (7%)	1	11
6	F	85/155 (55%)	61 (72%)	19 (22%)	5 (6%)	1	14
7	G	169/171 (99%)	120 (71%)	43 (25%)	6 (4%)	2	20
8	H	130/146 (89%)	90 (69%)	29 (22%)	11 (8%)	0	9
9	I	45/133 (34%)	31 (69%)	10 (22%)	4 (9%)	0	8
10	J	63/70 (90%)	43 (68%)	14 (22%)	6 (10%)	0	8
11	K	112/120 (93%)	79 (70%)	24 (21%)	9 (8%)	1	9
12	L	44/70 (63%)	21 (48%)	8 (18%)	15 (34%)	0	0
All	All	3785/4574 (83%)	2632 (70%)	795 (21%)	358 (10%)	0	8

All (358) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	35	ILE
1	A	42	ASP
1	A	48	ALA
1	A	54	ASN
1	A	57	ARG
1	A	67	CYS
1	A	68	GLN
1	A	74	MET
1	A	119	ASN
1	A	128	ILE
1	A	130	ASP
1	A	167	CYS
1	A	257	ARG
1	A	259	GLU
1	A	286	HIS
1	A	318	SER
1	A	330	LYS
1	A	332	LYS
1	A	361	LEU
1	A	424	ILE
1	A	439	ASN
1	A	465	TYR
1	A	525	GLN
1	A	527	THR
1	A	536	LEU
1	A	557	ASP

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Mol	Chain	Res	Type
1	A	567	LYS
1	A	591	PHE
1	A	597	LEU
1	A	598	LEU
1	A	666	ILE
1	A	790	ASP
1	A	868	TYR
1	A	986	ILE
1	A	1002	GLY
1	A	1036	ARG
1	A	1098	VAL
1	A	1139	GLU
1	A	1223	ASP
1	A	1229	SER
1	A	1232	ASN
1	A	1255	GLU
1	A	1270	ASN
1	A	1339	LEU
1	A	1365	TYR
1	A	1392	SER
1	A	1394	THR
1	A	1405	THR
1	A	1424	VAL
2	B	108	VAL
2	B	184	ALA
2	B	206	ASN
2	B	258	LEU
2	B	364	ILE
2	B	367	LEU
2	B	509	ALA
2	B	732	SER
2	B	735	ALA
2	B	738	PHE
2	B	1036	ALA
2	B	1046	PRO
2	B	1096	ARG
2	B	1103	ILE
2	B	1132	GLU
2	B	1155	SER
2	B	1156	ASP
2	B	1183	LYS
3	C	110	THR

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Mol	Chain	Res	Type
3	C	121	VAL
3	C	128	ASN
3	C	137	LYS
3	C	184	ASN
3	C	227	THR
4	D	19	GLU
4	D	119	ARG
5	E	45	LYS
5	E	105	PHE
5	E	168	TYR
5	E	179	GLN
6	F	128	LYS
6	F	139	PRO
6	F	140	ASP
7	G	41	LYS
7	G	139	ILE
8	H	82	PRO
8	H	128	ASN
8	H	140	ALA
10	J	2	ILE
10	J	42	LYS
10	J	64	ASN
11	K	28	PRO
11	K	64	GLU
12	L	35	SER
12	L	39	SER
12	L	46	VAL
12	L	50	ASP
12	L	59	ALA
12	L	60	ARG
12	L	63	ARG
1	A	52	GLY
1	A	66	LYS
1	A	76	GLU
1	A	166	GLY
1	A	204	THR
1	A	253	ASN
1	A	308	ILE
1	A	310	GLY
1	A	314	ALA
1	A	322	VAL
1	A	418	SER

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Mol	Chain	Res	Type
1	A	421	ALA
1	A	438	ASP
1	A	592	ASP
1	A	605	MET
1	A	628	GLY
1	A	636	GLU
1	A	653	VAL
1	A	693	VAL
1	A	780	VAL
1	A	875	ALA
1	A	912	LEU
1	A	985	ASP
1	A	999	VAL
1	A	1067	LEU
1	A	1120	LEU
1	A	1140	HIS
1	A	1188	GLN
1	A	1206	ASP
1	A	1221	LYS
1	A	1233	ASP
1	A	1242	VAL
1	A	1314	SER
1	A	1316	VAL
1	A	1367	HIS
1	A	1437	GLY
2	B	28	GLU
2	B	65	GLU
2	B	68	THR
2	B	369	GLY
2	B	435	THR
2	B	447	ALA
2	B	449	ASN
2	B	483	LEU
2	B	513	GLN
2	B	591	ARG
2	B	731	VAL
2	B	751	VAL
2	B	777	ALA
2	B	786	ASN
2	B	838	SER
2	B	887	HIS
2	B	937	ALA

