



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 02:09 PM UTC

PDB ID : 1M1C / pdb_00001m1c
Title : Structure of the L-A virus
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Deposited on : 2002-06-18
Resolution : 3.50 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

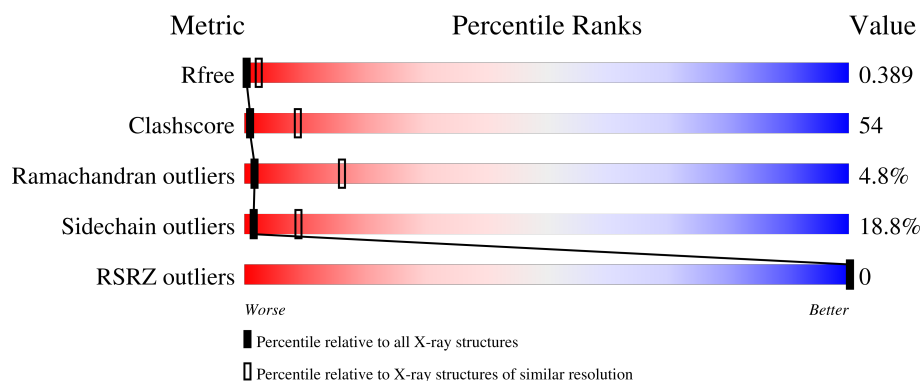
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

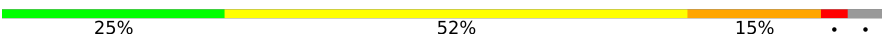
The reported resolution of this entry is 3.50 Å.

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	1085 (3.54-3.46)
Clashscore	190562	1140 (3.54-3.46)
Ramachandran outliers	187476	1113 (3.54-3.46)
Sidechain outliers	187428	1114 (3.54-3.46)
RSRZ outliers	180081	1084 (3.54-3.46)

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	680	
1	B	680	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 10302 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.

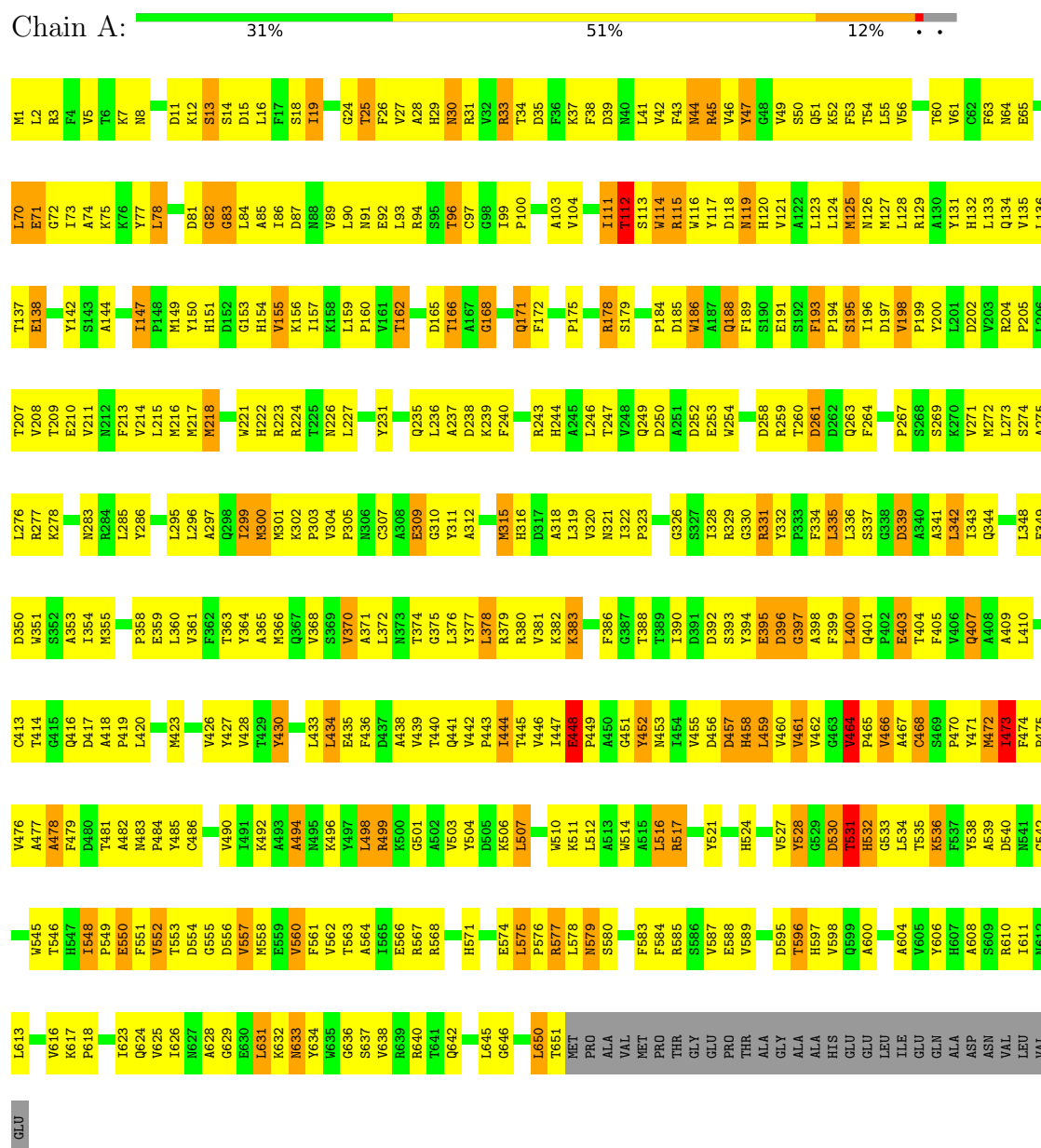
- Molecule 1 is a protein called Major coat protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	651	Total	C	N	O	S	0	0	0
			5151	3302	871	955	23			
1	B	651	Total	C	N	O	S	0	0	0
			5151	3302	871	955	23			

3 Residue-property plots

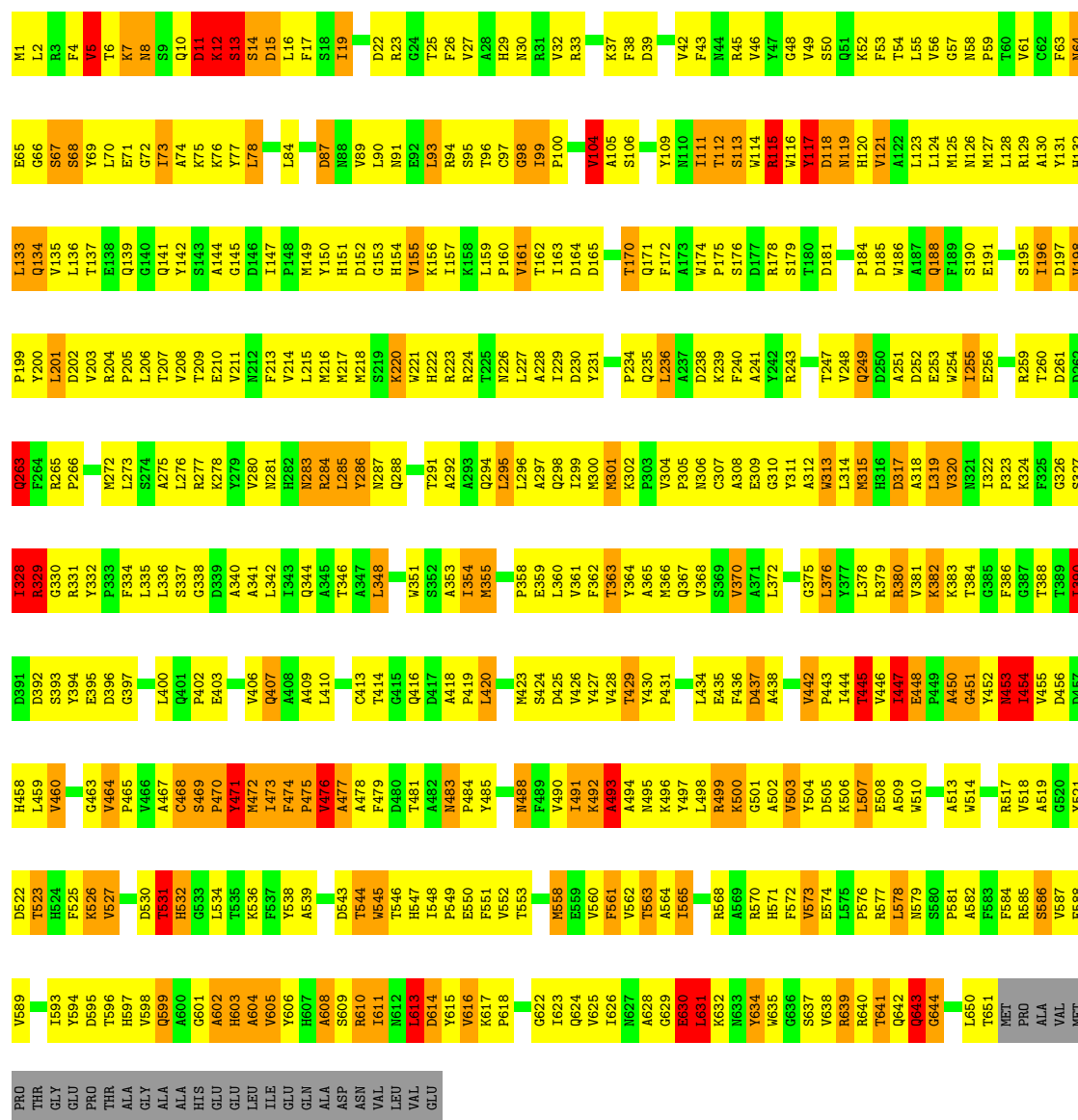
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Major coat protein



- Molecule 1: Major coat protein

Chain B: 25% 52% 15% . .



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	407.00Å 403.20Å 572.00Å 90.00° 90.46° 90.00°	Depositor
Resolution (Å)	30.00 – 3.50 30.00 – 3.50	Depositor EDS
% Data completeness (in resolution range)	(Not available) (30.00-3.50) 26.4 (30.00-3.50)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.31 (at 3.25Å)	Xtriage
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.266 , 0.268 0.389 , 0.389	Depositor DCC
R_{free} test set	96853 reflections (9.99%)	wwPDB-VP
Wilson B-factor (Å ²)	57.0	Xtriage
Anisotropy	0.158	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 0.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.41$, $\langle L^2 \rangle = 0.23$	Xtriage
Estimated twinning fraction	0.064 for k,h,-l 0.065 for -k,-h,-l 0.067 for h,-k,-l	Xtriage
F_o, F_c correlation	0.30	EDS
Total number of atoms	10302	wwPDB-VP
Average B, all atoms (Å ²)	62.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.47% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.35	0/5289	0.96	29/7212 (0.4%)
1	B	0.34	0/5289	0.96	27/7212 (0.4%)
All	All	0.35	0/10578	0.96	56/14424 (0.4%)

There are no bond length outliers.

All (56) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	613	LEU	N-CA-C	13.81	126.02	110.97
1	A	532	HIS	N-CA-C	-12.56	88.42	108.90
1	A	452	TYR	N-CA-C	-12.26	92.06	108.24
1	B	493	ALA	N-CA-C	9.23	123.72	109.60
1	B	615	TYR	N-CA-C	-9.20	101.82	113.23
1	B	492	LYS	N-CA-C	-8.57	95.27	109.07
1	A	448	GLU	CA-C-N	8.29	130.92	120.89
1	A	448	GLU	C-N-CA	8.29	130.92	120.89
1	A	499	ARG	N-CA-C	8.10	121.22	111.82
1	A	162	THR	N-CA-C	7.97	122.71	112.34
1	B	445	THR	N-CA-C	-7.62	101.44	108.75
1	B	111	ILE	N-CA-C	7.55	117.61	110.74
1	A	533	GLY	N-CA-C	-7.48	105.11	114.16
1	B	474	PHE	N-CA-C	7.44	126.25	109.81
1	B	643	GLN	N-CA-C	-7.25	104.41	113.55
1	A	631	LEU	N-CA-C	7.11	120.07	111.82
1	A	494	ALA	N-CA-C	-7.02	99.65	110.10
1	B	468	CYS	N-CA-C	6.76	117.17	108.24
1	B	614	ASP	N-CA-C	6.58	122.46	113.37
1	B	531	THR	N-CA-C	-6.57	104.89	113.17
1	B	450	ALA	N-CA-C	-6.26	99.89	109.41
1	B	5	VAL	N-CA-C	6.17	116.92	110.62
1	A	575	LEU	N-CA-C	6.14	115.20	108.14
1	B	471	TYR	N-CA-C	6.11	123.82	110.80
1	B	390	ILE	N-CA-C	5.98	116.10	110.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	448	GLU	N-CA-C	5.97	119.48	109.09
1	A	557	VAL	N-CA-C	-5.97	97.22	106.72
1	B	447	ILE	N-CA-C	-5.86	97.14	109.34
1	A	168	GLY	CA-C-N	5.83	126.71	120.13
1	A	168	GLY	C-N-CA	5.83	126.71	120.13
1	A	579	ASN	N-CA-C	-5.77	106.46	112.93
1	B	469	SER	CA-C-N	5.67	126.93	119.84
1	B	469	SER	C-N-CA	5.67	126.93	119.84
1	B	631	LEU	N-CA-C	-5.64	106.06	114.64
1	A	464	VAL	N-CA-C	5.60	113.42	107.76
1	B	396	ASP	N-CA-C	-5.55	106.36	113.02
1	A	531	THR	N-CA-C	-5.53	99.03	110.80
1	B	117	TYR	N-CA-C	5.48	122.48	110.80
1	B	491	ILE	N-CA-C	5.40	114.52	106.85
1	A	166	THR	N-CA-C	5.39	119.96	112.90
1	B	453	ASN	N-CA-C	5.32	122.14	110.80
1	B	454	ILE	N-CA-C	5.31	120.38	109.34
1	A	112	THR	N-CA-C	5.27	122.02	110.80
1	A	398	ALA	N-CA-C	5.21	119.12	112.87
1	A	430	TYR	CA-C-N	5.18	125.23	119.32
1	A	430	TYR	C-N-CA	5.18	125.23	119.32
1	B	585	ARG	N-CA-C	-5.18	100.01	108.76
1	B	13	SER	N-CA-C	-5.16	99.81	110.80
1	A	633	ASN	N-CA-C	-5.16	106.77	113.16
1	A	193	PHE	CA-C-N	5.12	125.03	119.76
1	A	193	PHE	C-N-CA	5.12	125.03	119.76
1	A	193	PHE	N-CA-C	5.10	114.69	108.11
1	A	250	ASP	N-CA-C	5.08	118.25	111.75
1	A	530	ASP	N-CA-C	5.05	116.53	107.80
1	B	476	VAL	N-CA-C	5.03	119.79	109.34
1	A	168	GLY	N-CA-C	5.01	122.57	112.34

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	5151	0	5012	503	0
1	B	5151	0	5012	603	0
All	All	10302	0	10024	1092	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 54.

