



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 08:35 PM UTC

PDB ID : 2R29 / pdb_00002r29
Title : Neutralization of dengue virus by a serotype cross-reactive antibody elucidated by cryoelectron microscopy and x-ray crystallography
Authors : Lok, S.M.; Kostyuchenko, V.K.; Nybakken, G.E.; Holdaway, H.A.; Battisti, A.J.; Sukupolvi-petty, S.; Sedlak, D.; Fremont, D.H.; Chipman, P.R.; Roehrig, J.T.; Diamond, M.S.; Kuhn, R.J.; Rossmann, M.G.
Deposited on : 2007-08-24
Resolution : 3.00 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

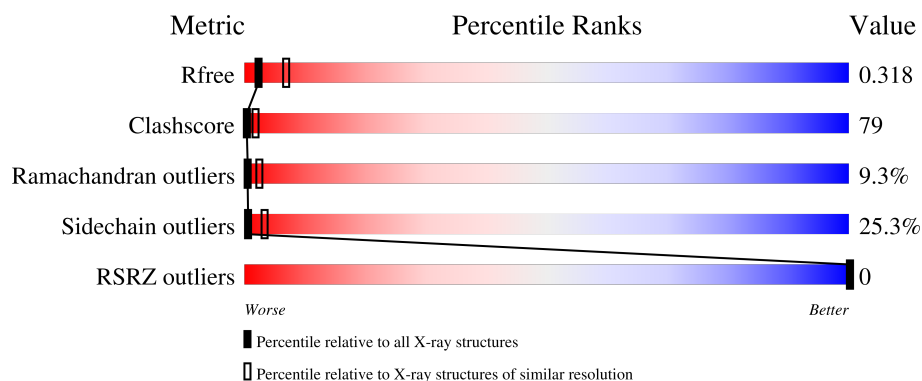
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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	97	
2	H	216	
3	L	217	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4234 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 Envelope protein E.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	97	Total	C	N	O	S	0	0	0
			768	495	126	143	4			

- Molecule 2 is a protein called Heavy chain of Fab 1A1D-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	216	Total	C	N	O	S	0	0	0
			1625	1028	266	326	5			

- Molecule 3 is a protein called Light chain of Fab 1A1D-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	L	204	Total	C	N	O	S	0	0	0
			1587	994	271	316	6			

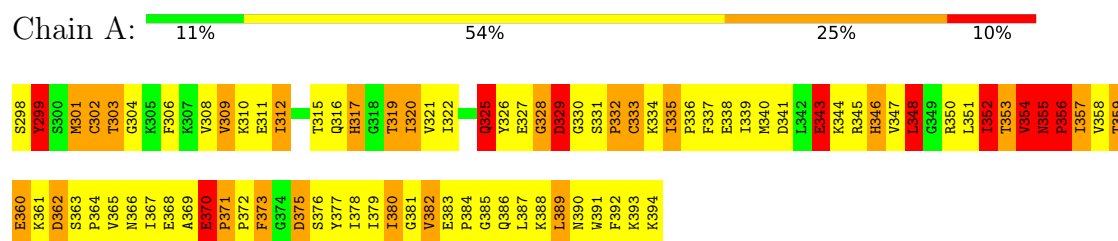
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	59	Total	O	0	0
			59	59		
4	H	97	Total	O	0	0
			97	97		
4	L	98	Total	O	0	0
			98	98		