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Mol	Chain	Res	Type
2	B	938	SER
2	B	1041	GLU
2	B	1079	LYS
2	B	1108	ARG
2	B	1112	GLN
2	B	1157	ALA
2	B	1158	PHE
2	B	1167	GLY
3	C	5	GLY
3	C	11	ARG
3	C	126	GLY
3	C	141	GLY
3	C	149	LYS
3	C	168	ALA
3	C	213	PRO
3	C	240	VAL
4	D	5	THR
4	D	14	ARG
4	D	16	LYS
4	D	168	LYS
4	D	171	GLY
5	E	3	GLN
5	E	75	MET
5	E	106	GLN
6	F	112	GLU
7	G	154	VAL
8	H	95	TYR
8	H	109	LYS
9	I	3	THR
9	I	8	ARG
10	J	6	ARG
11	K	68	PHE
12	L	26	THR
12	L	47	ARG
12	L	48	CYS
12	L	55	ILE
12	L	56	LEU
1	A	28	ARG
1	A	59	GLY
1	A	169	ASN
1	A	219	PHE
1	A	255	SER

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Mol	Chain	Res	Type
1	A	312	PRO
1	A	317	LYS
1	A	382	PRO
1	A	556	TRP
1	A	596	THR
1	A	785	PRO
1	A	869	GLY
1	A	910	PRO
1	A	1062	GLU
1	A	1114	PRO
1	A	1388	GLY
1	A	1403	GLU
1	A	1422	ARG
1	A	1448	GLU
2	B	27	ALA
2	B	45	SER
2	B	171	PRO
2	B	180	TYR
2	B	249	ARG
2	B	250	PHE
2	B	291	ILE
2	B	409	ALA
2	B	468	GLU
2	B	562	GLY
2	B	575	PRO
2	B	746	SER
2	B	760	ASP
2	B	775	LYS
2	B	802	PRO
2	B	834	ASN
2	B	907	GLY
2	B	959	ASP
2	B	1175	LEU
2	B	1176	ASN
2	B	1223	ASP
3	C	90	ASP
3	C	122	SER
3	C	142	VAL
3	C	167	HIS
3	C	214	ASN
3	C	241	ASP
4	D	31	GLN

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Mol	Chain	Res	Type
4	D	134	THR
4	D	191	ALA
4	D	198	LEU
5	E	73	PRO
5	E	76	GLY
5	E	130	ALA
5	E	151	PRO
7	G	28	THR
7	G	67	SER
8	H	59	ILE
8	H	60	ALA
8	H	81	PRO
9	I	47	GLU
11	K	8	GLU
12	L	45	ALA
12	L	64	LEU
1	A	170	THR
1	A	196	GLU
1	A	420	ARG
1	A	649	ILE
1	A	654	ASN
1	A	958	VAL
1	A	1016	THR
1	A	1124	HIS
1	A	1266	THR
1	A	1377	THR
1	A	1435	PRO
2	B	106	ASP
2	B	186	GLU
2	B	459	TYR
2	B	510	LYS
2	B	524	PRO
2	B	566	LEU
2	B	711	GLU
2	B	906	SER
2	B	1017	ILE
2	B	1169	MET
3	C	13	ALA
3	C	25	VAL
3	C	174	ALA
3	C	224	GLN
5	E	97	VAL

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Mol	Chain	Res	Type
6	F	141	GLY
8	H	83	GLN
10	J	13	VAL
11	K	54	ARG
12	L	40	LEU
1	A	45	GLN
1	A	62	ASP
1	A	75	ASN
1	A	155	GLU
1	A	205	GLU
1	A	396	PRO
1	A	579	SER
1	A	583	PRO
1	A	602	ASP
1	A	647	GLY
1	A	920	LEU
1	A	1080	THR
1	A	1345	ARG
2	B	46	GLN
2	B	124	TYR
2	B	304	ASP
2	B	636	PRO
2	B	707	PRO
2	B	758	PHE
2	B	889	THR
2	B	974	PRO
2	B	987	LYS
2	B	996	ARG
3	C	127	ARG
3	C	132	PRO
4	D	218	GLU
8	H	107	VAL
9	I	9	ASP
10	J	14	VAL
11	K	7	PHE
11	K	10	PHE
11	K	83	PRO
1	A	4	GLN
1	A	492	PRO
1	A	533	LYS
1	A	662	PHE
1	A	976	THR