All (1092) 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:B:16:LEU:HB2	1:B:363:THR:HG21	1.23	1.14
1:B:445:THR:HG23	1:B:446:VAL:H	1.05	1.14
1:A:15:ASP:HB2	1:A:608:ALA:HB1	1.25	1.13
1:A:144:ALA:HB3	1:A:165:ASP:HA	1.26	1.10
1:B:376:LEU:HB3	1:B:464:VAL:HG21	1.31	1.10
1:A:506:LYS:HE3	1:A:552:VAL:HG23	1.31	1.08
1:B:109:TYR:HA	1:B:113:SER:HB2	1.35	1.08
1:A:310:GLY:HA2	1:A:465:PRO:HB2	1.35	1.08
1:B:502:ALA:HB1	1:B:562:VAL:H	1.13	1.07
1:B:506:LYS:HB3	1:B:552:VAL:HG13	1.37	1.07
1:A:154:HIS:HA	1:A:448:GLU:HG3	1.38	1.05
1:B:198:VAL:HG22	1:B:199:PRO:HD2	1.33	1.05
1:A:154:HIS:HA	1:A:448:GLU:CG	1.88	1.04
1:B:10:GLN:C	1:B:12:LYS:H	1.64	1.02
1:B:390:ILE:HD12	1:B:390:ILE:H	1.21	1.01
1:B:310:GLY:HA2	1:B:465:PRO:HB2	1.43	1.01
1:A:506:LYS:HD2	1:A:558:MET:HA	1.41	1.00
1:B:14:SER:O	1:B:610:ARG:HA	1.63	0.97
1:B:450:ALA:O	1:B:452:TYR:N	1.97	0.97
1:B:299:ILE:HG22	1:B:428:VAL:HG21	1.44	0.96
1:B:154:HIS:HA	1:B:447:ILE:CG2	1.96	0.96
1:A:198:VAL:CG1	1:A:240:PHE:HA	1.95	0.94
1:B:445:THR:HG23	1:B:446:VAL:N	1.71	0.94
1:B:451:GLY:HA2	1:B:532:HIS:O	1.69	0.93
1:A:472:MET:HE3	1:A:472:MET:HA	1.50	0.92
1:B:99:ILE:HG21	1:B:104:VAL:HA	1.51	0.92
1:B:280:VAL:HA	1:B:285:LEU:HD12	1.48	0.92
1:B:613:LEU:O	1:B:616:VAL:HG12	1.69	0.91
1:A:15:ASP:HB3	1:A:610:ARG:CZ	2.00	0.91
1:B:99:ILE:HG21	1:B:104:VAL:CA	2.01	0.91
1:B:509:ALA:HB2	1:B:560:VAL:HG11	1.51	0.91
1:A:395:GLU:HG3	1:A:396:ASP:N	1.84	0.90