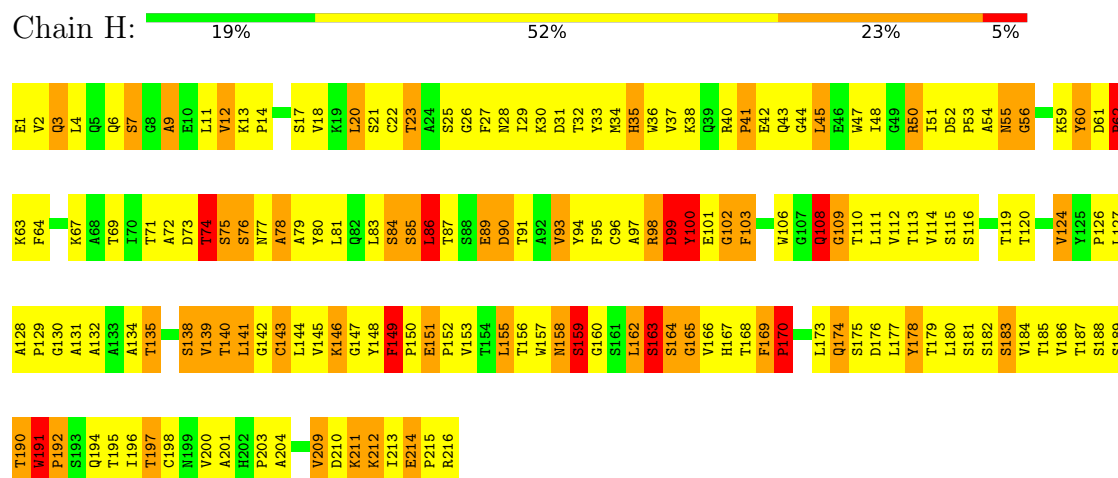
3 Residue-property plots

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.

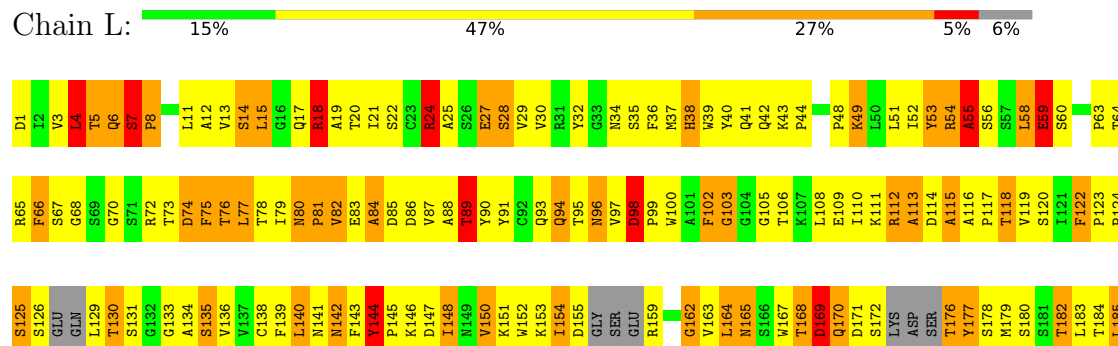
• Molecule 1: Envelope protein E



• Molecule 2: Heavy chain of Fab 1A1D-2



• Molecule 3: Light chain of Fab 1A1D-2



T186	K187	D188	E189	Y190	E191	R192	H193	M194	S195	Y196	T197	C198	E199	A200	T201	HIS	LYS	THR	SER	THR	S203	P204	I205	V206	K207	S208	F209	N210	R211	N212	E213
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4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	168.12Å 58.66Å 78.19Å 90.00° 114.41° 90.00°	Depositor
Resolution (Å)	50.00 – 3.00 50.00 – 3.00	Depositor EDS
% Data completeness (in resolution range)	(Not available) (50.00-3.00) 63.2 (50.00-3.00)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.10 (at 2.81Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.259 , 0.321 0.258 , 0.318	Depositor DCC
R_{free} test set	1125 reflections (10.28%)	wwPDB-VP
Wilson B-factor (Å ²)	45.3	Xtriage
Anisotropy	0.324	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 74.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4234	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.00% 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	0.83	1/785 (0.1%)	1.62	21/1060 (2.0%)
2	H	0.62	0/1667	1.43	36/2279 (1.6%)
3	L	0.59	1/1621 (0.1%)	1.58	44/2200 (2.0%)
All	All	0.65	2/4073 (0.0%)	1.53	101/5539 (1.8%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
3	L	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	355	ASN	CA-C	8.96	1.64	1.52
3	L	115	ALA	CA-CB	-5.31	1.45	1.53

All (101) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	L	7	SER	CA-C-N	-14.60	104.72	119.76
3	L	7	SER	C-N-CA	-14.60	104.72	119.76
3	L	154	ILE	N-CA-C	13.07	123.95	110.62
3	L	115	ALA	N-CA-C	13.07	129.43	109.96
3	L	142	ASN	N-CA-C	11.38	124.23	110.91
2	H	214	GLU	CA-C-N	11.17	133.80	119.84
2	H	214	GLU	C-N-CA	11.17	133.80	119.84
3	L	170	GLN	N-CA-C	11.02	126.53	113.20
1	A	361	LYS	N-CA-C	-10.38	97.99	114.09
1	A	346	HIS	N-CA-C	10.35	129.46	113.72
1	A	352	ILE	O-C-N	-10.22	112.89	122.09