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Mol	Chain	Res	Type
1	A	1302	PRO
1	A	1428	VAL
2	B	218	SER
2	B	219	ALA
2	B	463	THR
2	B	902	GLY
3	C	175	ALA
4	D	9	GLN
4	D	52	LEU
5	E	118	PRO
5	E	124	VAL
7	G	84	GLY
1	A	311	GLN
1	A	410	GLY
1	A	730	GLY
2	B	323	VAL
2	B	478	GLY
3	C	216	GLY
4	D	18	VAL
8	H	17	PRO
1	A	258	GLY
1	A	385	ILE
1	A	1107	VAL
2	B	24	PRO
1	A	197	PRO
1	A	392	VAL
2	B	100	PRO
2	B	260	GLY
1	A	250	ILE
1	A	800	VAL
1	A	937	VAL
1	A	1122	PRO
3	C	172	PRO
1	A	1190	PRO
2	B	114	PRO
2	B	1014	PRO
11	K	66	PRO
5	E	129	PRO

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	1239/1520 (82%)	1008 (81%)	231 (19%)	1	10
2	B	957/1061 (90%)	791 (83%)	166 (17%)	2	11
3	C	234/274 (85%)	196 (84%)	38 (16%)	2	12
4	D	160/198 (81%)	127 (79%)	33 (21%)	1	7
5	E	196/197 (100%)	169 (86%)	27 (14%)	3	15
6	F	77/137 (56%)	64 (83%)	13 (17%)	2	12
7	G	152/152 (100%)	122 (80%)	30 (20%)	1	8
8	H	118/128 (92%)	95 (80%)	23 (20%)	1	8
9	I	45/122 (37%)	38 (84%)	7 (16%)	2	13
10	J	60/65 (92%)	52 (87%)	8 (13%)	4	16
11	K	99/102 (97%)	80 (81%)	19 (19%)	1	9
12	L	40/57 (70%)	29 (72%)	11 (28%)	0	3
All	All	3377/4013 (84%)	2771 (82%)	606 (18%)	2	11

All (606) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	VAL
1	A	8	SER
1	A	11	LEU
1	A	17	VAL
1	A	22	PHE
1	A	32	VAL
1	A	34	LYS
1	A	40	THR
1	A	54	ASN
1	A	58	LEU
1	A	60	SER
1	A	62	ASP
1	A	64	ASN
1	A	68	GLN

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Mol	Chain	Res	Type
1	A	84	ILE
1	A	96	ILE
1	A	107	CYS
1	A	109	HIS
1	A	121	LEU
1	A	126	LEU
1	A	145	LYS
1	A	151	ASP
1	A	160	GLN
1	A	162	VAL
1	A	169	ASN
1	A	179	LEU
1	A	182	VAL
1	A	186	LYS
1	A	215	SER
1	A	216	VAL
1	A	222	LEU
1	A	225	ASN
1	A	232	GLU
1	A	236	LEU
1	A	237	THR
1	A	249	SER
1	A	261	ASP
1	A	262	LEU
1	A	263	THR
1	A	264	PHE
1	A	265	LYS
1	A	278	THR
1	A	300	VAL
1	A	302	THR
1	A	304	MET
1	A	308	ILE
1	A	316	GLN
1	A	320	ARG
1	A	322	VAL
1	A	337	ARG
1	A	344	ARG
1	A	345	VAL
1	A	351	THR
1	A	352	VAL
1	A	353	ILE
1	A	359	LEU

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Mol	Chain	Res	Type
1	A	368	LYS
1	A	370	ILE
1	A	375	THR
1	A	379	VAL
1	A	383	TYR
1	A	386	ASP
1	A	399	HIS
1	A	404	TYR
1	A	406	ILE
1	A	415	LEU
1	A	424	ILE
1	A	425	GLN
1	A	434	ARG
1	A	438	ASP
1	A	440	ASP
1	A	442	VAL
1	A	443	LEU
1	A	445	ASN
1	A	451	HIS
1	A	461	LYS
1	A	468	PHE
1	A	485	ASP
1	A	491	VAL
1	A	493	GLN
1	A	494	SER
1	A	507	VAL
1	A	509	LEU
1	A	512	VAL
1	A	513	SER
1	A	516	SER
1	A	524	VAL
1	A	528	LEU
1	A	533	LYS
1	A	534	LEU
1	A	538	ASP
1	A	542	GLU
1	A	566	ILE
1	A	571	LEU
1	A	577	ILE
1	A	578	LEU
1	A	582	ILE
1	A	586	ILE