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:154:HIS:HA	1:B:447:ILE:HG21	1.52	0.89
1:B:394:TYR:O	1:B:644:GLY:HA3	1.72	0.89
1:B:125:MET:HE1	1:B:295:LEU:HB2	1.55	0.88
1:B:310:GLY:HA2	1:B:465:PRO:CB	2.04	0.88
1:A:260:THR:HG23	1:A:263:GLN:H	1.38	0.88
1:B:474:PHE:C	1:B:476:VAL:H	1.82	0.88
1:B:72:GLY:HA3	1:B:329:ARG:HH12	1.39	0.87
1:A:473:ILE:O	1:A:476:VAL:HG22	1.74	0.87
1:A:15:ASP:HB2	1:A:608:ALA:CB	2.06	0.86
1:B:630:GLU:HG2	1:B:632:LYS:H	1.40	0.86
1:A:332:TYR:HB2	1:A:335:LEU:HD13	1.57	0.86
1:B:507:LEU:HD22	1:B:507:LEU:H	1.40	0.86
1:B:573:VAL:HG12	1:B:574:GLU:H	1.40	0.85
1:A:396:ASP:OD2	1:B:277:ARG:HD3	1.75	0.85
1:B:526:LYS:O	1:B:561:PHE:HB2	1.77	0.85
1:B:498:LEU:H	1:B:501:GLY:HA2	1.40	0.84
1:B:420:LEU:HD13	1:B:637:SER:HA	1.58	0.83
1:B:531:THR:HG22	1:B:534:LEU:HB2	1.57	0.83
1:A:440:THR:HB	1:A:461:VAL:HG23	1.60	0.83
1:B:570:ARG:HG2	1:B:570:ARG:HH11	1.43	0.83
1:B:22:ASP:HB2	1:B:601:GLY:HA2	1.60	0.82
1:A:200:TYR:HB2	1:A:331:ARG:HG2	1.62	0.82
1:B:445:THR:CG2	1:B:446:VAL:N	2.40	0.82
1:A:73:ILE:HB	1:A:78:LEU:HD21	1.62	0.82
1:A:236:LEU:HD22	1:A:236:LEU:H	1.43	0.82
1:A:52:LYS:HG3	1:A:304:VAL:HB	1.61	0.81
1:A:147:ILE:HG21	1:A:159:LEU:HB2	1.63	0.81
1:B:72:GLY:HA3	1:B:329:ARG:NH1	1.95	0.81
1:B:128:LEU:HD22	1:B:296:LEU:HD23	1.62	0.81
1:A:198:VAL:HG11	1:A:240:PHE:HA	1.64	0.80
1:B:446:VAL:O	1:B:447:ILE:HG12	1.81	0.80
1:B:496:LYS:HG2	1:B:503:VAL:HG11	1.63	0.80
1:B:22:ASP:HB2	1:B:601:GLY:CA	2.11	0.80
1:B:199:PRO:HG2	1:B:240:PHE:HB3	1.63	0.80
1:A:123:LEU:HD13	1:A:218:MET:HE3	1.63	0.79
1:A:312:ALA:HB2	1:A:461:VAL:HG21	1.65	0.78
1:B:277:ARG:HA	1:B:351:TRP:HH2	1.48	0.78
1:B:126:ASN:HA	1:B:129:ARG:HD3	1.65	0.78
1:B:353:ALA:HA	1:B:616:VAL:CG2	2.14	0.78
1:B:418:ALA:HB3	1:B:638:VAL:HG12	1.66	0.78
1:B:5:VAL:HG11	1:B:360:LEU:HD11	1.63	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:577:ARG:HD2	1:A:579:ASN:OD1	1.84	0.77
1:A:285:LEU:HD21	1:A:329:ARG:HD2	1.67	0.77
1:A:277:ARG:HB2	1:A:355:MET:HE2	1.65	0.77
1:A:506:LYS:CE	1:A:552:VAL:HG23	2.14	0.77
1:A:260:THR:CG2	1:A:263:GLN:H	1.98	0.77
1:B:471:TYR:HA	1:B:474:PHE:HB2	1.68	0.77
1:A:381:VAL:HG22	1:A:553:THR:HG22	1.67	0.76
1:B:121:VAL:HG13	1:B:292:ALA:HB2	1.65	0.76
1:A:123:LEU:HB2	1:A:218:MET:HE1	1.66	0.76
1:B:153:GLY:O	1:B:447:ILE:HG21	1.84	0.76
1:B:502:ALA:HB1	1:B:562:VAL:N	1.97	0.76
1:B:496:LYS:O	1:B:503:VAL:HB	1.86	0.75
1:A:202:ASP:OD2	1:A:204:ARG:HD3	1.86	0.75
1:A:127:MET:HE3	1:A:276:LEU:HG	1.69	0.75
1:A:216:MET:HA	1:A:216:MET:HE2	1.67	0.75
1:B:318:ALA:HB2	1:B:430:TYR:CZ	2.22	0.75
1:B:10:GLN:C	1:B:12:LYS:N	2.39	0.75
1:A:443:PRO:HA	1:A:458:HIS:HB3	1.69	0.74
1:B:355:MET:HE3	1:B:355:MET:HA	1.67	0.74
1:B:144:ALA:HB3	1:B:165:ASP:HA	1.69	0.74
1:B:156:LYS:HG3	1:B:445:THR:HG21	1.69	0.74
1:A:137:THR:HG21	1:A:172:PHE:H	1.52	0.74
1:A:393:SER:OG	1:A:395:GLU:HG2	1.88	0.74
1:B:14:SER:HB2	1:B:611:ILE:O	1.87	0.74
1:B:14:SER:O	1:B:610:ARG:CA	2.36	0.74
1:B:223:ARG:NH1	1:B:231:TYR:HA	2.01	0.74
1:B:423:MET:HB2	1:B:426:VAL:HG21	1.70	0.74
1:A:249:GLN:O	1:A:253:GLU:HG3	1.88	0.73
1:B:150:TYR:CE2	1:B:308:ALA:HB2	2.23	0.73
1:A:70:LEU:HD22	1:A:328:ILE:HD11	1.68	0.73
1:A:267:PRO:HB2	1:A:271:VAL:CG1	2.18	0.73
1:A:576:PRO:O	1:A:578:LEU:HG	1.88	0.73
1:B:526:LYS:HA	1:B:526:LYS:HZ3	1.51	0.73
1:A:574:GLU:O	1:A:575:LEU:HD23	1.87	0.73
1:B:22:ASP:CB	1:B:601:GLY:HA2	2.18	0.73
1:B:630:GLU:HG2	1:B:632:LYS:N	2.03	0.73
1:B:1:MET:HE1	1:B:361:VAL:HA	1.71	0.73
1:B:65:GLU:HG3	1:B:117:TYR:CE1	2.23	0.73
1:B:526:LYS:HZ2	1:B:527:VAL:H	1.37	0.73
1:A:55:LEU:HB2	1:A:227:LEU:HD12	1.70	0.73
1:B:120:HIS:O	1:B:124:LEU:HG	1.88	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:75:LYS:HG2	1:A:341:ALA:HB2	1.71	0.72
1:B:284:ARG:HD3	1:B:340:ALA:HB2	1.71	0.72
1:A:632:LYS:HG3	1:B:334:PHE:HB2	1.71	0.72
1:A:198:VAL:HG13	1:A:199:PRO:HD2	1.72	0.72
1:B:283:ASN:O	1:B:284:ARG:HB2	1.90	0.71
1:A:155:VAL:HA	1:A:447:ILE:HG12	1.70	0.71
1:A:444:ILE:HD13	1:A:457:ASP:O	1.90	0.71
1:B:390:ILE:H	1:B:390:ILE:CD1	1.94	0.71
1:A:531:THR:HG23	1:A:531:THR:O	1.89	0.71
1:A:423:MET:HB3	1:A:426:VAL:HG21	1.72	0.71
1:B:22:ASP:CG	1:B:601:GLY:HA2	2.16	0.71
1:B:328:ILE:HG22	1:B:328:ILE:O	1.89	0.71
1:A:189:PHE:CZ	1:A:191:GLU:HB2	2.26	0.71
1:B:469:SER:HB2	1:B:470:PRO:HD2	1.72	0.71
1:B:493:ALA:HB2	1:B:584:PHE:CD1	2.26	0.71
1:B:368:VAL:HG13	1:B:410:LEU:HD21	1.72	0.71
1:B:410:LEU:O	1:B:414:THR:HG22	1.91	0.71
1:A:299:ILE:HG22	1:A:428:VAL:HG11	1.74	0.70
1:B:125:MET:HE2	1:B:292:ALA:HA	1.74	0.70
1:B:348:LEU:HD12	1:B:618:PRO:HB3	1.71	0.70
1:A:549:PRO:O	1:A:552:VAL:HG12	1.91	0.70
1:A:395:GLU:CB	1:A:436:PHE:HE1	2.04	0.70
1:A:557:VAL:HG13	1:A:557:VAL:O	1.91	0.70
1:A:15:ASP:HB3	1:A:610:ARG:NH2	2.06	0.70
1:A:157:ILE:HG22	1:A:444:ILE:HA	1.74	0.70
1:B:548:ILE:HD12	1:B:548:ILE:H	1.57	0.70
1:A:150:TYR:HB3	1:A:157:ILE:HG13	1.72	0.70
1:A:267:PRO:HB2	1:A:271:VAL:HG11	1.74	0.70
1:B:259:ARG:HG3	1:B:263:GLN:OE1	1.91	0.70
1:A:5:VAL:HG11	1:A:360:LEU:HD11	1.74	0.69
1:B:630:GLU:HG2	1:B:631:LEU:N	2.06	0.69
1:A:441:GLN:HE22	1:A:460:VAL:HG12	1.57	0.69
1:B:380:ARG:HA	1:B:436:PHE:CD2	2.27	0.69
1:B:91:ASN:HA	1:B:94:ARG:HD2	1.74	0.69
1:B:156:LYS:HE3	1:B:445:THR:HB	1.75	0.69
1:B:55:LEU:HA	1:B:301:MET:HB2	1.75	0.69
1:A:3:ARG:O	1:A:7:LYS:HG2	1.93	0.69
1:A:199:PRO:HG2	1:A:240:PHE:HB3	1.74	0.69
1:A:260:THR:HG22	1:A:263:GLN:HB2	1.74	0.69
1:B:131:TYR:O	1:B:135:VAL:HG23	1.92	0.69
1:A:154:HIS:HA	1:A:448:GLU:HG2	1.71	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:526:LYS:HA	1:B:526:LYS:NZ	2.07	0.69
1:A:629:GLY:H	1:A:650:LEU:HD11	1.58	0.69
1:B:523:THR:HB	1:B:565:ILE:HG22	1.74	0.69
1:A:378:LEU:HG	1:A:478:ALA:HB1	1.75	0.69
1:B:299:ILE:HG22	1:B:428:VAL:CG2	2.22	0.69
1:B:70:LEU:O	1:B:73:ILE:HG23	1.93	0.68
1:B:175:PRO:O	1:B:222:HIS:HB2	1.94	0.68
1:B:198:VAL:CG2	1:B:199:PRO:HD2	2.19	0.68
1:A:452:TYR:HB3	1:A:459:LEU:HD22	1.74	0.68
1:A:494:ALA:HB2	1:A:504:TYR:CZ	2.27	0.68
1:B:277:ARG:HA	1:B:351:TRP:CH2	2.28	0.68
1:B:598:VAL:O	1:B:606:TYR:HB2	1.93	0.68
1:A:320:VAL:HG12	1:A:428:VAL:HG22	1.73	0.68
1:B:474:PHE:O	1:B:476:VAL:HG23	1.93	0.68
1:B:211:VAL:O	1:B:215:LEU:HD12	1.94	0.68
1:A:342:LEU:N	1:A:342:LEU:HD23	2.09	0.67
1:A:395:GLU:HB2	1:A:436:PHE:CE1	2.30	0.67
1:A:576:PRO:HG2	1:A:578:LEU:HD21	1.74	0.67
1:B:137:THR:HG23	1:B:171:GLN:HA	1.76	0.67
1:A:532:HIS:ND1	1:A:532:HIS:O	2.27	0.67
1:A:210:GLU:HG2	1:A:275:ALA:HB2	1.76	0.67
1:A:73:ILE:HG23	1:A:335:LEU:O	1.94	0.67
1:A:126:ASN:HB3	1:A:217:MET:HE3	1.75	0.67
1:A:379:ARG:NH2	1:A:395:GLU:O	2.22	0.67
1:B:499:ARG:HB2	1:B:500:LYS:HD2	1.77	0.67
1:A:49:VAL:HA	1:A:542:GLY:O	1.95	0.67
1:B:154:HIS:HD2	1:B:448:GLU:HG3	1.59	0.67
1:A:254:TRP:HA	1:A:259:ARG:HB3	1.77	0.67
1:B:525:PHE:HB2	1:B:561:PHE:O	1.94	0.67
1:A:94:ARG:HE	1:A:104:VAL:HG21	1.60	0.67
1:A:496:LYS:HB3	1:A:503:VAL:HG22	1.77	0.67
1:B:602:ALA:O	1:B:603:HIS:HB3	1.93	0.67
1:A:490:VAL:HG22	1:A:588:GLU:HB2	1.76	0.66
1:B:513:ALA:HA	1:B:523:THR:HG21	1.77	0.66
1:A:37:LYS:HA	1:A:41:LEU:O	1.95	0.66
1:A:70:LEU:HB3	1:A:73:ILE:HD12	1.76	0.66
1:A:184:PRO:HB3	1:A:237:ALA:O	1.95	0.66
1:A:482:ALA:O	1:A:511:LYS:HD2	1.95	0.66
1:B:291:THR:HG21	1:B:423:MET:CG	2.27	0.65
1:A:375:GLY:HA2	1:A:413:CYS:HB2	1.78	0.65
1:B:490:VAL:HG22	1:B:491:ILE:N	2.12	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:35:ASP:O	1:A:589:VAL:HA	1.96	0.65
1:A:383:LYS:HG2	1:A:383:LYS:O	1.97	0.65
1:A:535:THR:O	1:A:536:LYS:CB	2.44	0.65
1:B:1:MET:CE	1:B:361:VAL:HA	2.26	0.65
1:A:348:LEU:HD11	1:A:623:ILE:HG21	1.79	0.65
1:A:354:ILE:HD13	1:A:360:LEU:HB3	1.77	0.65
1:B:526:LYS:HD3	1:B:527:VAL:N	2.12	0.65
1:A:535:THR:O	1:A:536:LYS:HB2	1.95	0.64
1:A:403:GLU:HG3	1:A:634:TYR:CZ	2.33	0.64
1:B:254:TRP:HE1	1:B:261:ASP:HA	1.59	0.64
1:B:217:MET:HA	1:B:217:MET:HE2	1.80	0.64
1:B:10:GLN:O	1:B:12:LYS:N	2.28	0.64
1:B:117:TYR:O	1:B:118:ASP:O	2.14	0.64
1:B:351:TRP:O	1:B:355:MET:HB2	1.98	0.64
1:A:395:GLU:CB	1:A:436:PHE:CE1	2.81	0.64
1:B:291:THR:HG21	1:B:423:MET:HG2	1.78	0.64
1:B:400:LEU:HD23	1:B:434:LEU:HD11	1.78	0.64
1:B:565:ILE:HD13	1:B:565:ILE:H	1.61	0.64
1:A:149:MET:HE2	1:A:156:LYS:HZ2	1.63	0.64
1:B:57:GLY:O	1:B:59:PRO:HD3	1.97	0.64
1:B:407:GLN:HB3	1:B:638:VAL:HG12	1.79	0.64
1:B:527:VAL:CG1	1:B:531:THR:HG23	2.28	0.64
1:A:149:MET:SD	1:A:151:HIS:HB3	2.38	0.64
1:A:27:VAL:HG23	1:A:51:GLN:C	2.23	0.64
1:B:379:ARG:CZ	1:B:436:PHE:HB3	2.28	0.64
1:B:284:ARG:NH2	1:B:336:LEU:O	2.31	0.63
1:A:358:PRO:HD2	1:A:359:GLU:OE2	1.98	0.63
1:A:300:MET:HG2	1:A:301:MET:HE3	1.80	0.63
1:A:420:LEU:HG	1:A:636:GLY:O	1.99	0.63
1:B:531:THR:HG22	1:B:534:LEU:CB	2.25	0.63
1:B:593:ILE:HD12	1:B:593:ILE:H	1.64	0.63
1:A:379:ARG:HD2	1:A:399:PHE:CE1	2.33	0.63
1:B:375:GLY:HA3	1:B:409:ALA:O	1.98	0.63
1:A:82:GLY:HA3	1:B:95:SER:HB3	1.80	0.63
1:A:186:TRP:HD1	1:A:215:LEU:HD12	1.64	0.63
1:A:307:CYS:HB2	1:A:309:GLU:OE1	1.98	0.63
1:B:68:SER:HB2	1:B:117:TYR:C	2.24	0.63
1:B:72:GLY:H	1:B:329:ARG:HH22	1.45	0.62
1:B:353:ALA:HA	1:B:616:VAL:HG21	1.79	0.62
1:B:438:ALA:O	1:B:463:GLY:N	2.33	0.62
1:A:377:TYR:CE1	1:A:466:VAL:HG22	2.34	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:150:TYR:HB2	1:B:311:TYR:CE1	2.35	0.62
1:B:156:LYS:HG3	1:B:445:THR:CG2	2.29	0.62
1:B:327:SER:O	1:B:328:ILE:HD12	2.00	0.62
1:A:8:ASN:O	1:A:475:PRO:HG2	2.00	0.62
1:A:650:LEU:CD1	1:A:651:THR:H	2.13	0.62
1:B:22:ASP:HB2	1:B:601:GLY:N	2.15	0.62
1:B:213:PHE:CE2	1:B:272:MET:HE3	2.35	0.62
1:A:29:HIS:HB2	1:A:597:HIS:CE1	2.35	0.62
1:B:75:LYS:HA	1:B:341:ALA:HB2	1.81	0.62
1:B:294:GLN:HE22	1:B:419:PRO:HB2	1.65	0.62
1:B:418:ALA:HB3	1:B:638:VAL:CG1	2.29	0.62
1:A:483:ASN:O	1:A:486:CYS:HB2	1.99	0.61
1:B:117:TYR:O	1:B:117:TYR:CD1	2.53	0.61
1:B:332:TYR:HB2	1:B:335:LEU:HD13	1.82	0.61
1:A:235:GLN:HE21	1:A:238:ASP:HA	1.64	0.61
1:A:371:ALA:HB3	1:A:410:LEU:HD23	1.82	0.61
1:B:506:LYS:CB	1:B:552:VAL:HG13	2.22	0.61
1:B:73:ILE:HG13	1:B:74:ALA:N	2.16	0.61
1:B:453:ASN:OD1	1:B:453:ASN:O	2.18	0.61
1:A:186:TRP:CD1	1:A:215:LEU:HD12	2.35	0.61
1:B:493:ALA:HB2	1:B:584:PHE:CE1	2.34	0.61
1:A:126:ASN:HB3	1:A:217:MET:CE	2.30	0.61
1:A:218:MET:HE2	1:A:236:LEU:CB	2.30	0.61
1:B:155:VAL:N	1:B:447:ILE:HB	2.15	0.61
1:A:28:ALA:O	1:A:50:SER:HA	2.00	0.61
1:A:379:ARG:HG3	1:A:394:TYR:HB3	1.81	0.61
1:A:439:VAL:HG12	1:A:462:VAL:HG22	1.83	0.61
1:B:348:LEU:HD12	1:B:618:PRO:CB	2.30	0.61
1:B:109:TYR:O	1:B:114:TRP:HD1	1.84	0.61
1:B:283:ASN:HA	1:B:336:LEU:HD13	1.83	0.61
1:B:359:GLU:O	1:B:363:THR:HG23	2.01	0.61
1:A:310:GLY:HA2	1:A:465:PRO:CB	2.21	0.61
1:A:472:MET:HA	1:A:472:MET:CE	2.27	0.61
1:B:476:VAL:HG12	1:B:477:ALA:N	2.16	0.61
1:B:7:LYS:HD2	1:B:10:GLN:NE2	2.15	0.60
1:B:112:THR:CG2	1:B:197:ASP:HA	2.31	0.60
1:B:358:PRO:HD2	1:B:359:GLU:OE2	2.01	0.60
1:A:194:PRO:HD3	1:A:331:ARG:O	2.01	0.60
1:A:204:ARG:HD2	1:A:243:ARG:HG2	1.82	0.60
1:B:381:VAL:HG22	1:B:553:THR:CG2	2.31	0.60
1:A:506:LYS:HE3	1:A:552:VAL:CG2	2.21	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:401:GLN:HB3	1:A:404:THR:OG1	2.01	0.60
1:B:15:ASP:HB3	1:B:610:ARG:HG2	1.82	0.60
1:B:453:ASN:OD1	1:B:453:ASN:C	2.43	0.60
1:A:8:ASN:O	1:A:475:PRO:CG	2.50	0.60
1:A:524:HIS:HB2	1:A:563:THR:O	2.01	0.60
1:B:455:VAL:HG12	1:B:456:ASP:OD1	2.02	0.60
1:B:476:VAL:O	1:B:477:ALA:HB3	1.99	0.60
1:A:37:LYS:HG2	1:A:42:VAL:HG22	1.84	0.60
1:A:75:LYS:HD3	1:A:624:GLN:NE2	2.16	0.60
1:A:56:VAL:HA	1:A:227:LEU:HA	1.82	0.60
1:B:277:ARG:HB2	1:B:355:MET:SD	2.42	0.60
1:A:35:ASP:OD2	1:A:44:ASN:HB3	2.02	0.60
1:A:129:ARG:O	1:A:133:LEU:HD23	2.00	0.60
1:B:58:ASN:O	1:B:317:ASP:HB3	2.01	0.60
1:B:509:ALA:CB	1:B:560:VAL:HG11	2.27	0.60
1:A:100:PRO:HG2	1:A:103:ALA:HB2	1.83	0.60
1:B:68:SER:HB2	1:B:117:TYR:O	2.02	0.60
1:B:341:ALA:HA	1:B:624:GLN:HB3	1.81	0.60