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	343	GLU	N-CA-C	-10.03	94.68	109.63
1	A	355	ASN	CA-C-N	8.85	130.91	119.84
1	A	355	ASN	C-N-CA	8.85	130.91	119.84
3	L	189	GLU	N-CA-C	-8.74	102.78	113.97
2	H	98	ARG	N-CA-C	8.54	123.56	112.12
1	A	317	HIS	N-CA-C	-8.52	101.21	111.69
2	H	214	GLU	N-CA-C	8.46	120.39	109.64
2	H	100	TYR	N-CA-C	-8.44	101.95	113.18
3	L	54	ARG	N-CA-C	-8.30	98.19	110.48
3	L	66	PHE	N-CA-C	-8.20	93.33	110.80
2	H	84	SER	N-CA-C	8.17	121.83	109.41
2	H	190	THR	N-CA-C	8.14	122.93	112.92
2	H	169	PHE	CA-C-N	7.92	129.74	119.84
2	H	169	PHE	C-N-CA	7.92	129.74	119.84
2	H	211	LYS	N-CA-C	7.91	122.35	109.85
3	L	6	GLN	N-CA-C	-7.87	97.24	109.76
3	L	64	THR	N-CA-C	7.76	124.10	111.37
3	L	8	PRO	N-CA-C	-7.70	99.14	111.14
3	L	113	ALA	N-CA-C	-7.45	99.99	110.50
2	H	149	PHE	CA-C-N	-7.45	111.96	119.93
2	H	149	PHE	C-N-CA	-7.45	111.96	119.93
2	H	163	SER	N-CA-C	-7.36	95.13	110.80
3	L	141	ASN	N-CA-C	7.30	122.23	109.96
3	L	118	THR	N-CA-C	-7.22	100.81	110.55
3	L	130	THR	N-CA-C	-7.18	104.65	113.19
1	A	328	GLY	N-CA-C	7.11	122.36	110.56
2	H	76	SER	N-CA-C	-6.86	99.12	110.17
2	H	86	LEU	N-CA-C	6.86	119.67	110.35
1	A	391	TRP	N-CA-C	-6.85	105.70	112.97
1	A	362	ASP	N-CA-C	-6.84	105.09	113.50
2	H	146	LYS	N-CA-C	6.83	120.30	108.76
1	A	299	TYR	N-CA-C	6.82	125.32	110.80
2	H	99	ASP	N-CA-C	6.80	125.28	110.80
3	L	98	ASP	CA-C-N	-6.77	103.46	120.53
3	L	98	ASP	C-N-CA	-6.77	103.46	120.53
3	L	122	PHE	CA-C-N	6.72	127.30	120.38
3	L	122	PHE	C-N-CA	6.72	127.30	120.38
3	L	7	SER	N-CA-C	6.67	124.56	109.81
1	A	345	ARG	N-CA-C	-6.67	104.17	113.37
2	H	191	TRP	N-CA-C	6.62	124.45	109.81
1	A	381	GLY	N-CA-C	6.62	121.05	112.18
3	L	28	SER	N-CA-C	6.57	120.01	110.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	L	148	ILE	N-CA-C	-6.49	99.14	108.42
2	H	56	GLY	N-CA-C	-6.48	105.98	115.63
2	H	93	VAL	N-CA-C	-6.46	97.88	107.51
2	H	75	SER	N-CA-C	-6.36	103.27	111.02
2	H	74	THR	N-CA-C	-6.35	105.66	113.41
2	H	138	SER	N-CA-C	6.29	119.74	109.24
2	H	159	SER	N-CA-C	-6.28	104.84	113.56
1	A	352	ILE	CA-C-N	-6.27	111.54	122.64
1	A	352	ILE	C-N-CA	-6.27	111.54	122.64
3	L	24	ARG	N-CA-C	6.22	119.53	107.71
2	H	165	GLY	N-CA-C	-6.20	98.50	113.18
3	L	89	THR	N-CA-C	-6.18	99.57	109.96
2	H	164	SER	N-CA-C	-6.16	102.12	110.68
1	A	319	THR	N-CA-C	6.14	119.56	109.85
3	L	117	PRO	N-CA-C	6.10	120.02	110.80
3	L	17	GLN	N-CA-C	6.02	118.11	109.14
3	L	206	VAL	N-CA-C	6.00	117.33	108.45
3	L	116	ALA	CA-C-N	6.00	126.20	119.90
3	L	116	ALA	C-N-CA	6.00	126.20	119.90
3	L	147	ASP	N-CA-C	5.97	119.21	109.24
3	L	64	THR	CA-C-N	-5.89	112.64	122.17
3	L	64	THR	C-N-CA	-5.89	112.64	122.17
3	L	3	VAL	N-CA-C	5.83	117.08	108.45
3	L	125	SER	N-CA-C	5.79	118.17	110.53
3	L	75	PHE	N-CA-C	5.78	119.34	108.65
3	L	18	ARG	N-CA-C	5.75	123.05	110.80
3	L	98	ASP	N-CA-C	5.70	122.40	109.81
3	L	122	PHE	N-CA-C	5.58	119.34	108.85
2	H	151	GLU	CA-C-N	-5.46	113.51	120.23
2	H	151	GLU	C-N-CA	-5.46	113.51	120.23
3	L	36	PHE	N-CA-C	5.44	117.51	110.65
2	H	174	GLN	CA-C-N	-5.43	115.12	121.64
2	H	174	GLN	C-N-CA	-5.43	115.12	121.64
2	H	142	GLY	N-CA-C	5.42	118.44	110.80
3	L	55	ALA	N-CA-C	5.39	122.28	110.80
2	H	7	SER	N-CA-C	5.38	118.59	109.76
1	A	360	GLU	N-CA-C	5.36	115.61	108.38
2	H	78	ALA	N-CA-C	5.24	117.78	109.50
3	L	150	VAL	N-CA-C	5.24	116.45	108.80
2	H	174	GLN	CA-CB-CG	-5.19	103.72	114.10
1	A	320	ILE	N-CA-C	-5.19	100.98	108.71
2	H	134	ALA	N-CA-C	5.15	116.56	109.15