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Mol	Chain	Res	Type
1	A	595	THR
1	A	596	THR
1	A	602	ASP
1	A	613	ILE
1	A	617	VAL
1	A	618	GLU
1	A	622	VAL
1	A	629	LEU
1	A	630	ILE
1	A	635	ARG
1	A	640	GLN
1	A	652	VAL
1	A	657	LEU
1	A	666	ILE
1	A	670	ILE
1	A	680	THR
1	A	690	VAL
1	A	691	LEU
1	A	738	LYS
1	A	739	ASP
1	A	741	ASN
1	A	742	ASN
1	A	743	VAL
1	A	745	GLN
1	A	762	SER
1	A	768	GLN
1	A	774	ARG
1	A	805	LEU
1	A	827	THR
1	A	831	THR
1	A	833	GLU
1	A	838	GLN
1	A	841	LEU
1	A	845	LEU
1	A	852	TYR
1	A	858	ASN
1	A	860	LEU
1	A	864	ILE
1	A	886	ILE
1	A	899	VAL
1	A	902	LEU
1	A	903	ASN

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Mol	Chain	Res	Type
1	A	908	LEU
1	A	919	ILE
1	A	933	TYR
1	A	936	LEU
1	A	948	VAL
1	A	969	GLN
1	A	970	THR
1	A	976	THR
1	A	983	ILE
1	A	987	VAL
1	A	988	LEU
1	A	990	VAL
1	A	999	VAL
1	A	1000	LEU
1	A	1001	ARG
1	A	1011	GLN
1	A	1017	LEU
1	A	1019	CYS
1	A	1029	ARG
1	A	1033	GLN
1	A	1036	ARG
1	A	1037	LEU
1	A	1057	VAL
1	A	1058	VAL
1	A	1067	LEU
1	A	1072	ILE
1	A	1078	GLN
1	A	1095	THR
1	A	1103	GLU
1	A	1110	ASN
1	A	1111	MET
1	A	1116	LEU
1	A	1117	THR
1	A	1120	LEU
1	A	1121	GLU
1	A	1124	HIS
1	A	1134	ILE
1	A	1142	THR
1	A	1148	ILE
1	A	1152	ILE
1	A	1159	ARG
1	A	1167	GLU

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Mol	Chain	Res	Type
1	A	1171	GLN
1	A	1217	LYS
1	A	1219	THR
1	A	1228	TRP
1	A	1229	SER
1	A	1241	ARG
1	A	1255	GLU
1	A	1259	MET
1	A	1263	ILE
1	A	1264	GLU
1	A	1271	ILE
1	A	1272	THR
1	A	1274	ARG
1	A	1276	VAL
1	A	1280	GLU
1	A	1289	ARG
1	A	1291	VAL
1	A	1295	THR
1	A	1297	GLU
1	A	1299	VAL
1	A	1305	VAL
1	A	1308	THR
1	A	1309	ASP
1	A	1315	GLU
1	A	1325	THR
1	A	1329	THR
1	A	1333	ILE
1	A	1336	MET
1	A	1349	TYR
1	A	1368	MET
1	A	1370	LEU
1	A	1373	ASP
1	A	1377	THR
1	A	1378	GLN
1	A	1382	THR
1	A	1384	VAL
1	A	1385	THR
1	A	1397	LEU
1	A	1401	SER
1	A	1403	GLU
1	A	1405	THR
1	A	1411	GLU

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Mol	Chain	Res	Type
1	A	1424	VAL
1	A	1426	GLU
1	A	1428	VAL
1	A	1429	ILE
1	A	1442	ASP
1	A	1443	VAL
1	A	1451	VAL
2	B	43	LEU
2	B	44	VAL
2	B	46	GLN
2	B	47	GLN
2	B	57	TYR
2	B	61	ASP
2	B	63	ILE
2	B	90	ILE
2	B	95	ILE
2	B	102	VAL
2	B	115	GLN
2	B	165	VAL
2	B	167	ILE
2	B	181	LEU
2	B	185	THR
2	B	186	GLU
2	B	188	ASP
2	B	191	LYS
2	B	194	GLU
2	B	202	TYR
2	B	217	ARG
2	B	218	SER
2	B	222	ILE
2	B	225	VAL
2	B	237	VAL
2	B	249	ARG
2	B	259	TYR
2	B	261	ARG
2	B	272	THR
2	B	276	ILE
2	B	282	ILE
2	B	283	VAL
2	B	284	ILE
2	B	289	LEU
2	B	305	VAL