1:A:197:ASP:HA	1:A:331:ARG:HH12	1.67	0.59
1:A:549:PRO:O	1:A:552:VAL:CG1	2.50	0.59
1:A:236:LEU:H	1:A:236:LEU:CD2	2.15	0.59
1:A:150:TYR:HB2	1:A:311:TYR:CZ	2.37	0.59
1:A:223:ARG:HE	1:A:231:TYR:HA	1.67	0.59
1:A:227:LEU:HD22	1:A:227:LEU:H	1.66	0.59
1:A:366:MET:O	1:A:370:VAL:HG12	2.02	0.59
1:B:470:PRO:O	1:B:473:ILE:HD13	2.01	0.59
1:B:55:LEU:HD12	1:B:55:LEU:H	1.67	0.59
1:B:444:ILE:HG22	1:B:445:THR:N	2.17	0.59
1:B:577:ARG:HG3	1:B:579:ASN:OD1	2.02	0.59
1:B:394:TYR:CE1	1:B:644:GLY:HA2	2.37	0.59
1:A:55:LEU:H	1:A:55:LEU:HD22	1.67	0.59
1:A:74:ALA:HB2	1:A:337:SER:O	2.02	0.59
1:A:111:ILE:HG23	1:A:111:ILE:O	2.00	0.59
1:B:38:PHE:O	1:B:39:ASP:HB2	2.02	0.59
1:B:381:VAL:HG11	1:B:386:PHE:CZ	2.38	0.59
1:A:441:GLN:NE2	1:A:460:VAL:HG12	2.18	0.59
1:B:17:PHE:O	1:B:359:GLU:HB3	2.03	0.59
1:B:199:PRO:HG2	1:B:240:PHE:CB	2.32	0.59
1:A:417:ASP:HB3	1:A:637:SER:OG	2.03	0.59
1:B:52:LYS:HB2	1:B:304:VAL:HB	1.82	0.59
1:A:443:PRO:CA	1:A:458:HIS:HB3	2.33	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:150:TYR:HB3	1:A:157:ILE:CG1	2.33	0.58
1:B:474:PHE:O	1:B:476:VAL:N	2.36	0.58
1:B:235:GLN:HE21	1:B:238:ASP:HA	1.68	0.58
1:A:56:VAL:HG13	1:A:226:ASN:HB3	1.85	0.58
1:A:33:ARG:HB2	1:A:46:VAL:HG12	1.84	0.58
1:A:198:VAL:HG13	1:A:240:PHE:HA	1.83	0.58
1:A:638:VAL:HG23	1:A:646:GLY:C	2.28	0.58
1:A:414:THR:HB	1:A:416:GLN:HE21	1.69	0.58
1:A:442:VAL:C	1:A:458:HIS:HB3	2.28	0.58
1:A:610:ARG:HH11	1:A:610:ARG:HG2	1.68	0.58
1:B:13:SER:O	1:B:14:SER:OG	2.12	0.58
1:B:364:TYR:O	1:B:368:VAL:HG23	2.04	0.58
1:B:23:ARG:HH11	1:B:23:ARG:HB2	1.69	0.58
1:B:149:MET:HE1	1:B:156:LYS:HB2	1.86	0.58
1:B:376:LEU:HD23	1:B:464:VAL:HG11	1.86	0.58
1:A:199:PRO:O	1:A:240:PHE:HB2	2.03	0.58
1:A:323:PRO:HD3	1:A:426:VAL:HG12	1.86	0.58
1:B:201:LEU:N	1:B:201:LEU:HD23	2.19	0.58
1:A:267:PRO:O	1:A:272:MET:HE3	2.02	0.58
1:A:632:LYS:HG3	1:B:334:PHE:CB	2.33	0.58
1:B:59:PRO:HD2	1:B:223:ARG:HH22	1.66	0.58
1:A:119:ASN:HD21	1:A:323:PRO:HG2	1.69	0.58
1:A:650:LEU:HD12	1:A:651:THR:H	1.69	0.58
1:B:15:ASP:HB2	1:B:608:ALA:HB1	1.85	0.58
1:A:371:ALA:CB	1:A:410:LEU:HD23	2.34	0.57
1:A:459:LEU:C	1:A:459:LEU:HD12	2.29	0.57
1:B:154:HIS:HA	1:B:447:ILE:CB	2.34	0.57
1:A:149:MET:SD	1:A:156:LYS:HG3	2.45	0.57
1:B:476:VAL:O	1:B:477:ALA:CB	2.51	0.57
1:A:483:ASN:HB2	1:A:484:PRO:HD2	1.86	0.57
1:B:491:ILE:O	1:B:586:SER:HB2	2.04	0.57
1:B:498:LEU:N	1:B:501:GLY:HA2	2.17	0.57
1:B:117:TYR:CD2	1:B:326:GLY:HA3	2.40	0.57
1:B:15:ASP:HB2	1:B:608:ALA:CB	2.34	0.57
1:B:249:GLN:O	1:B:253:GLU:HG3	2.04	0.57
1:A:120:HIS:O	1:A:124:LEU:HG	2.04	0.57
1:A:386:PHE:CE2	1:A:507:LEU:HG	2.39	0.57
1:B:207:THR:O	1:B:211:VAL:HG23	2.05	0.57
1:A:92:GLU:O	1:A:96:THR:HG22	2.03	0.57
1:A:550:GLU:C	1:A:550:GLU:CD	2.72	0.57
1:B:517:ARG:NE	1:B:539:ALA:HB3	2.19	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:218:MET:HG2	1:A:240:PHE:CE2	2.40	0.57
1:A:577:ARG:NH2	1:A:580:SER:HB2	2.19	0.57
1:A:178:ARG:HG3	1:A:254:TRP:CH2	2.39	0.57
1:B:380:ARG:NH1	1:B:380:ARG:HG3	2.19	0.57
1:B:483:ASN:HB2	1:B:484:PRO:CD	2.35	0.57
1:B:514:TRP:CD1	1:B:546:THR:H	2.22	0.57
1:B:490:VAL:HG22	1:B:491:ILE:H	1.69	0.56
1:B:126:ASN:HA	1:B:129:ARG:CD	2.34	0.56
1:B:381:VAL:HG22	1:B:553:THR:HG23	1.85	0.56
1:B:506:LYS:HE2	1:B:552:VAL:O	2.05	0.56
1:A:375:GLY:HA3	1:A:409:ALA:O	2.04	0.56
1:A:440:THR:HB	1:A:461:VAL:CG2	2.32	0.56
1:A:645:LEU:HD23	1:A:645:LEU:H	1.70	0.56
1:B:155:VAL:HG13	1:B:444:ILE:HG23	1.87	0.56
1:B:305:PRO:HA	1:B:467:ALA:O	2.05	0.56
1:B:117:TYR:O	1:B:117:TYR:CG	2.58	0.56
1:B:423:MET:HB2	1:B:426:VAL:CG2	2.35	0.56
1:B:552:VAL:O	1:B:552:VAL:HG12	2.05	0.56
1:A:442:VAL:O	1:A:458:HIS:CB	2.53	0.56
1:B:253:GLU:CB	1:B:259:ARG:HB2	2.36	0.56
1:B:283:ASN:O	1:B:284:ARG:CB	2.53	0.56
1:B:380:ARG:HG3	1:B:380:ARG:HH11	1.69	0.56
1:B:428:VAL:HG13	1:B:428:VAL:O	2.04	0.56
1:A:403:GLU:HG3	1:A:634:TYR:CE2	2.40	0.56
1:B:69:TYR:HB2	1:B:288:GLN:HE21	1.71	0.56
1:B:473:ILE:HG12	1:B:473:ILE:O	2.05	0.56
1:B:473:ILE:HD13	1:B:473:ILE:H	1.71	0.56
1:A:131:TYR:O	1:A:135:VAL:HG23	2.06	0.56
1:A:133:LEU:HB3	1:A:172:PHE:CD1	2.40	0.56
1:A:476:VAL:O	1:A:485:TYR:O	2.24	0.56
1:B:10:GLN:O	1:B:11:ASP:CG	2.49	0.56
1:B:150:TYR:O	1:B:156:LYS:HA	2.05	0.56
1:B:19:ILE:HB	1:B:136:LEU:HD11	1.88	0.56
1:B:161:VAL:HG22	1:B:226:ASN:CG	2.30	0.56
1:A:305:PRO:HG2	1:A:311:TYR:CD1	2.41	0.55
1:A:470:PRO:HG2	1:A:545:TRP:CH2	2.40	0.55
1:A:484:PRO:HD2	1:A:510:TRP:CZ3	2.41	0.55
1:A:532:HIS:ND1	1:A:532:HIS:C	2.65	0.55
1:B:55:LEU:HD12	1:B:55:LEU:N	2.22	0.55
1:A:254:TRP:HE1	1:A:261:ASP:HA	1.71	0.55
1:A:423:MET:HB3	1:A:426:VAL:CG2	2.35	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:328:ILE:O	1:B:328:ILE:CG2	2.55	0.55
1:A:557:VAL:O	1:A:557:VAL:CG1	2.55	0.55
1:B:355:MET:HA	1:B:355:MET:CE	2.36	0.55
1:B:510:TRP:CE2	1:B:548:ILE:HG13	2.42	0.55
1:B:2:LEU:HD23	1:B:360:LEU:HD22	1.89	0.55
1:B:150:TYR:CE2	1:B:152:ASP:HB3	2.40	0.55
1:B:628:ALA:HA	1:B:651:THR:O	2.07	0.55
1:B:16:LEU:HD13	1:B:359:GLU:HB2	1.88	0.55
1:B:125:MET:HE1	1:B:295:LEU:CB	2.34	0.55
1:B:151:HIS:ND1	1:B:156:LYS:HB3	2.21	0.55
1:B:283:ASN:N	1:B:283:ASN:HD22	2.03	0.55
1:A:386:PHE:CE1	1:A:552:VAL:HG21	2.41	0.55
1:A:419:PRO:HA	1:A:637:SER:HB3	1.87	0.55
1:B:152:ASP:C	1:B:154:HIS:H	2.14	0.55
1:A:218:MET:HE2	1:A:236:LEU:HB2	1.88	0.55
1:B:14:SER:O	1:B:611:ILE:N	2.39	0.55
1:A:577:ARG:HH21	1:A:580:SER:HB2	1.72	0.55
1:B:93:LEU:HA	1:B:96:THR:OG1	2.06	0.55
1:B:295:LEU:O	1:B:299:ILE:HG23	2.07	0.55
1:B:446:VAL:HG22	1:B:447:ILE:N	2.21	0.55
1:A:259:ARG:HD2	1:A:263:GLN:HB3	1.88	0.54
1:A:516:LEU:CD1	1:A:521:TYR:HB2	2.37	0.54
1:B:483:ASN:HB2	1:B:484:PRO:HD2	1.88	0.54
1:B:514:TRP:O	1:B:518:VAL:HG23	2.06	0.54
1:A:7:LYS:HD3	1:A:11:ASP:OD1	2.07	0.54
1:A:55:LEU:HD22	1:A:55:LEU:N	2.22	0.54
1:B:507:LEU:HD22	1:B:507:LEU:N	2.18	0.54
1:A:277:ARG:HB2	1:A:355:MET:CE	2.37	0.54
1:B:104:VAL:HG22	1:B:105:ALA:N	2.23	0.54
1:A:188:GLN:HG2	1:A:247:THR:OG1	2.07	0.54
1:B:57:GLY:O	1:B:228:ALA:HA	2.06	0.54
1:A:604:ALA:HB1	1:A:606:TYR:CE1	2.42	0.54
1:B:442:VAL:HG23	1:B:459:LEU:O	2.07	0.54
1:A:45:ARG:C	1:A:45:ARG:HD2	2.33	0.54
1:A:55:LEU:CB	1:A:227:LEU:HD12	2.37	0.54
1:A:70:LEU:CD2	1:A:328:ILE:HD11	2.38	0.54
1:B:210:GLU:O	1:B:214:VAL:HG23	2.08	0.54
1:A:159:LEU:HD11	1:A:315:MET:HE3	1.90	0.54
1:B:611:ILE:O	1:B:611:ILE:HG23	2.07	0.54
1:A:61:VAL:HA	1:A:320:VAL:HG22	1.90	0.54
1:A:477:ALA:HA	1:A:485:TYR:HB3	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:454:ILE:N	1:B:454:ILE:HD12	2.22	0.54
1:B:527:VAL:HG13	1:B:531:THR:HG23	1.90	0.54
1:B:204:ARG:HG3	1:B:243:ARG:O	2.08	0.53
1:B:571:HIS:CE1	1:B:573:VAL:O	2.61	0.53
1:A:200:TYR:CD1	1:A:331:ARG:HD3	2.43	0.53
1:B:126:ASN:HA	1:B:129:ARG:HG2	1.91	0.53
1:B:260:THR:OG1	1:B:263:GLN:HG3	2.07	0.53
1:B:544:THR:HG22	1:B:544:THR:O	2.09	0.53
1:A:506:LYS:NZ	1:A:557:VAL:O	2.39	0.53
1:B:71:GLU:HG2	1:B:287:ASN:HD22	1.73	0.53
1:B:521:TYR:HA	1:B:568:ARG:HG3	1.90	0.53
1:A:260:THR:HG22	1:A:263:GLN:CB	2.37	0.53
1:B:99:ILE:CG2	1:B:104:VAL:HB	2.38	0.53
1:B:99:ILE:HB	1:B:104:VAL:HB	1.90	0.53
1:A:490:VAL:HG22	1:A:588:GLU:CB	2.37	0.53
1:B:603:HIS:O	1:B:604:ALA:HB3	2.07	0.53
1:A:86:ILE:HG13	1:B:100:PRO:HB3	1.91	0.53
1:A:368:VAL:HG13	1:A:410:LEU:HD11	1.91	0.53
1:A:512:LEU:O	1:A:516:LEU:HB2	2.08	0.53
1:B:402:PRO:CB	1:B:429:THR:HG21	2.39	0.53
1:B:469:SER:CB	1:B:470:PRO:HD2	2.39	0.53
1:B:500:LYS:HD2	1:B:500:LYS:N	2.23	0.53
1:A:178:ARG:HG3	1:A:254:TRP:CZ2	2.43	0.53
1:A:355:MET:HE3	1:A:355:MET:HA	1.90	0.53
1:B:513:ALA:HA	1:B:523:THR:CG2	2.38	0.53
1:B:147:ILE:HD12	1:B:147:ILE:H	1.74	0.53
1:B:307:CYS:HB3	1:B:538:TYR:CE2	2.43	0.53
1:B:382:LYS:HB3	1:B:394:TYR:CZ	2.44	0.53
1:B:113:SER:O	1:B:116:TRP:NE1	2.42	0.53
1:B:490:VAL:HG23	1:B:587:VAL:O	2.09	0.53
1:B:56:VAL:HG22	1:B:145:GLY:HA3	1.91	0.52
1:B:493:ALA:HB1	1:B:497:TYR:OH	2.08	0.52
1:A:64:ASN:O	1:A:323:PRO:HA	2.10	0.52
1:A:207:THR:HG22	1:A:210:GLU:CG	2.39	0.52
1:B:622:GLY:O	1:B:623:ILE:HG13	2.09	0.52
1:B:630:GLU:HG2	1:B:631:LEU:H	1.71	0.52
1:A:364:TYR:O	1:A:368:VAL:HG23	2.10	0.52
1:B:254:TRP:HA	1:B:259:ARG:HB3	1.91	0.52
1:B:310:GLY:CA	1:B:465:PRO:HB2	2.28	0.52
1:B:614:ASP:HB3	1:B:617:LYS:HE3	1.90	0.52
1:A:312:ALA:CB	1:A:461:VAL:HG21	2.37	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:283:ASN:HA	1:B:336:LEU:CD1	2.40	0.52
1:B:568:ARG:HH11	1:B:568:ARG:HB3	1.74	0.52
1:A:16:LEU:HA	1:A:363:THR:HG21	1.92	0.52
1:A:227:LEU:HD22	1:A:227:LEU:N	2.24	0.52
1:A:375:GLY:CA	1:A:413:CYS:HB2	2.38	0.52
1:A:483:ASN:ND2	1:A:486:CYS:SG	2.82	0.52
1:B:155:VAL:CG1	1:B:444:ILE:HG23	2.40	0.52
1:B:406:VAL:O	1:B:410:LEU:HD13	2.10	0.52
1:B:442:VAL:O	1:B:458:HIS:HA	2.09	0.52
1:A:111:ILE:O	1:A:111:ILE:CG2	2.58	0.52
1:A:527:VAL:HG12	1:A:560:VAL:HA	1.90	0.52
1:A:638:VAL:HG23	1:A:646:GLY:O	2.10	0.52
1:B:17:PHE:H	1:B:363:THR:CG2	2.22	0.52
1:B:175:PRO:HB2	1:B:221:TRP:HA	1.91	0.52
1:A:441:GLN:HE22	1:A:460:VAL:CG1	2.21	0.52
1:B:280:VAL:HB	1:B:351:TRP:CH2	2.45	0.52
1:B:570:ARG:HG2	1:B:570:ARG:NH1	2.14	0.52
1:A:195:SER:C	1:A:196:ILE:HG13	2.34	0.52
1:A:342:LEU:HD23	1:A:342:LEU:H	1.74	0.52
1:B:493:ALA:HB2	1:B:584:PHE:CG	2.44	0.52
1:A:273:LEU:HD23	1:A:277:ARG:NH1	2.25	0.52
1:A:459:LEU:HD12	1:A:459:LEU:O	2.10	0.52
1:B:573:VAL:HG12	1:B:574:GLU:N	2.19	0.52
1:A:154:HIS:HB3	1:A:448:GLU:HB2	1.91	0.52
1:A:610:ARG:HG2	1:A:610:ARG:NH1	2.25	0.52
1:B:134:GLN:HG3	1:B:272:MET:SD	2.50	0.52
1:B:394:TYR:HA	1:B:643:GLN:O	2.09	0.52
1:B:650:LEU:HD23	1:B:650:LEU:H	1.76	0.51
1:A:125:MET:HE1	1:A:295:LEU:CB	2.41	0.51
1:A:277:ARG:HD3	1:A:355:MET:HE2	1.91	0.51
1:A:377:TYR:HD1	1:A:464:VAL:O	1.93	0.51
1:A:386:PHE:CD1	1:A:552:VAL:HG21	2.46	0.51
1:A:575:LEU:HB3	1:A:576:PRO:CD	2.40	0.51
1:B:152:ASP:C	1:B:154:HIS:N	2.67	0.51
1:B:61:VAL:HG11	1:B:231:TYR:HD1	1.75	0.51
1:B:156:LYS:H	1:B:445:THR:HG22	1.74	0.51
1:B:402:PRO:HB3	1:B:429:THR:HG21	1.92	0.51
1:A:41:LEU:HB3	1:A:43:PHE:HE2	1.75	0.51
1:A:115:ARG:H	1:A:115:ARG:HD2	1.75	0.51
1:B:147:ILE:HD12	1:B:147:ILE:N	2.26	0.51
1:B:184:PRO:HG3	1:B:235:GLN:HB2	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:593:ILE:HD12	1:B:593:ILE:N	2.25	0.51
1:B:601:GLY:O	1:B:602:ALA:HB2	2.09	0.51
1:B:629:GLY:O	1:B:630:GLU:O	2.29	0.51
1:A:29:HIS:HA	1:A:49:VAL:O	2.11	0.51
1:A:124:LEU:HD23	1:A:127:MET:HE2	1.93	0.51
1:A:218:MET:HG2	1:A:240:PHE:CZ	2.45	0.51
1:A:457:ASP:OD1	1:A:457:ASP:N	2.44	0.51
1:B:116:TRP:HE3	1:B:328:ILE:HG12	1.75	0.51
1:B:156:LYS:C	1:B:156:LYS:HD2	2.35	0.51
1:B:218:MET:SD	1:B:236:LEU:HD12	2.51	0.51
1:B:285:LEU:O	1:B:288:GLN:N	2.43	0.51
1:A:81:ASP:HB2	1:B:95:SER:OG	2.11	0.51
1:B:318:ALA:HB2	1:B:430:TYR:CE1	2.46	0.51
1:B:78:LEU:HD12	1:B:342:LEU:HD12	1.93	0.51
1:B:99:ILE:HG21	1:B:104:VAL:CB	2.41	0.51
1:B:254:TRP:CA	1:B:259:ARG:HB3	2.41	0.51
1:B:498:LEU:H	1:B:501:GLY:CA	2.20	0.51
1:A:407:GLN:HG3	1:A:420:LEU:CD2	2.40	0.51
1:B:61:VAL:HB	1:B:320:VAL:HG13	1.93	0.51
1:B:112:THR:HG23	1:B:197:ASP:HA	1.93	0.51
1:B:353:ALA:CA	1:B:616:VAL:CG2	2.89	0.51
1:B:407:GLN:HB3	1:B:638:VAL:CG1	2.39	0.51
1:A:16:LEU:HD13	1:A:359:GLU:HB2	1.93	0.51
1:A:123:LEU:O	1:A:127:MET:HG3	2.10	0.51
1:A:137:THR:CG2	1:A:171:GLN:HA	2.41	0.51
1:B:251:ALA:O	1:B:255:ILE:HG12	2.11	0.51
1:A:399:PHE:HB3	1:A:405:PHE:CD1	2.46	0.50
1:A:640:ARG:HA	1:A:645:LEU:HA	1.93	0.50
1:B:185:ASP:HB3	1:B:239:LYS:HG3	1.93	0.50
1:B:383:LYS:NZ	1:B:392:ASP:HB3	2.26	0.50
1:B:490:VAL:HG23	1:B:587:VAL:C	2.36	0.50
1:B:599:GLN:HG3	1:B:599:GLN:O	2.10	0.50
1:A:30:ASN:HD22	1:A:596:THR:HB	1.77	0.50
1:A:613:LEU:HD12	1:A:613:LEU:H	1.75	0.50
1:B:61:VAL:HG11	1:B:231:TYR:CD1	2.47	0.50
1:B:97:CYS:O	1:B:99:ILE:N	2.41	0.50
1:A:123:LEU:HD13	1:A:218:MET:CE	2.39	0.50
1:A:254:TRP:CD1	1:A:264:PHE:HB3	2.46	0.50