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	356	PRO	N-CA-C	5.13	123.03	112.47
3	L	4	LEU	CA-CB-CG	-5.12	98.36	116.30
2	H	174	GLN	CB-CA-C	5.06	119.22	110.86
1	A	325	GLN	N-CA-C	-5.06	101.63	109.72
1	A	312	ILE	N-CA-C	-5.04	98.86	109.34
3	L	15	LEU	N-CA-C	-5.01	105.49	112.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	L	144	TYR	Sidechain

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	768	0	780	125	0
2	H	1625	0	1584	246	0
3	L	1587	0	1527	284	0
4	A	59	0	0	9	0
4	H	97	0	0	4	0
4	L	98	0	0	10	0
All	All	4234	0	3891	619	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 79.

All (619) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:4:LEU:HD13	3:L:103:GLY:HA2	1.22	1.17
3:L:58:LEU:H	3:L:58:LEU:HD22	1.12	1.15
3:L:115:ALA:CB	3:L:144:TYR:H	1.60	1.15
3:L:115:ALA:HB2	3:L:144:TYR:N	1.63	1.12

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:353:THR:HG22	1:A:356:PRO:HG2	1.35	1.08
2:H:40:ARG:HB3	2:H:41:PRO:HD2	1.32	1.06
1:A:355:ASN:N	1:A:356:PRO:HD2	1.69	1.04
3:L:115:ALA:HB2	3:L:144:TYR:H	0.90	1.04
1:A:343:GLU:HG3	1:A:344:LYS:H	1.17	1.04
3:L:96:ASN:HD22	3:L:96:ASN:C	1.61	1.03
3:L:77:LEU:HD12	3:L:78:THR:H	1.22	1.02
3:L:112:ARG:NH2	3:L:113:ALA:HB3	1.74	1.01
3:L:55:ALA:O	3:L:68:GLY:HA3	1.61	1.00
1:A:353:THR:CG2	1:A:356:PRO:HG2	1.90	1.00
2:H:191:TRP:HB3	2:H:192:PRO:HD3	1.46	0.98
2:H:170:PRO:HD2	3:L:167:TRP:O	1.64	0.97
3:L:58:LEU:HD22	3:L:58:LEU:N	1.79	0.97
3:L:58:LEU:H	3:L:58:LEU:CD2	1.79	0.96
3:L:140:LEU:H	3:L:140:LEU:HD12	1.31	0.95
3:L:4:LEU:CD1	3:L:103:GLY:HA2	1.97	0.94
2:H:100:TYR:O	3:L:95:THR:HG21	1.68	0.94
2:H:187:THR:HB	2:H:190:THR:OG1	1.67	0.94
2:H:212:LYS:HA	4:H:281:HOH:O	1.67	0.92
3:L:150:VAL:HG13	3:L:199:GLU:O	1.68	0.92
2:H:140:THR:HB	2:H:185:THR:HG22	1.52	0.91
3:L:4:LEU:HD13	3:L:103:GLY:CA	1.99	0.91
3:L:30:VAL:HG23	3:L:72:ARG:HA	1.49	0.91
3:L:41:GLN:HB2	3:L:51:LEU:HD21	1.50	0.91
3:L:77:LEU:HD12	3:L:78:THR:N	1.85	0.90
3:L:12:ALA:HB1	3:L:111:LYS:HE3	1.51	0.90