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Mol	Chain	Res	Type
2	B	324	ILE
2	B	328	GLU
2	B	329	THR
2	B	363	HIS
2	B	365	THR
2	B	368	GLU
2	B	378	LEU
2	B	387	LEU
2	B	401	PHE
2	B	408	LEU
2	B	413	LEU
2	B	416	LEU
2	B	427	ASP
2	B	429	PHE
2	B	430	ARG
2	B	463	THR
2	B	476	ARG
2	B	479	VAL
2	B	482	VAL
2	B	485	ARG
2	B	489	SER
2	B	496	ARG
2	B	500	THR
2	B	516	ASN
2	B	531	GLN
2	B	538	ASN
2	B	539	LEU
2	B	552	MET
2	B	558	LEU
2	B	559	SER
2	B	563	MET
2	B	568	ASP
2	B	570	VAL
2	B	576	ASP
2	B	580	VAL
2	B	597	MET
2	B	602	THR
2	B	615	MET
2	B	616	ILE
2	B	619	ILE
2	B	626	ILE
2	B	628	THR

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Mol	Chain	Res	Type
2	B	635	ARG
2	B	641	GLU
2	B	646	LEU
2	B	678	GLU
2	B	684	LEU
2	B	708	GLU
2	B	709	ASP
2	B	714	GLU
2	B	730	ARG
2	B	737	THR
2	B	739	THR
2	B	743	ILE
2	B	748	ILE
2	B	751	VAL
2	B	773	MET
2	B	780	VAL
2	B	786	ASN
2	B	787	VAL
2	B	791	THR
2	B	797	TYR
2	B	807	ARG
2	B	815	ARG
2	B	816	GLU
2	B	822	ASN
2	B	830	TYR
2	B	839	MET
2	B	841	MET
2	B	842	ASN
2	B	844	SER
2	B	856	PHE
2	B	859	TYR
2	B	862	GLN
2	B	869	SER
2	B	878	GLN
2	B	879	ARG
2	B	885	MET
2	B	890	TYR
2	B	895	ASP
2	B	904	ARG
2	B	914	LYS
2	B	942	ARG
2	B	944	THR

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Mol	Chain	Res	Type
2	B	946	ASN
2	B	952	VAL
2	B	953	LEU
2	B	957	ASN
2	B	959	ASP
2	B	969	ARG
2	B	976	ILE
2	B	987	LYS
2	B	989	THR
2	B	996	ARG
2	B	998	ASP
2	B	999	MET
2	B	1006	ILE
2	B	1007	VAL
2	B	1010	LEU
2	B	1012	ILE
2	B	1020	ARG
2	B	1026	LEU
2	B	1029	CYS
2	B	1034	VAL
2	B	1037	LEU
2	B	1051	THR
2	B	1071	VAL
2	B	1076	HIS
2	B	1084	GLN
2	B	1095	LEU
2	B	1113	VAL
2	B	1115	THR
2	B	1135	ARG
2	B	1145	SER
2	B	1147	LEU
2	B	1149	GLU
2	B	1150	ARG
2	B	1156	ASP
2	B	1159	ARG
2	B	1162	ILE
2	B	1170	THR
2	B	1176	ASN
2	B	1182	CYS
2	B	1183	LYS
2	B	1185	CYS
2	B	1194	ILE

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Mol	Chain	Res	Type
2	B	1196	ILE
2	B	1210	MET
2	B	1212	ILE
2	B	1216	LEU
2	B	1218	THR
3	C	7	GLN
3	C	22	LEU
3	C	25	VAL
3	C	43	THR
3	C	48	SER
3	C	51	VAL
3	C	53	THR
3	C	56	THR
3	C	63	ILE
3	C	70	ILE
3	C	72	LEU
3	C	77	ILE
3	C	83	SER
3	C	89	GLU
3	C	100	THR
3	C	106	GLU
3	C	112	ASN
3	C	118	LEU
3	C	119	VAL
3	C	123	ASN
3	C	124	LEU
3	C	134	ILE
3	C	143	LEU
3	C	147	LEU
3	C	155	LEU
3	C	166	GLU
3	C	167	HIS
3	C	178	PHE
3	C	190	ASP
3	C	193	TYR
3	C	231	ASN
3	C	232	VAL
3	C	235	VAL
3	C	240	VAL
3	C	245	VAL
3	C	254	LYS
3	C	258	ILE