1:B:218:MET:HA	1:B:236:LEU:HD12	1.93	0.50
1:A:96:THR:CG2	1:A:97:CYS:N	2.75	0.50
1:A:399:PHE:CB	1:A:434:LEU:HD21	2.41	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:15:ASP:HA	1:B:610:ARG:HA	1.93	0.50
1:B:63:PHE:HA	1:B:322:ILE:O	2.12	0.50
1:B:354:ILE:HG12	1:B:361:VAL:CG2	2.42	0.50
1:B:502:ALA:O	1:B:503:VAL:HG23	2.12	0.50
1:A:26:PHE:CE2	1:A:53:PHE:HB2	2.47	0.50
1:A:61:VAL:HG13	1:A:320:VAL:CG2	2.42	0.50
1:B:453:ASN:OD1	1:B:460:VAL:HG22	2.12	0.50
1:B:549:PRO:HB2	1:B:551:PHE:CD2	2.47	0.50
1:A:154:HIS:C	1:A:447:ILE:HB	2.36	0.50
1:A:378:LEU:O	1:A:382:LYS:HB2	2.11	0.50
1:A:442:VAL:O	1:A:458:HIS:HB3	2.12	0.50
1:B:157:ILE:HD11	1:B:308:ALA:HB1	1.92	0.50
1:B:307:CYS:HB3	1:B:538:TYR:HE2	1.76	0.50
1:B:616:VAL:O	1:B:617:LYS:C	2.54	0.50
1:A:49:VAL:HG11	1:A:471:TYR:CD1	2.46	0.50
1:A:194:PRO:HD2	1:A:332:TYR:CE2	2.46	0.50
1:A:453:ASN:HB3	1:A:460:VAL:HG22	1.93	0.50
1:A:483:ASN:HB2	1:A:484:PRO:CD	2.41	0.50
1:A:496:LYS:HB3	1:A:503:VAL:CG2	2.41	0.50
1:B:378:LEU:HD11	1:B:478:ALA:HB1	1.94	0.50
1:A:118:ASP:OD2	1:A:328:ILE:HG12	2.12	0.50
1:A:598:VAL:HG13	1:A:606:TYR:HB2	1.94	0.50
1:B:5:VAL:HG11	1:B:360:LEU:CD1	2.36	0.50
1:B:45:ARG:HG2	1:B:46:VAL:N	2.27	0.50
1:B:218:MET:HG2	1:B:240:PHE:CE2	2.46	0.50
1:A:198:VAL:HG13	1:A:199:PRO:CD	2.40	0.49
1:A:315:MET:HG3	1:A:442:VAL:HG12	1.94	0.49
1:A:548:ILE:HD11	1:A:552:VAL:HG11	1.93	0.49
1:A:549:PRO:HB2	1:A:551:PHE:CE2	2.47	0.49
1:B:208:VAL:HG23	1:B:248:VAL:HG23	1.94	0.49
1:A:350:ASP:O	1:A:354:ILE:HG12	2.12	0.49
1:B:576:PRO:O	1:B:578:LEU:HD23	2.12	0.49
1:B:634:TYR:N	1:B:634:TYR:CD1	2.80	0.49
1:A:89:VAL:HG13	1:A:334:PHE:CZ	2.48	0.49
1:A:453:ASN:O	1:A:459:LEU:HB2	2.11	0.49
1:B:6:THR:O	1:B:7:LYS:HB2	2.12	0.49
1:B:27:VAL:HG12	1:B:599:GLN:O	2.12	0.49
1:A:516:LEU:HD12	1:A:521:TYR:HB2	1.94	0.49
1:B:7:LYS:HD2	1:B:10:GLN:CD	2.37	0.49
1:B:22:ASP:O	1:B:23:ARG:HB2	2.13	0.49
1:B:161:VAL:HG22	1:B:226:ASN:OD1	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:353:ALA:CA	1:B:616:VAL:HG21	2.42	0.49
1:A:73:ILE:HG13	1:A:329:ARG:HH21	1.76	0.49
1:A:207:THR:CG2	1:A:210:GLU:H	2.26	0.49
1:A:299:ILE:HG22	1:A:428:VAL:CG1	2.42	0.49
1:A:584:PHE:C	1:A:585:ARG:HG2	2.36	0.49
1:B:527:VAL:HG13	1:B:531:THR:CG2	2.42	0.49
1:A:186:TRP:HE3	1:A:186:TRP:O	1.95	0.49
1:A:531:THR:O	1:A:531:THR:CG2	2.60	0.49
1:B:6:THR:O	1:B:7:LYS:CB	2.60	0.49
1:B:378:LEU:HD12	1:B:413:CYS:HA	1.95	0.49
1:A:73:ILE:HB	1:A:78:LEU:CD2	2.39	0.49
1:B:22:ASP:HB2	1:B:601:GLY:H	1.78	0.49
1:B:491:ILE:O	1:B:584:PHE:HZ	1.95	0.49
1:A:41:LEU:HB3	1:A:43:PHE:CE2	2.47	0.49
1:A:210:GLU:OE2	1:A:278:LYS:HD2	2.13	0.49
1:A:305:PRO:HB3	1:A:467:ALA:CB	2.43	0.49
1:A:27:VAL:HA	1:A:52:LYS:HA	1.95	0.49
1:A:37:LYS:O	1:A:587:VAL:HA	2.12	0.49
1:A:64:ASN:HD22	1:A:321:ASN:HD21	1.60	0.49
1:A:382:LYS:HB3	1:A:394:TYR:CE2	2.48	0.49
1:B:37:LYS:HB3	1:B:588:GLU:HB2	1.95	0.49
1:B:150:TYR:CD2	1:B:308:ALA:HB2	2.47	0.49
1:A:54:THR:HG22	1:A:302:LYS:O	2.13	0.49
1:A:354:ILE:HD12	1:A:361:VAL:CG2	2.43	0.49
1:A:571:HIS:CE1	1:A:574:GLU:HA	2.48	0.49
1:B:99:ILE:HG21	1:B:104:VAL:N	2.28	0.49
1:B:208:VAL:HG13	1:B:209:THR:N	2.28	0.49
1:A:175:PRO:HB2	1:A:221:TRP:HA	1.95	0.48
1:B:142:TYR:CE1	1:B:172:PHE:HB2	2.48	0.48
1:B:155:VAL:H	1:B:447:ILE:HB	1.76	0.48
1:B:370:VAL:HG23	1:B:468:CYS:SG	2.53	0.48
1:A:332:TYR:CB	1:A:335:LEU:HD13	2.37	0.48
1:A:378:LEU:HG	1:A:478:ALA:CB	2.43	0.48
1:A:407:GLN:HB2	1:A:638:VAL:HG11	1.95	0.48
1:B:123:LEU:HD13	1:B:218:MET:HE2	1.94	0.48
1:B:126:ASN:HA	1:B:129:ARG:CG	2.43	0.48
1:A:38:PHE:HB3	1:A:587:VAL:HG22	1.94	0.48
1:A:349:GLU:HA	1:A:618:PRO:HG3	1.96	0.48
1:A:399:PHE:O	1:A:434:LEU:HD21	2.13	0.48
1:B:505:ASP:HB2	1:B:508:GLU:HG2	1.96	0.48
1:A:19:ILE:H	1:A:19:ILE:HG13	1.39	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:213:PHE:O	1:A:217:MET:HG2	2.13	0.48
1:A:534:LEU:HG	1:A:535:THR:N	2.28	0.48
1:A:613:LEU:HD12	1:A:613:LEU:N	2.29	0.48
1:B:429:THR:O	1:B:431:PRO:HD3	2.13	0.48
1:B:123:LEU:HD22	1:B:218:MET:HE1	1.95	0.48
1:B:195:SER:C	1:B:196:ILE:HG13	2.39	0.48
1:B:342:LEU:HA	1:B:626:ILE:HG12	1.94	0.48
1:B:568:ARG:HB3	1:B:568:ARG:NH1	2.29	0.48
1:A:395:GLU:HG3	1:A:396:ASP:H	1.70	0.48
1:A:528:TYR:CE2	1:A:557:VAL:HG23	2.49	0.48
1:B:13:SER:OG	1:B:610:ARG:NH1	2.46	0.48
1:B:163:ILE:HG23	1:B:226:ASN:ND2	2.29	0.48
1:B:284:ARG:HD3	1:B:340:ALA:CB	2.40	0.48
1:B:309:GLU:CD	1:B:309:GLU:H	2.21	0.48
1:A:29:HIS:ND1	1:A:50:SER:HB3	2.28	0.48
1:A:30:ASN:HA	1:A:595:ASP:O	2.14	0.48
1:A:47:TYR:CD2	1:A:47:TYR:N	2.81	0.48
1:A:27:VAL:HG13	1:A:29:HIS:NE2	2.29	0.48
1:A:27:VAL:HG13	1:A:29:HIS:HE2	1.78	0.48
1:B:129:ARG:O	1:B:133:LEU:HD22	2.13	0.48
1:B:142:TYR:O	1:B:170:THR:HA	2.13	0.48
1:B:149:MET:CE	1:B:156:LYS:HD3	2.43	0.48
1:B:384:THR:HG23	1:B:386:PHE:H	1.78	0.48
1:B:479:PHE:C	1:B:481:THR:H	2.19	0.48
1:B:517:ARG:HE	1:B:539:ALA:HB3	1.78	0.48
1:B:639:ARG:HG3	1:B:640:ARG:N	2.28	0.48
1:A:24:GLY:C	1:A:25:THR:HG22	2.38	0.48
1:A:82:GLY:CA	1:B:95:SER:HB3	2.43	0.48
1:A:271:VAL:HA	1:A:274:SER:OG	2.14	0.48
1:A:283:ASN:HA	1:A:336:LEU:HD22	1.96	0.48
1:B:57:GLY:O	1:B:223:ARG:NH2	2.47	0.48
1:B:301:MET:SD	1:B:366:MET:HA	2.53	0.48
1:B:634:TYR:N	1:B:634:TYR:HD1	2.12	0.48
1:A:218:MET:HE3	1:A:218:MET:HA	1.95	0.47
1:B:162:THR:HG22	1:B:162:THR:O	2.14	0.47
1:B:204:ARG:HB2	1:B:205:PRO:HD3	1.95	0.47
1:B:291:THR:HG21	1:B:423:MET:HG3	1.97	0.47
1:B:493:ALA:CB	1:B:497:TYR:OH	2.62	0.47
1:B:390:ILE:HD12	1:B:390:ILE:N	2.06	0.47
1:B:631:LEU:HD13	1:B:631:LEU:O	2.15	0.47
1:B:1:MET:HG3	1:B:364:TYR:CD1	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1:MET:O	1:B:5:VAL:HG12	2.14	0.47
1:B:64:ASN:HA	1:B:324:LYS:NZ	2.30	0.47
1:B:220:LYS:HG3	1:B:234:PRO:HA	1.95	0.47
1:B:284:ARG:NH2	1:B:338:GLY:O	2.48	0.47
1:B:630:GLU:CG	1:B:631:LEU:N	2.71	0.47
1:A:296:LEU:O	1:A:299:ILE:HD13	2.15	0.47
1:A:376:LEU:C	1:A:464:VAL:HG21	2.39	0.47
1:B:154:HIS:O	1:B:155:VAL:O	2.33	0.47
1:B:444:ILE:HG22	1:B:445:THR:H	1.79	0.47
1:A:99:ILE:HG22	1:A:104:VAL:HG13	1.96	0.47
1:A:1:MET:HG3	1:A:364:TYR:CE1	2.50	0.47
1:A:73:ILE:HG22	1:A:74:ALA:N	2.29	0.47
1:A:254:TRP:CA	1:A:259:ARG:HB3	2.45	0.47
1:A:354:ILE:HD12	1:A:361:VAL:HG23	1.96	0.47
1:A:371:ALA:HB3	1:A:410:LEU:CD2	2.45	0.47
1:A:399:PHE:HB3	1:A:434:LEU:HD21	1.96	0.47
1:A:400:LEU:CD1	1:B:281:ASN:HB3	2.45	0.47
1:B:38:PHE:HB3	1:B:587:VAL:HG22	1.97	0.47
1:B:56:VAL:HA	1:B:227:LEU:HA	1.96	0.47
1:B:114:TRP:H	1:B:114:TRP:CD1	2.33	0.47
1:B:353:ALA:HA	1:B:616:VAL:HG23	1.93	0.47
1:B:451:GLY:HA2	1:B:532:HIS:C	2.39	0.47
1:B:506:LYS:HG3	1:B:558:MET:O	2.15	0.47
1:B:527:VAL:HG11	1:B:531:THR:HG23	1.96	0.47
1:A:96:THR:HG23	1:A:97:CYS:N	2.30	0.47
1:A:236:LEU:HD22	1:A:236:LEU:N	2.23	0.47
1:A:315:MET:SD	1:A:442:VAL:HG12	2.55	0.47
1:B:61:VAL:HA	1:B:320:VAL:HG13	1.96	0.47
1:B:129:ARG:NH2	1:B:231:TYR:O	2.44	0.47
1:B:284:ARG:CZ	1:B:336:LEU:O	2.63	0.47
1:B:331:ARG:HB3	1:B:332:TYR:CE1	2.49	0.47
1:B:531:THR:O	1:B:534:LEU:HB2	2.15	0.47
1:B:372:LEU:HD12	1:B:409:ALA:HB3	1.96	0.47
1:A:61:VAL:HG21	1:A:231:TYR:CG	2.50	0.47
1:A:211:VAL:O	1:A:215:LEU:HD23	2.14	0.47
1:B:14:SER:HB2	1:B:611:ILE:HG23	1.96	0.47
1:B:297:ALA:O	1:B:365:ALA:HB1	2.15	0.47
1:B:302:LYS:HE2	1:B:313:TRP:CZ2	2.49	0.47
1:B:407:GLN:C	1:B:638:VAL:HG11	2.40	0.47
1:B:650:LEU:HD23	1:B:650:LEU:N	2.30	0.47
1:A:194:PRO:HB3	1:A:331:ARG:HD2	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:283:ASN:HB3	1:A:329:ARG:HD3	1.97	0.46
1:A:353:ALA:HA	1:A:616:VAL:HG21	1.96	0.46
1:A:396:ASP:O	1:A:397:GLY:O	2.34	0.46
1:A:451:GLY:C	1:A:452:TYR:HD1	2.23	0.46
1:B:116:TRP:CE3	1:B:328:ILE:HG12	2.50	0.46
1:B:378:LEU:CD1	1:B:413:CYS:HA	2.45	0.46
1:A:414:THR:HB	1:A:416:GLN:NE2	2.30	0.46
1:A:446:VAL:O	1:A:447:ILE:HD13	2.14	0.46
1:B:253:GLU:HB3	1:B:259:ARG:HB2	1.95	0.46
1:B:66:GLY:O	1:B:117:TYR:HE1	1.99	0.46
1:B:174:TRP:CD2	1:B:216:MET:HE2	2.50	0.46
1:B:493:ALA:N	1:B:584:PHE:CZ	2.83	0.46
1:B:519:ALA:O	1:B:572:PHE:O	2.32	0.46
1:A:115:ARG:HD2	1:A:115:ARG:N	2.31	0.46
1:A:210:GLU:O	1:A:214:VAL:HG23	2.15	0.46
1:A:342:LEU:N	1:A:342:LEU:CD2	2.79	0.46
1:B:7:LYS:HB2	1:B:10:GLN:HG2	1.97	0.46
1:B:344:GLN:HG3	1:B:635:TRP:CZ3	2.51	0.46
1:B:381:VAL:O	1:B:384:THR:N	2.48	0.46
1:B:420:LEU:HD13	1:B:637:SER:CA	2.39	0.46
1:A:316:HIS:CD2	1:A:433:LEU:HD11	2.50	0.46
1:A:318:ALA:HB2	1:A:430:TYR:CZ	2.50	0.46
1:A:442:VAL:C	1:A:458:HIS:CB	2.88	0.46
1:A:512:LEU:C	1:A:512:LEU:HD23	2.41	0.46
1:B:59:PRO:CD	1:B:223:ARG:HH22	2.28	0.46
1:B:331:ARG:HB3	1:B:332:TYR:CD1	2.51	0.46
1:B:437:ASP:O	1:B:437:ASP:CG	2.59	0.46
1:B:497:TYR:CG	1:B:581:PRO:HG3	2.50	0.46
1:B:64:ASN:O	1:B:323:PRO:HA	2.15	0.46
1:B:380:ARG:HH11	1:B:380:ARG:CG	2.28	0.46
1:B:424:SER:O	1:B:425:ASP:HB2	2.16	0.46
1:B:446:VAL:C	1:B:447:ILE:HG12	2.41	0.46
1:B:531:THR:HG21	1:B:534:LEU:O	2.16	0.46
1:B:611:ILE:HD13	1:B:611:ILE:C	2.41	0.46
1:A:300:MET:HG2	1:A:301:MET:CE	2.46	0.46
1:B:1:MET:HG3	1:B:364:TYR:CE1	2.50	0.46
1:B:117:TYR:O	1:B:118:ASP:C	2.57	0.46
1:B:363:THR:O	1:B:367:GLN:HG3	2.16	0.46
1:B:570:ARG:NH1	1:B:570:ARG:CG	2.79	0.46
1:B:611:ILE:O	1:B:611:ILE:HD13	2.15	0.46
1:A:18:SER:HB3	1:A:359:GLU:CG	2.45	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:204:ARG:HD2	1:A:243:ARG:CD	2.46	0.46
1:A:297:ALA:HA	1:A:300:MET:HE2	1.97	0.46
1:B:32:VAL:HG12	1:B:593:ILE:HG13	1.98	0.46
1:B:74:ALA:HB2	1:B:337:SER:O	2.15	0.46
1:B:514:TRP:HD1	1:B:546:THR:H	1.63	0.46
1:B:650:LEU:H	1:B:650:LEU:CD2	2.29	0.46
1:A:77:TYR:C	1:A:78:LEU:HD22	2.40	0.46
1:A:133:LEU:HD12	1:A:142:TYR:HE2	1.81	0.46
1:A:445:THR:HG22	1:A:445:THR:O	2.16	0.46
1:B:130:ALA:C	1:B:272:MET:HE1	2.40	0.46
1:B:154:HIS:CD2	1:B:448:GLU:HG3	2.46	0.46
1:B:603:HIS:O	1:B:604:ALA:O	2.33	0.46
1:A:386:PHE:C	1:A:388:THR:H	2.24	0.45
1:A:403:GLU:CD	1:A:403:GLU:H	2.23	0.45
1:B:76:LYS:HD2	1:B:77:TYR:CE2	2.51	0.45
1:B:118:ASP:O	1:B:119:ASN:HB2	2.17	0.45
1:B:319:LEU:O	1:B:319:LEU:HD23	2.17	0.45
1:A:175:PRO:O	1:A:222:HIS:HB2	2.15	0.45
1:A:198:VAL:O	1:A:200:TYR:CD1	2.68	0.45
1:A:218:MET:HE2	1:A:236:LEU:HG	1.97	0.45
1:A:613:LEU:HA	1:A:616:VAL:HG12	1.98	0.45
1:A:625:VAL:HG13	1:A:625:VAL:O	2.14	0.45
1:B:99:ILE:HG21	1:B:104:VAL:HB	1.99	0.45
1:B:563:THR:O	1:B:564:ALA:HB2	2.16	0.45
1:A:137:THR:HG23	1:A:171:GLN:HA	1.98	0.45
1:B:16:LEU:HD13	1:B:359:GLU:CB	2.47	0.45
1:B:283:ASN:N	1:B:283:ASN:ND2	2.64	0.45
1:B:581:PRO:O	1:B:582:ALA:HB3	2.17	0.45
1:A:562:VAL:O	1:A:562:VAL:HG13	2.17	0.45
1:B:17:PHE:O	1:B:363:THR:HG22	2.15	0.45
1:B:154:HIS:HA	1:B:447:ILE:HB	1.99	0.45
1:B:220:LYS:CG	1:B:234:PRO:HA	2.46	0.45
1:B:294:GLN:NE2	1:B:419:PRO:HB2	2.29	0.45
1:B:299:ILE:C	1:B:299:ILE:HD12	2.41	0.45
1:B:454:ILE:HD12	1:B:454:ILE:H	1.80	0.45
1:A:312:ALA:HA	1:A:315:MET:HG3	1.97	0.45
1:A:376:LEU:HB3	1:A:464:VAL:HG21	1.98	0.45
1:B:476:VAL:HB	1:B:485:TYR:O	2.16	0.45
1:A:198:VAL:O	1:A:200:TYR:HD1	1.99	0.45
1:A:300:MET:HB3	1:A:365:ALA:HB1	1.99	0.45
1:B:59:PRO:HD2	1:B:223:ARG:NH2	2.32	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:638:VAL:HG13	1:B:638:VAL:O	2.17	0.45
1:A:207:THR:HG23	1:A:210:GLU:H	1.82	0.45
1:A:218:MET:HE2	1:A:236:LEU:CG	2.47	0.45
1:A:125:MET:HE1	1:A:295:LEU:HB3	1.98	0.45
1:A:157:ILE:HG22	1:A:443:PRO:O	2.17	0.45
1:A:583:PHE:CD1	1:A:583:PHE:N	2.84	0.45
1:B:19:ILE:HG23	1:B:362:PHE:CD1	2.52	0.45
1:B:42:VAL:O	1:B:42:VAL:HG13	2.17	0.45
1:B:499:ARG:H	1:B:499:ARG:HG3	1.42	0.45
1:A:117:TYR:CD2	1:A:326:GLY:HA3	2.52	0.45
1:B:29:HIS:HE1	1:B:48:GLY:HA3	1.81	0.45
1:B:99:ILE:CB	1:B:104:VAL:HB	2.46	0.45
1:B:163:ILE:HG23	1:B:226:ASN:HD21	1.82	0.45
1:B:418:ALA:O	1:B:637:SER:HB2	2.17	0.45
1:A:123:LEU:HB2	1:A:218:MET:CE	2.42	0.45
1:A:185:ASP:HB3	1:A:239:LYS:CG	2.47	0.45
1:A:204:ARG:HA	1:A:244:HIS:HA	1.99	0.45