3:L:80:ASN:HB3	3:L:81:PRO:HD3	1.52	0.90
1:A:352:ILE:HG22	1:A:352:ILE:O	1.71	0.89
1:A:355:ASN:O	1:A:355:ASN:ND2	2.05	0.89
1:A:355:ASN:N	1:A:356:PRO:CD	2.35	0.88
3:L:18:ARG:NH1	3:L:18:ARG:HB3	1.88	0.88
2:H:129:PRO:HG3	3:L:122:PHE:HE2	1.38	0.87
2:H:200:VAL:HB	2:H:209:VAL:HG13	1.54	0.87
3:L:37:MET:HE3	3:L:37:MET:HA	1.56	0.87
3:L:200:ALA:HB3	3:L:205:ILE:HG23	1.56	0.86
1:A:343:GLU:HG3	1:A:344:LYS:N	1.85	0.86
1:A:353:THR:HG22	1:A:356:PRO:CG	2.06	0.85
1:A:344:LYS:C	1:A:346:HIS:H	1.84	0.84
3:L:97:VAL:HG12	3:L:99:PRO:HD2	1.59	0.84
3:L:4:LEU:HD12	3:L:4:LEU:O	1.78	0.84
3:L:179:MET:HG2	3:L:180:SER:H	1.43	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:102:PHE:N	3:L:102:PHE:HD2	1.74	0.84
3:L:119:VAL:HB	3:L:207:LYS:HD3	1.57	0.84
2:H:50:ARG:HG3	2:H:50:ARG:HH11	1.40	0.83
3:L:5:THR:HG23	3:L:24:ARG:CB	2.08	0.83
2:H:11:LEU:HD11	2:H:115:SER:HB3	1.61	0.83
2:H:200:VAL:HB	2:H:209:VAL:CG1	2.08	0.83
3:L:115:ALA:HB1	3:L:143:PHE:HA	1.61	0.82
1:A:340:MET:HE3	1:A:346:HIS:HA	1.61	0.82
3:L:40:TYR:HB2	3:L:91:TYR:HB2	1.62	0.82
2:H:37:VAL:HG11	2:H:45:LEU:HD23	1.61	0.82
2:H:166:VAL:HG13	2:H:183:SER:O	1.80	0.81
3:L:163:VAL:C	3:L:164:LEU:HD23	2.05	0.81
2:H:14:PRO:HG3	2:H:114:VAL:HG12	1.63	0.81
3:L:102:PHE:N	3:L:102:PHE:CD2	2.48	0.81
2:H:6:GLN:HG2	2:H:96:CYS:SG	2.21	0.81
2:H:174:GLN:OE1	3:L:164:LEU:HD22	1.80	0.81
3:L:200:ALA:HB3	3:L:205:ILE:CG2	2.10	0.80
3:L:162:GLY:O	3:L:183:LEU:HA	1.80	0.80
1:A:306:PHE:O	1:A:387:LEU:HD21	1.83	0.79
2:H:45:LEU:HD21	3:L:48:PRO:HG2	1.64	0.79
3:L:28:SER:HA	3:L:73:THR:OG1	1.83	0.79
3:L:159:ARG:NH2	3:L:189:GLU:HG2	1.96	0.79
3:L:96:ASN:HD22	3:L:97:VAL:N	1.81	0.78
2:H:4:LEU:HD21	2:H:27:PHE:HZ	1.49	0.78
2:H:9:ALA:HB3	2:H:203:PRO:HB2	1.66	0.78
1:A:319:THR:HG21	1:A:368:GLU:OE2	1.84	0.78
1:A:366:ASN:HD22	1:A:367:ILE:H	1.32	0.78
2:H:29:ILE:HG12	2:H:53:PRO:HG2	1.64	0.77
2:H:40:ARG:HB3	2:H:41:PRO:CD	2.12	0.77
1:A:354:VAL:C	1:A:356:PRO:CD	2.57	0.77
1:A:350:ARG:HB3	1:A:370:GLU:HB3	1.64	0.77
1:A:352:ILE:O	1:A:352:ILE:CG2	2.33	0.77