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Mol	Chain	Res	Type
3	C	259	LEU
4	D	5	THR
4	D	11	ARG
4	D	13	ARG
4	D	17	LYS
4	D	18	VAL
4	D	20	GLU
4	D	22	GLU
4	D	23	ASN
4	D	26	THR
4	D	29	LEU
4	D	34	GLN
4	D	40	HIS
4	D	47	LEU
4	D	51	ASN
4	D	59	ILE
4	D	65	GLU
4	D	70	PHE
4	D	118	THR
4	D	120	GLU
4	D	123	LEU
4	D	132	GLN
4	D	151	PHE
4	D	156	ASP
4	D	180	LEU
4	D	183	LEU
4	D	194	LEU
4	D	195	ILE
4	D	198	LEU
4	D	202	ILE
4	D	206	GLU
4	D	214	LEU
4	D	220	LEU
4	D	221	TYR
5	E	6	GLU
5	E	7	ARG
5	E	12	LEU
5	E	37	LEU
5	E	72	PHE
5	E	85	GLU
5	E	90	VAL
5	E	100	ILE

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Mol	Chain	Res	Type
5	E	102	GLU
5	E	106	GLN
5	E	107	THR
5	E	113	GLN
5	E	114	ASN
5	E	115	ASN
5	E	132	ILE
5	E	144	ILE
5	E	150	VAL
5	E	153	HIS
5	E	165	LEU
5	E	168	TYR
5	E	171	LYS
5	E	173	SER
5	E	175	LEU
5	E	188	LEU
5	E	191	LYS
5	E	196	VAL
5	E	214	CYS
6	F	71	GLU
6	F	79	ARG
6	F	86	THR
6	F	93	ILE
6	F	94	LEU
6	F	101	ILE
6	F	107	VAL
6	F	112	GLU
6	F	120	ILE
6	F	138	LEU
6	F	140	ASP
6	F	147	SER
6	F	152	ILE
7	G	1	MET
7	G	7	LEU
7	G	9	LEU
7	G	13	LEU
7	G	21	ARG
7	G	42	PHE
7	G	45	ILE
7	G	48	VAL
7	G	53	ASN
7	G	56	ILE

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Mol	Chain	Res	Type
7	G	60	ARG
7	G	61	ILE
7	G	67	SER
7	G	74	TYR
7	G	75	ARG
7	G	87	VAL
7	G	88	ASP
7	G	96	GLN
7	G	97	HIS
7	G	110	VAL
7	G	111	THR
7	G	113	HIS
7	G	114	LEU
7	G	138	THR
7	G	139	ILE
7	G	150	CYS
7	G	154	VAL
7	G	162	SER
7	G	165	GLU
7	G	168	LEU
8	H	7	ASP
8	H	27	GLU
8	H	33	GLN
8	H	35	GLN
8	H	38	LEU
8	H	39	THR
8	H	44	VAL
8	H	49	VAL
8	H	53	ASP
8	H	57	VAL
8	H	59	ILE
8	H	61	SER
8	H	64	ASN
8	H	89	LEU
8	H	102	TYR
8	H	103	LYS
8	H	107	VAL
8	H	128	ASN
8	H	129	TYR
8	H	130	ARG
8	H	135	LEU
8	H	138	GLU

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Mol	Chain	Res	Type
8	H	146	ARG
9	I	2	THR
9	I	6	PHE
9	I	8	ARG
9	I	15	TYR
9	I	19	ASP
9	I	26	LEU
9	I	28	GLU
10	J	2	ILE
10	J	7	CYS
10	J	8	PHE
10	J	28	ASP
10	J	41	LEU
10	J	44	TYR
10	J	48	ARG
10	J	55	ASP
11	K	5	ASP
11	K	11	LEU
11	K	32	VAL
11	K	34	THR
11	K	41	THR
11	K	42	LEU
11	K	47	ARG
11	K	50	LEU
11	K	51	LEU
11	K	53	ASP
11	K	56	VAL
11	K	57	LEU
11	K	61	TYR
11	K	65	HIS
11	K	68	PHE
11	K	75	ILE
11	K	81	TYR
11	K	111	LEU
11	K	114	LEU
12	L	27	LEU
12	L	38	LEU
12	L	46	VAL
12	L	47	ARG
12	L	48	CYS
12	L	54	ARG
12	L	55	ILE