1:A:435:GLU:HB2	1:A:438:ALA:HB2	1.99	0.45
1:A:448:GLU:HA	1:A:449:PRO:HD3	1.63	0.45
1:A:470:PRO:HG2	1:A:545:TRP:CZ3	2.52	0.45
1:A:473:ILE:HD12	1:A:476:VAL:HG21	1.98	0.45
1:B:61:VAL:CB	1:B:320:VAL:HG13	2.47	0.45
1:B:455:VAL:O	1:B:458:HIS:HB2	2.17	0.45
1:B:504:TYR:C	1:B:560:VAL:HG12	2.42	0.45
1:A:73:ILE:HG13	1:A:329:ARG:NH2	2.32	0.44
1:A:78:LEU:HD13	1:A:84:LEU:HA	1.98	0.44
1:A:200:TYR:HB3	1:A:330:GLY:C	2.42	0.44
1:B:109:TYR:C	1:B:109:TYR:CD2	2.95	0.44
1:B:470:PRO:HG2	1:B:471:TYR:H	1.82	0.44
1:A:81:ASP:N	1:A:81:ASP:OD1	2.49	0.44
1:A:204:ARG:HD2	1:A:243:ARG:CG	2.46	0.44
1:A:447:ILE:C	1:A:448:GLU:HG2	2.42	0.44
1:B:223:ARG:HH11	1:B:231:TYR:HA	1.79	0.44
1:B:344:GLN:HG3	1:B:635:TRP:CH2	2.52	0.44
1:A:285:LEU:O	1:A:286:TYR:C	2.60	0.44
1:A:320:VAL:CG1	1:A:428:VAL:HG22	2.43	0.44
1:A:631:LEU:O	1:A:634:TYR:N	2.47	0.44
1:B:123:LEU:HB2	1:B:236:LEU:HD13	1.98	0.44
1:A:137:THR:HA	1:A:142:TYR:HB2	2.00	0.44
1:A:521:TYR:CA	1:A:568:ARG:HG3	2.47	0.44
1:B:23:ARG:HB2	1:B:23:ARG:NH1	2.32	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:30:ASN:OD1	1:B:596:THR:HB	2.17	0.44
1:B:66:GLY:O	1:B:67:SER:C	2.61	0.44
1:B:130:ALA:HB2	1:B:221:TRP:CH2	2.52	0.44
1:B:442:VAL:O	1:B:444:ILE:HG12	2.17	0.44
1:A:303:PRO:HB2	1:A:468:CYS:HB3	1.99	0.44
1:A:355:MET:CE	1:A:355:MET:HA	2.46	0.44
1:B:115:ARG:HH11	1:B:115:ARG:CG	2.31	0.44
1:A:207:THR:HG22	1:A:210:GLU:HB2	2.00	0.44
1:A:370:VAL:HA	1:A:468:CYS:SG	2.57	0.44
1:A:407:GLN:H	1:A:407:GLN:HG2	1.58	0.44
1:B:115:ARG:HA	1:B:115:ARG:HD2	1.61	0.44
1:B:127:MET:HB2	1:B:276:LEU:HD13	2.00	0.44
1:B:435:GLU:O	1:B:438:ALA:N	2.48	0.44
1:A:114:TRP:H	1:A:114:TRP:CD1	2.35	0.44
1:A:204:ARG:N	1:A:205:PRO:CD	2.81	0.44
1:B:118:ASP:HB2	1:B:328:ILE:HB	2.00	0.44
1:B:126:ASN:O	1:B:129:ARG:HG2	2.18	0.44
1:B:479:PHE:C	1:B:481:THR:N	2.76	0.44
1:B:547:HIS:O	1:B:549:PRO:HD3	2.18	0.44
1:A:125:MET:HE1	1:A:295:LEU:HB2	2.00	0.44
1:A:283:ASN:N	1:A:283:ASN:HD22	2.15	0.44
1:A:328:ILE:HA	1:A:331:ARG:HG3	1.99	0.44
1:A:332:TYR:HB2	1:A:335:LEU:CD1	2.39	0.44
1:A:481:THR:HG22	1:A:482:ALA:N	2.33	0.44
1:B:91:ASN:HD22	1:B:94:ARG:HD2	1.82	0.44
1:B:129:ARG:HG3	1:B:221:TRP:CH2	2.52	0.44
1:A:131:TYR:CE1	1:A:269:SER:HB2	2.53	0.44
1:A:399:PHE:HB3	1:A:405:PHE:HD1	1.82	0.44
1:A:459:LEU:C	1:A:459:LEU:CD1	2.91	0.44
1:B:252:ASP:O	1:B:255:ILE:HB	2.18	0.44
1:B:380:ARG:HA	1:B:436:PHE:HD2	1.78	0.44
1:B:306:ASN:HD21	1:B:547:HIS:HE1	1.66	0.43
1:B:470:PRO:O	1:B:472:MET:N	2.50	0.43
1:B:525:PHE:CD1	1:B:525:PHE:C	2.96	0.43
1:A:395:GLU:HB3	1:A:436:PHE:HE1	1.80	0.43
1:B:2:LEU:HD13	1:B:2:LEU:C	2.43	0.43
1:B:61:VAL:CA	1:B:320:VAL:HG13	2.48	0.43
1:B:99:ILE:HA	1:B:100:PRO:HD3	1.77	0.43
1:B:223:ARG:HG3	1:B:230:ASP:O	2.18	0.43
1:B:474:PHE:C	1:B:476:VAL:N	2.54	0.43
1:A:442:VAL:HA	1:A:443:PRO:HD3	1.85	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:473:ILE:C	1:A:475:PRO:HD2	2.43	0.43
1:B:58:ASN:HA	1:B:223:ARG:NH2	2.34	0.43
1:B:111:ILE:O	1:B:197:ASP:HB2	2.18	0.43
1:B:154:HIS:O	1:B:155:VAL:C	2.61	0.43
1:A:194:PRO:HD2	1:A:332:TYR:CZ	2.52	0.43
1:A:207:THR:HG22	1:A:210:GLU:HG3	1.99	0.43
1:B:43:PHE:HB3	1:B:573:VAL:HG13	2.00	0.43
1:B:312:ALA:O	1:B:315:MET:HG2	2.19	0.43
1:B:495:ASN:HB2	1:B:496:LYS:HE3	1.99	0.43
1:A:116:TRP:H	1:B:98:GLY:HA3	1.84	0.43
1:A:207:THR:HG22	1:A:210:GLU:CB	2.49	0.43
1:A:354:ILE:HD13	1:A:360:LEU:HD23	2.00	0.43
1:A:539:ALA:HB2	1:A:546:THR:HA	2.00	0.43
1:A:85:ALA:O	1:A:89:VAL:HG23	2.19	0.43
1:A:283:ASN:HA	1:A:336:LEU:HD13	2.00	0.43
1:A:320:VAL:HA	1:A:427:TYR:O	2.19	0.43
1:A:616:VAL:HG22	1:A:616:VAL:O	2.19	0.43
1:B:75:LYS:CA	1:B:341:ALA:HB2	2.47	0.43
1:A:277:ARG:CB	1:A:355:MET:HE2	2.41	0.43
1:A:380:ARG:O	1:A:383:LYS:HE3	2.19	0.43
1:A:418:ALA:C	1:A:637:SER:HB2	2.43	0.43
1:A:434:LEU:HD22	1:A:434:LEU:HA	1.84	0.43
1:B:71:GLU:OE1	1:B:71:GLU:N	2.51	0.43
1:B:211:VAL:HG12	1:B:215:LEU:HD11	1.99	0.43
1:A:89:VAL:O	1:A:93:LEU:HD13	2.19	0.43
1:A:137:THR:CG2	1:A:172:PHE:H	2.27	0.43
1:A:204:ARG:HB2	1:A:205:PRO:HD3	2.01	0.43
1:B:14:SER:CB	1:B:611:ILE:HG23	2.48	0.43
1:B:112:THR:O	1:B:114:TRP:N	2.51	0.43
1:B:188:GLN:HG2	1:B:247:THR:OG1	2.18	0.43
1:B:447:ILE:O	1:B:448:GLU:C	2.61	0.43
1:B:562:VAL:HG22	1:B:562:VAL:O	2.18	0.43
1:A:82:GLY:N	1:B:95:SER:HB3	2.34	0.43
1:A:157:ILE:HG21	1:A:444:ILE:HG22	2.01	0.43
1:A:477:ALA:C	1:A:479:PHE:H	2.27	0.43
1:A:532:HIS:O	1:A:532:HIS:CG	2.72	0.43
1:B:112:THR:HG22	1:B:197:ASP:HA	2.01	0.43
1:B:200:TYR:CD1	1:B:241:ALA:HB3	2.54	0.43
1:B:210:GLU:HG2	1:B:275:ALA:HB2	2.00	0.43
1:B:474:PHE:CD2	1:B:593:ILE:HD11	2.54	0.43
1:A:65:GLU:CD	1:B:191:GLU:HA	2.44	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:154:HIS:CA	1:A:448:GLU:HG3	2.27	0.43
1:A:386:PHE:C	1:A:388:THR:N	2.77	0.43
1:B:186:TRP:HA	1:B:240:PHE:O	2.19	0.43
1:B:510:TRP:NE1	1:B:546:THR:O	2.47	0.43
1:A:134:GLN:O	1:A:138:GLU:HB2	2.19	0.42
1:B:25:THR:HA	1:B:53:PHE:O	2.19	0.42
1:B:372:LEU:HD13	1:B:406:VAL:HG13	2.01	0.42
1:B:491:ILE:O	1:B:586:SER:HA	2.18	0.42
1:B:507:LEU:N	1:B:507:LEU:HD13	2.34	0.42
1:A:155:VAL:CG2	1:A:156:LYS:N	2.81	0.42
1:A:443:PRO:N	1:A:458:HIS:HB3	2.34	0.42
1:A:617:LYS:HA	1:A:618:PRO:HD3	1.87	0.42
1:B:150:TYR:CZ	1:B:152:ASP:HB3	2.54	0.42
1:A:74:ALA:O	1:A:78:LEU:HD23	2.19	0.42
1:A:93:LEU:HD21	1:A:193:PHE:HB3	2.01	0.42
1:A:344:GLN:OE1	1:A:626:ILE:HD11	2.19	0.42
1:A:396:ASP:OD2	1:B:277:ARG:CD	2.60	0.42
1:B:112:THR:HB	1:B:331:ARG:CZ	2.49	0.42
1:B:202:ASP:OD1	1:B:204:ARG:HD3	2.18	0.42
1:B:229:ILE:HB	1:B:296:LEU:HD11	2.01	0.42
1:B:285:LEU:O	1:B:286:TYR:C	2.63	0.42
1:A:517:ARG:HD3	1:A:517:ARG:O	2.20	0.42
1:A:575:LEU:HB3	1:A:576:PRO:HD2	2.00	0.42
1:B:37:LYS:HD3	1:B:588:GLU:CD	2.44	0.42
1:B:119:ASN:N	1:B:326:GLY:O	2.51	0.42
1:B:156:LYS:CG	1:B:445:THR:HG21	2.46	0.42
1:B:372:LEU:HD12	1:B:409:ALA:CB	2.49	0.42
1:B:381:VAL:O	1:B:382:LYS:C	2.62	0.42
1:A:150:TYR:HB2	1:A:311:TYR:CE1	2.54	0.42
1:B:43:PHE:CD1	1:B:43:PHE:N	2.87	0.42
1:B:342:LEU:H	1:B:624:GLN:HG2	1.85	0.42
1:B:560:VAL:HG22	1:B:561:PHE:N	2.34	0.42
1:B:64:ASN:HA	1:B:324:LYS:HG3	2.02	0.42
1:B:93:LEU:H	1:B:93:LEU:HG	1.53	0.42
1:B:403:GLU:HB2	1:B:634:TYR:HE2	1.83	0.42
1:A:159:LEU:HA	1:A:160:PRO:HD3	1.88	0.42
1:A:501:GLY:C	1:A:561:PHE:HE2	2.28	0.42
1:A:578:LEU:C	1:A:580:SER:H	2.27	0.42
1:B:16:LEU:HD22	1:B:360:LEU:HA	2.01	0.42
1:B:19:ILE:HG12	1:B:132:HIS:HD2	1.85	0.42
1:B:305:PRO:HB2	1:B:311:TYR:CD1	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:504:TYR:O	1:B:560:VAL:HG12	2.20	0.42
1:A:81:ASP:O	1:A:82:GLY:C	2.61	0.42
1:A:208:VAL:HG13	1:A:209:THR:N	2.33	0.42
1:A:498:LEU:HD11	1:A:561:PHE:CZ	2.55	0.42
1:A:610:ARG:HD3	1:A:610:ARG:HA	1.89	0.42
1:B:16:LEU:HD12	1:B:16:LEU:C	2.45	0.42
1:B:32:VAL:HG12	1:B:593:ILE:HA	2.01	0.42
1:B:33:ARG:NH1	1:B:594:TYR:HE1	2.18	0.42
1:B:442:VAL:HB	1:B:459:LEU:HB3	2.01	0.42
1:A:449:PRO:C	1:A:451:GLY:N	2.76	0.42
1:B:504:TYR:HB2	1:B:560:VAL:CG1	2.50	0.42
1:B:526:LYS:HB3	1:B:561:PHE:CD2	2.54	0.42
1:A:33:ARG:HG2	1:A:34:THR:N	2.35	0.42
1:A:91:ASN:HA	1:A:94:ARG:NH1	2.35	0.42
1:A:128:LEU:O	1:A:131:TYR:HB3	2.20	0.42
1:A:178:ARG:NH1	1:A:254:TRP:O	2.53	0.42
1:A:435:GLU:OE2	1:B:278:LYS:HE3	2.19	0.42
1:A:470:PRO:O	1:A:474:PHE:HA	2.20	0.42
1:A:554:ASP:O	1:A:557:VAL:HG12	2.20	0.42
1:B:562:VAL:O	1:B:562:VAL:HG13	2.20	0.42
1:B:641:THR:HB	1:B:643:GLN:OE1	2.20	0.42
1:A:63:PHE:HA	1:A:322:ILE:O	2.20	0.41
1:A:121:VAL:O	1:A:125:MET:HB2	2.20	0.41
1:A:442:VAL:O	1:A:444:ILE:HG23	2.20	0.41
1:B:188:GLN:O	1:B:188:GLN:HG3	2.19	0.41
1:B:474:PHE:HB3	1:B:475:PRO:HD3	2.01	0.41
1:B:643:GLN:HE21	1:B:643:GLN:HB2	1.69	0.41
1:A:61:VAL:HG13	1:A:320:VAL:HG23	2.02	0.41
1:A:343:ILE:CG2	1:A:348:LEU:HG	2.50	0.41
1:A:354:ILE:CD1	1:A:360:LEU:HD23	2.51	0.41
1:A:413:CYS:SG	1:A:478:ALA:HB2	2.61	0.41
1:B:12:LYS:HB3	1:B:12:LYS:HE2	1.91	0.41
1:B:200:TYR:HB3	1:B:330:GLY:C	2.45	0.41
1:B:218:MET:HG2	1:B:240:PHE:CZ	2.55	0.41
1:B:403:GLU:HG3	1:B:427:TYR:HD2	1.85	0.41
1:A:71:GLU:H	1:A:71:GLU:HG2	1.58	0.41
1:A:344:GLN:CD	1:A:626:ILE:HD11	2.45	0.41
1:A:514:TRP:HA	1:A:546:THR:CB	2.51	0.41
1:B:10:GLN:C	1:B:11:ASP:CG	2.88	0.41
1:B:19:ILE:HG12	1:B:132:HIS:CD2	2.56	0.41
1:B:505:ASP:O	1:B:508:GLU:N	2.52	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:526:LYS:HB3	1:B:561:PHE:HD2	1.85	0.41
1:B:543:ASP:C	1:B:545:TRP:H	2.27	0.41
1:A:132:HIS:O	1:A:136:LEU:HD23	2.20	0.41
1:A:528:TYR:CZ	1:A:557:VAL:HG23	2.55	0.41
1:A:555:GLY:C	1:A:557:VAL:H	2.28	0.41
1:A:600:ALA:O	1:A:604:ALA:HB3	2.19	0.41
1:B:394:TYR:CD1	1:B:644:GLY:HA2	2.54	0.41
1:B:488:ASN:HA	1:B:589:VAL:O	2.20	0.41
1:A:78:LEU:HD12	1:A:82:GLY:O	2.20	0.41
1:B:2:LEU:HD23	1:B:360:LEU:CD2	2.50	0.41
1:B:139:GLN:HB3	1:B:141:GLN:NE2	2.35	0.41
1:B:294:GLN:O	1:B:298:GLN:HG3	2.20	0.41
1:A:407:GLN:HB2	1:A:638:VAL:CG1	2.50	0.41
1:A:423:MET:N	1:A:633:ASN:O	2.54	0.41
1:A:550:GLU:O	1:A:553:THR:OG1	2.39	0.41
1:B:54:THR:HG23	1:B:54:THR:O	2.20	0.41
1:B:159:LEU:HA	1:B:160:PRO:HD3	1.79	0.41
1:A:372:LEU:O	1:A:376:LEU:HG	2.20	0.41
1:A:374:THR:O	1:A:378:LEU:HD23	2.21	0.41
1:B:354:ILE:HG13	1:B:360:LEU:HD23	2.02	0.41
1:A:99:ILE:CG2	1:A:104:VAL:HG13	2.51	0.41
1:A:221:TRP:HZ3	1:A:223:ARG:HA	1.86	0.41
1:A:305:PRO:HB3	1:A:467:ALA:HB1	2.03	0.41
1:A:589:VAL:O	1:A:589:VAL:HG23	2.20	0.41
1:B:4:PHE:HB2	1:B:419:PRO:HG3	2.03	0.41
1:B:172:PHE:O	1:B:266:PRO:HG2	2.20	0.41
1:B:203:VAL:HB	1:B:206:LEU:HD12	2.03	0.41
1:B:470:PRO:HG3	1:B:545:TRP:CZ2	2.56	0.41
1:B:493:ALA:HB1	1:B:494:ALA:H	1.61	0.41
1:B:531:THR:C	1:B:532:HIS:CG	2.98	0.41
1:A:12:LYS:HB3	1:A:12:LYS:HE2	1.87	0.41
1:A:100:PRO:HG2	1:A:103:ALA:CB	2.48	0.41
1:A:113:SER:O	1:A:114:TRP:C	2.63	0.41
1:A:645:LEU:H	1:A:645:LEU:CD2	2.34	0.41
1:B:1:MET:HE2	1:B:360:LEU:HG	2.02	0.41
1:B:33:ARG:HA	1:B:45:ARG:O	2.20	0.41
1:B:61:VAL:HG23	1:B:320:VAL:HG22	2.02	0.41
1:B:150:TYR:HB2	1:B:311:TYR:CZ	2.56	0.41
1:B:273:LEU:HD11	1:B:358:PRO:HG3	2.03	0.41
1:B:442:VAL:HA	1:B:443:PRO:HD3	1.95	0.41
1:B:525:PHE:CB	1:B:561:PHE:O	2.65	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:577:ARG:HH11	1:B:577:ARG:HG2	1.86	0.41
1:B:597:HIS:HD2	1:B:605:VAL:HG12	1.86	0.41
1:B:630:GLU:OE1	1:B:635:TRP:HD1	2.04	0.41
1:A:216:MET:HA	1:A:216:MET:CE	2.44	0.41
1:A:376:LEU:O	1:A:379:ARG:HB3	2.20	0.41
1:A:388:THR:O	1:A:388:THR:HG23	2.20	0.41
1:A:434:LEU:O	1:A:434:LEU:HD13	2.21	0.41
1:A:442:VAL:HG22	1:A:459:LEU:O	2.20	0.41
1:A:538:TYR:OH	1:A:540:ASP:HB2	2.21	0.41
1:A:632:LYS:HE3	1:A:632:LYS:HB3	1.86	0.41
1:B:491:ILE:HG22	1:B:584:PHE:HE1	1.86	0.41
1:A:82:GLY:O	1:A:83:GLY:C	2.63	0.40
1:B:59:PRO:HD2	1:B:223:ARG:HH12	1.86	0.40
1:B:277:ARG:O	1:B:281:ASN:HB2	2.21	0.40
1:A:142:TYR:CE2	1:A:172:PHE:HB2	2.56	0.40
1:B:64:ASN:HA	1:B:324:LYS:HZ2	1.87	0.40
1:B:89:VAL:HG13	1:B:334:PHE:CZ	2.56	0.40
1:B:473:ILE:HD13	1:B:473:ILE:N	2.36	0.40
1:B:483:ASN:CB	1:B:484:PRO:CD	2.98	0.40
1:A:426:VAL:O	1:A:426:VAL:HG23	2.20	0.40
1:A:474:PHE:N	1:A:475:PRO:CD	2.84	0.40
1:A:626:ILE:HD13	1:A:628:ALA:O	2.22	0.40
1:B:87:ASP:O	1:B:91:ASN:HB2	2.21	0.40
1:B:176:SER:OG	1:B:216:MET:HE1	2.21	0.40
1:B:315:MET:HG2	1:B:315:MET:H	1.58	0.40
1:A:7:LYS:HD3	1:A:7:LYS:HA	1.87	0.40
1:A:315:MET:CG	1:A:442:VAL:HG12	2.51	0.40
1:B:26:PHE:CZ	1:B:366:MET:HE2	2.56	0.40
1:A:64:ASN:HB3	1:B:190:SER:CB	2.52	0.40
1:A:74:ALA:HA	1:A:339:ASP:O	2.21	0.40
1:A:86:ILE:HG22	1:A:90:LEU:HD11	2.04	0.40
1:A:273:LEU:HD11	1:A:358:PRO:HG3	2.03	0.40
1:A:418:ALA:HA	1:A:419:PRO:HD2	1.87	0.40
1:B:201:LEU:HD23	1:B:201:LEU:H	1.83	0.40
1:B:526:LYS:NZ	1:B:527:VAL:H	2.13	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	649/680 (95%)	597 (92%)	36 (6%)	16 (2%)	4	28
1	B	649/680 (95%)	540 (83%)	63 (10%)	46 (7%)	1	10
All	All	1298/1360 (95%)	1137 (88%)	99 (8%)	62 (5%)	2	16