3:L:112:ARG:HH21	3:L:113:ALA:HB3	1.49	0.77
1:A:366:ASN:C	1:A:367:ILE:HD12	2.09	0.77
1:A:344:LYS:HE3	4:A:425:HOH:O	1.84	0.77
3:L:140:LEU:HD12	3:L:140:LEU:N	1.99	0.77
3:L:183:LEU:HG	3:L:185:LEU:HD21	1.67	0.76
3:L:155:ASP:HA	3:L:195:SER:O	1.86	0.76
2:H:29:ILE:HG13	2:H:34:MET:CE	2.16	0.75
3:L:84:ALA:HA	3:L:110:ILE:HD12	1.68	0.75
2:H:170:PRO:CD	3:L:167:TRP:O	2.34	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:140:LEU:HD13	3:L:179:MET:HB3	1.67	0.75
1:A:366:ASN:HD22	1:A:367:ILE:N	1.85	0.75
2:H:38:LYS:HB2	2:H:48:ILE:HD11	1.68	0.75
3:L:192:ARG:HG3	3:L:192:ARG:HH11	1.52	0.75
2:H:127:LEU:HD11	2:H:144:LEU:HB2	1.66	0.74
2:H:40:ARG:CB	2:H:41:PRO:HD2	2.14	0.74
2:H:129:PRO:HG3	3:L:122:PHE:CE2	2.22	0.74
2:H:162:LEU:O	2:H:164:SER:N	2.21	0.74
3:L:95:THR:HG22	3:L:95:THR:O	1.88	0.74
2:H:211:LYS:HZ3	2:H:212:LYS:HE3	1.53	0.74
1:A:369:ALA:O	1:A:371:PRO:HD3	1.88	0.73
3:L:4:LEU:HD22	3:L:102:PHE:C	2.12	0.73
2:H:2:VAL:HG13	2:H:27:PHE:CD2	2.23	0.73
3:L:58:LEU:HD11	4:L:245:HOH:O	1.87	0.73
1:A:358:VAL:HG22	1:A:365:VAL:HG21	1.71	0.72
2:H:149:PHE:HB3	2:H:150:PRO:HD3	1.71	0.72
2:H:191:TRP:CD1	2:H:196:ILE:HG13	2.26	0.71
2:H:97:ALA:HB3	2:H:103:PHE:HD1	1.55	0.71
3:L:187:LYS:O	3:L:187:LYS:HG2	1.89	0.71
2:H:33:TYR:O	2:H:98:ARG:O	2.09	0.71
2:H:29:ILE:HG13	2:H:34:MET:HE1	1.73	0.70
2:H:124:VAL:HG22	2:H:211:LYS:HG3	1.73	0.70
2:H:139:VAL:HG21	4:H:283:HOH:O	1.91	0.70
2:H:169:PHE:CD2	3:L:180:SER:HB2	2.26	0.70
2:H:166:VAL:HG12	2:H:167:HIS:N	2.05	0.70
2:H:47:TRP:HB2	3:L:102:PHE:HE2	1.56	0.70
3:L:18:ARG:HA	3:L:80:ASN:HA	1.74	0.70
3:L:213:GLU:HB2	4:L:299:HOH:O	1.92	0.70
3:L:13:VAL:HG11	3:L:82:VAL:HG21	1.74	0.69
2:H:99:ASP:HB2	2:H:102:GLY:O	1.92	0.69
1:A:365:VAL:HA	4:A:450:HOH:O	1.92	0.69
2:H:2:VAL:HA	2:H:25:SER:OG	1.92	0.69
3:L:30:VAL:CG2	3:L:72:ARG:HA	2.22	0.69
2:H:29:ILE:HG12	2:H:53:PRO:CG	2.22	0.69
2:H:102:GLY:N	3:L:38:HIS:ND1	2.40	0.69
2:H:89:GLU:C	2:H:91:THR:H	2.00	0.69
2:H:190:THR:O	2:H:190:THR:HG22	1.93	0.69