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Mol	Chain	Res	Type
12	L	61	THR
12	L	64	LEU
12	L	65	VAL
12	L	68	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (120) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	54	ASN
1	A	68	GLN
1	A	75	ASN
1	A	83	HIS
1	A	92	HIS
1	A	225	ASN
1	A	282	ASN
1	A	286	HIS
1	A	297	GLN
1	A	306	ASN
1	A	311	GLN
1	A	339	ASN
1	A	394	ASN
1	A	425	GLN
1	A	435	HIS
1	A	445	ASN
1	A	451	HIS
1	A	471	ASN
1	A	479	ASN
1	A	515	GLN
1	A	548	ASN
1	A	603	ASN
1	A	631	HIS
1	A	660	ASN
1	A	706	HIS
1	A	736	ASN
1	A	741	ASN
1	A	742	ASN
1	A	745	GLN
1	A	760	GLN
1	A	767	GLN
1	A	786	HIS
1	A	838	GLN
1	A	851	HIS

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Mol	Chain	Res	Type
1	A	858	ASN
1	A	877	HIS
1	A	903	ASN
1	A	926	GLN
1	A	965	GLN
1	A	966	ASN
1	A	972	HIS
1	A	1004	ASN
1	A	1011	GLN
1	A	1048	ASN
1	A	1106	ASN
1	A	1110	ASN
1	A	1124	HIS
1	A	1140	HIS
1	A	1173	HIS
1	A	1211	GLN
1	A	1354	ASN
1	A	1432	GLN
2	B	47	GLN
2	B	53	GLN
2	B	60	GLN
2	B	121	ASN
2	B	178	ASN
2	B	236	HIS
2	B	465	ASN
2	B	515	HIS
2	B	516	ASN
2	B	518	HIS
2	B	531	GLN
2	B	538	ASN
2	B	667	GLN
2	B	706	GLN
2	B	734	HIS
2	B	744	HIS
2	B	761	HIS
2	B	786	ASN
2	B	821	GLN
2	B	835	GLN
2	B	842	ASN
2	B	862	GLN
2	B	887	HIS
2	B	946	ASN

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Mol	Chain	Res	Type
2	B	951	GLN
2	B	957	ASN
2	B	975	GLN
2	B	1015	HIS
2	B	1065	GLN
2	B	1076	HIS
2	B	1084	GLN
2	B	1097	HIS
2	B	1161	HIS
2	B	1193	GLN
2	B	1211	ASN
3	C	7	GLN
3	C	54	ASN
3	C	65	HIS
3	C	112	ASN
3	C	131	HIS
3	C	203	GLN
3	C	231	ASN
3	C	242	GLN
3	C	252	GLN
4	D	31	GLN
4	D	39	ASN
4	D	51	ASN
4	D	132	GLN
4	D	179	GLN
4	D	199	ASN
5	E	3	GLN
5	E	5	ASN
5	E	54	GLN
5	E	101	GLN
5	E	113	GLN
5	E	143	ASN
5	E	153	HIS
7	G	14	HIS
7	G	53	ASN
7	G	97	HIS
7	G	113	HIS
7	G	126	ASN
7	G	158	HIS
8	H	128	ASN
8	H	131	ASN
9	I	22	ASN

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Mol	Chain	Res	Type
11	K	65	HIS
12	L	66	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 8 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	1416/1733 (81%)	0.07	37 (2%) 57 44	31, 133, 223, 314	0
2	B	1103/1224 (90%)	0.14	39 (3%) 47 37	23, 141, 230, 357	0
3	C	266/318 (83%)	0.04	5 (1%) 66 51	58, 139, 209, 281	0
4	D	178/219 (81%)	0.00	7 (3%) 43 36	64, 146, 212, 263	0
5	E	214/215 (99%)	0.10	4 (1%) 66 51	71, 165, 246, 367	0
6	F	87/155 (56%)	0.07	2 (2%) 61 48	48, 95, 156, 221	0
7	G	171/171 (100%)	0.05	2 (1%) 76 61	43, 121, 199, 230	0
8	H	134/146 (91%)	0.38	10 (7%) 20 23	115, 185, 250, 406	0
9	I	47/133 (35%)	0.33	3 (6%) 25 26	93, 163, 204, 221	0
10	J	65/70 (92%)	0.23	2 (3%) 51 40	80, 147, 210, 261	0
11	K	114/120 (95%)	0.05	2 (1%) 67 53	63, 128, 194, 254	0
12	L	46/70 (65%)	0.53	3 (6%) 25 25	93, 159, 240, 306	0
All	All	3841/4574 (83%)	0.11	116 (3%) 52 40	23, 140, 227, 406	0

All (116) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	2	VAL	6.8
1	A	1080	THR	5.8
2	B	245	GLU	4.9
2	B	246	LYS	4.5
1	A	1317	MET	4.4
8	H	2	SER	4.1
1	A	480	ALA	4.0
2	B	715	ALA	3.8
4	D	171	GLY	3.8
5	E	93	MET	3.8
1	A	481	ASP	3.8