All (62) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	13	SER
1	A	397	GLY
1	A	473	ILE
1	A	531	THR
1	A	536	LYS
1	B	7	LYS
1	B	118	ASP
1	B	155	VAL
1	B	284	ARG
1	B	300	MET
1	B	451	GLY
1	B	454	ILE
1	B	471	TYR
1	B	476	VAL
1	B	477	ALA
1	B	573	VAL
1	B	602	ALA
1	B	630	GLU
1	A	83	GLY
1	A	168	GLY
1	A	556	ASP
1	A	564	ALA
1	B	12	LYS
1	B	14	SER
1	B	98	GLY

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Mol	Chain	Res	Type
1	B	113	SER
1	B	119	ASN
1	B	329	ARG
1	B	453	ASN
1	B	544	THR
1	B	603	HIS
1	A	72	GLY
1	A	114	TRP
1	B	13	SER
1	B	115	ARG
1	B	117	TYR
1	B	286	TYR
1	B	397	GLY
1	B	470	PRO
1	B	563	THR
1	B	604	ALA
1	B	608	ALA
1	B	634	TYR
1	A	112	THR
1	B	8	ASN
1	B	11	ASP
1	B	263	GLN
1	B	265	ARG
1	B	447	ILE
1	B	448	GLU
1	B	236	LEU
1	B	382	LYS
1	B	493	ALA
1	A	82	GLY
1	A	478	ALA
1	B	475	PRO
1	B	644	GLY
1	A	153	GLY
1	B	328	ILE
1	A	455	VAL
1	B	104	VAL
1	B	503	VAL

5.3.2 Protein sidechains ⓘ

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

resolution.

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	550/573 (96%)	457 (83%)	93 (17%)	2	13
1	B	550/573 (96%)	436 (79%)	114 (21%)	1	6
All	All	1100/1146 (96%)	893 (81%)	207 (19%)	1	9

All (207) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	LEU
1	A	13	SER
1	A	14	SER
1	A	19	ILE
1	A	25	THR
1	A	30	ASN
1	A	31	ARG
1	A	33	ARG
1	A	39	ASP
1	A	44	ASN
1	A	45	ARG
1	A	47	TYR
1	A	60	THR
1	A	70	LEU
1	A	71	GLU
1	A	78	LEU
1	A	87	ASP
1	A	96	THR
1	A	111	ILE
1	A	112	THR
1	A	115	ARG
1	A	119	ASN
1	A	125	MET
1	A	138	GLU
1	A	147	ILE
1	A	155	VAL
1	A	162	THR
1	A	166	THR
1	A	171	GLN
1	A	178	ARG
1	A	179	SER

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Mol	Chain	Res	Type
1	A	186	TRP
1	A	188	GLN
1	A	195	SER
1	A	198	VAL
1	A	218	MET
1	A	224	ARG
1	A	246	LEU
1	A	252	ASP
1	A	258	ASP
1	A	261	ASP
1	A	299	ILE
1	A	300	MET
1	A	309	GLU
1	A	315	MET
1	A	319	LEU
1	A	331	ARG
1	A	335	LEU
1	A	339	ASP
1	A	342	LEU
1	A	351	TRP
1	A	370	VAL
1	A	378	LEU
1	A	383	LYS
1	A	390	ILE
1	A	392	ASP
1	A	395	GLU
1	A	396	ASP
1	A	400	LEU
1	A	403	GLU
1	A	407	GLN
1	A	434	LEU
1	A	444	ILE
1	A	448	GLU
1	A	456	ASP
1	A	457	ASP
1	A	458	HIS
1	A	459	LEU
1	A	461	VAL
1	A	464	VAL
1	A	466	VAL
1	A	468	CYS
1	A	472	MET

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Mol	Chain	Res	Type
1	A	473	ILE
1	A	492	LYS
1	A	498	LEU
1	A	499	ARG
1	A	507	LEU
1	A	516	LEU
1	A	517	ARG
1	A	528	TYR
1	A	530	ASP
1	A	548	ILE
1	A	550	GLU
1	A	552	VAL
1	A	560	VAL
1	A	566	GLU
1	A	567	ARG
1	A	577	ARG
1	A	596	THR
1	A	611	ILE
1	A	642	GLN
1	A	650	LEU
1	B	5	VAL
1	B	8	ASN
1	B	11	ASP
1	B	12	LYS
1	B	15	ASP
1	B	19	ILE
1	B	49	VAL
1	B	50	SER
1	B	64	ASN
1	B	67	SER
1	B	68	SER
1	B	73	ILE
1	B	78	LEU
1	B	84	LEU
1	B	87	ASP
1	B	90	LEU
1	B	93	LEU
1	B	99	ILE
1	B	104	VAL
1	B	106	SER
1	B	112	THR
1	B	115	ARG

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Mol	Chain	Res	Type
1	B	121	VAL
1	B	133	LEU
1	B	134	GLN
1	B	161	VAL
1	B	164	ASP
1	B	170	THR
1	B	178	ARG
1	B	179	SER
1	B	181	ASP
1	B	188	GLN
1	B	196	ILE
1	B	198	VAL
1	B	201	LEU
1	B	220	LYS
1	B	224	ARG
1	B	249	GLN
1	B	255	ILE
1	B	256	GLU
1	B	263	GLN
1	B	283	ASN
1	B	285	LEU
1	B	295	LEU
1	B	301	MET
1	B	313	TRP
1	B	314	LEU
1	B	315	MET
1	B	317	ASP
1	B	319	LEU
1	B	320	VAL
1	B	328	ILE
1	B	329	ARG
1	B	346	THR
1	B	348	LEU
1	B	354	ILE
1	B	355	MET
1	B	363	THR
1	B	370	VAL
1	B	376	LEU
1	B	380	ARG
1	B	388	THR
1	B	390	ILE
1	B	393	SER

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Mol	Chain	Res	Type
1	B	395	GLU
1	B	407	GLN
1	B	416	GLN
1	B	420	LEU
1	B	429	THR
1	B	437	ASP
1	B	442	VAL
1	B	445	THR
1	B	453	ASN
1	B	454	ILE
1	B	460	VAL
1	B	464	VAL
1	B	472	MET
1	B	473	ILE
1	B	483	ASN
1	B	488	ASN
1	B	492	LYS
1	B	499	ARG
1	B	500	LYS
1	B	507	LEU
1	B	522	ASP
1	B	523	THR
1	B	526	LYS
1	B	527	VAL
1	B	530	ASP
1	B	531	THR
1	B	532	HIS
1	B	536	LYS
1	B	545	TRP
1	B	550	GLU
1	B	558	MET
1	B	561	PHE
1	B	565	ILE
1	B	578	LEU
1	B	586	SER
1	B	595	ASP
1	B	599	GLN
1	B	605	VAL
1	B	609	SER
1	B	610	ARG
1	B	611	ILE
1	B	613	LEU

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Mol	Chain	Res	Type
1	B	616	VAL
1	B	625	VAL
1	B	630	GLU
1	B	631	LEU
1	B	639	ARG
1	B	641	THR
1	B	642	GLN
1	B	643	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (44) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	8	ASN
1	A	30	ASN
1	A	51	GLN
1	A	58	ASN
1	A	119	ASN
1	A	212	ASN
1	A	235	GLN
1	A	281	ASN
1	A	283	ASN
1	A	306	ASN
1	A	316	HIS
1	A	321	ASN
1	A	373	ASN
1	A	407	GLN
1	A	416	GLN
1	A	441	GLN
1	A	483	ASN
1	A	495	ASN
1	A	571	HIS
1	A	597	HIS
1	A	624	GLN
1	B	8	ASN
1	B	10	GLN
1	B	29	HIS
1	B	44	ASN
1	B	58	ASN
1	B	91	ASN
1	B	102	ASN
1	B	134	GLN
1	B	154	HIS

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Mol	Chain	Res	Type
1	B	222	HIS
1	B	235	GLN
1	B	283	ASN
1	B	344	GLN
1	B	367	GLN
1	B	401	GLN
1	B	495	ASN
1	B	547	HIS
1	B	597	HIS
1	B	599	GLN
1	B	624	GLN
1	B	627	ASN
1	B	642	GLN
1	B	643	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](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

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

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

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	651/680 (95%)	-1.18	0 100 100	16, 54, 114, 169	0
1	B	651/680 (95%)	-1.20	0 100 100	14, 60, 127, 167	0
All	All	1302/1360 (95%)	-1.19	0 100 100	14, 57, 125, 169	0

There are no RSRZ outliers to report.

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.