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Mol	Chain	Res	Type	RSRZ
1	A	195	ASP	3.4
2	B	397	ASP	3.4
2	B	1109	GLY	3.4
8	H	139	ASN	3.3
1	A	1204	ASP	3.2
1	A	893	PHE	3.1
2	B	658	ILE	3.1
12	L	25	ALA	3.1
12	L	70	ARG	3.1
1	A	309	ALA	3.0
2	B	25	ILE	3.0
8	H	77	ARG	3.0
7	G	137	ILE	2.9
12	L	37	LYS	2.9
2	B	953	LEU	2.9
2	B	529	GLU	2.9
1	A	69	THR	2.9
1	A	251	SER	2.9
2	B	886	LYS	2.8
8	H	45	GLU	2.8
8	H	140	ALA	2.8
1	A	790	ASP	2.8
2	B	722	ASP	2.8
2	B	92	PHE	2.8
1	A	1455	PRO	2.7
3	C	204	SER	2.7
4	D	135	GLY	2.7
9	I	31	THR	2.7
2	B	726	ALA	2.7
11	K	105	PHE	2.7
3	C	178	PHE	2.6
2	B	813	LYS	2.6
2	B	597	MET	2.6
1	A	585	GLY	2.6
1	A	303	TYR	2.6
1	A	627	GLY	2.6
1	A	250	ILE	2.6
6	F	104	ASN	2.6
2	B	477	ALA	2.6
2	B	954	VAL	2.6
9	I	2	THR	2.6
2	B	1097	HIS	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	237	THR	2.6
1	A	1081	LEU	2.5
1	A	521	MET	2.5
2	B	737	THR	2.5
8	H	64	ASN	2.5
10	J	36	LEU	2.4
2	B	1153	GLU	2.4
2	B	731	VAL	2.4
8	H	76	THR	2.4
2	B	1175	LEU	2.4
2	B	1041	GLU	2.4
2	B	837	ASP	2.4
4	D	207	LEU	2.4
2	B	70	ILE	2.4
5	E	35	VAL	2.4
1	A	626	ASN	2.4
4	D	200	ASN	2.4
3	C	152	GLU	2.4
10	J	41	LEU	2.4
2	B	1184	GLY	2.3
1	A	1175	SER	2.3
4	D	189	ASP	2.3
1	A	179	LEU	2.3
1	A	659	HIS	2.3
6	F	116	ASP	2.3
1	A	542	GLU	2.3
1	A	114	LEU	2.3
11	K	87	LEU	2.3
1	A	1254	ALA	2.3
1	A	997	LEU	2.3
2	B	786	ASN	2.3
5	E	144	ILE	2.3
3	C	224	GLN	2.3
1	A	483	ASP	2.2
1	A	545	GLN	2.2
2	B	963	PHE	2.2
1	A	877	HIS	2.2
2	B	838	SER	2.2
1	A	710	LEU	2.2
8	H	134	ASN	2.2
4	D	38	ILE	2.2
5	E	58	MET	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	704	ALA	2.2
2	B	641	GLU	2.2
1	A	374	LEU	2.1
2	B	1071	VAL	2.1
3	C	115	SER	2.1
4	D	3	VAL	2.1
1	A	1123	GLY	2.1
2	B	991	GLY	2.1
2	B	964	VAL	2.1
1	A	666	ILE	2.1
8	H	57	VAL	2.1
7	G	22	MET	2.1
2	B	1223	ASP	2.1
8	H	59	ILE	2.1
2	B	130	VAL	2.1
1	A	447	GLN	2.0
9	I	3	THR	2.0
2	B	1067	ARG	2.0
2	B	918	ILE	2.0
2	B	533	CYS	2.0
2	B	1004	GLU	2.0

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

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

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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($\text{\AA}^2$)	Q<0.9
14	MG	A	3009	1/1	0.57	0.23	89,89,89,89	0
13	ZN	A	3006	1/1	0.97	0.04	274,274,274,274	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
13	ZN	L	3005	1/1	0.98	0.04	229,229,229,229	0
13	ZN	J	3001	1/1	0.99	0.06	153,153,153,153	0
13	ZN	A	3008	1/1	0.99	0.06	97,97,97,97	0
13	ZN	C	3002	1/1	0.99	0.08	129,129,129,129	0
13	ZN	I	3003	1/1	1.00	0.05	133,133,133,133	0
13	ZN	B	3007	1/1	1.00	0.05	116,116,116,116	0

6.5 Other polymers [i](#)

There are no such residues in this entry.