3:L:179:MET:HG2	3:L:180:SER:N	2.06	0.69
1:A:354:VAL:C	1:A:356:PRO:HD3	2.17	0.69
3:L:18:ARG:HB3	3:L:18:ARG:HH11	1.58	0.69
3:L:183:LEU:HG	3:L:185:LEU:CD2	2.23	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:184:VAL:HG13	2:H:184:VAL:O	1.92	0.68
3:L:80:ASN:O	3:L:82:VAL:N	2.27	0.68
3:L:150:VAL:HG12	3:L:151:LYS:N	2.09	0.68
3:L:15:LEU:HD23	3:L:83:GLU:HA	1.76	0.68
3:L:11:LEU:H	3:L:11:LEU:HD23	1.59	0.68
1:A:354:VAL:C	1:A:356:PRO:HD2	2.19	0.68
2:H:30:LYS:HA	2:H:53:PRO:HB2	1.74	0.68
3:L:5:THR:CG2	3:L:24:ARG:HB3	2.24	0.68
2:H:17:SER:HB2	2:H:84:SER:HB2	1.76	0.67
2:H:67:LYS:NZ	2:H:86:LEU:HA	2.10	0.67
3:L:70:GLY:HA3	3:L:75:PHE:CD1	2.30	0.67
3:L:13:VAL:HG12	3:L:14:SER:N	2.10	0.67
1:A:353:THR:HG21	1:A:356:PRO:HG2	1.77	0.67
2:H:73:ASP:O	2:H:76:SER:O	2.13	0.67
1:A:376:SER:H	1:A:392:PHE:HA	1.60	0.67
2:H:102:GLY:H	3:L:38:HIS:CE1	2.12	0.67
3:L:5:THR:HG22	3:L:24:ARG:O	1.95	0.67
3:L:6:GLN:NE2	3:L:90:TYR:O	2.27	0.67
3:L:96:ASN:C	3:L:96:ASN:ND2	2.38	0.66
1:A:315:THR:HG22	1:A:317:HIS:H	1.60	0.66
3:L:194:ASN:O	3:L:211:ARG:HG3	1.95	0.66
2:H:71:THR:HG22	2:H:72:ALA:N	2.10	0.66
3:L:115:ALA:CB	3:L:144:TYR:N	2.38	0.66
2:H:111:LEU:HD13	2:H:152:PRO:HG3	1.77	0.65
2:H:191:TRP:HB3	2:H:192:PRO:CD	2.24	0.65
3:L:49:LYS:NZ	3:L:49:LYS:HB2	2.12	0.65
3:L:95:THR:HG23	3:L:100:TRP:CE3	2.32	0.65
3:L:115:ALA:CB	3:L:143:PHE:HA	2.26	0.65
3:L:96:ASN:ND2	3:L:97:VAL:N	2.45	0.65
2:H:141:LEU:HA	3:L:122:PHE:CZ	2.31	0.65
2:H:216:ARG:CD	3:L:126:SER:OG	2.44	0.65
2:H:173:LEU:HD11	2:H:176:ASP:HA	1.77	0.65
3:L:191:GLU:C	3:L:193:HIS:H	2.05	0.65
3:L:98:ASP:O	3:L:100:TRP:N	2.28	0.65
3:L:148:ILE:HG23	3:L:201:THR:OG1	1.97	0.65
3:L:115:ALA:HB1	3:L:143:PHE:CA	2.27	0.65
3:L:155:ASP:OD1	3:L:194:ASN:ND2	2.30	0.64
2:H:140:THR:CB	2:H:185:THR:HG22	2.27	0.64
1:A:382:VAL:O	1:A:386:GLN:HB2	1.97	0.64
3:L:5:THR:HG23	3:L:24:ARG:HB3	1.79	0.64
1:A:359:THR:O	1:A:359:THR:CG2	2.45	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:169:PHE:O	2:H:180:LEU:HG	1.98	0.64
1:A:343:GLU:CG	1:A:344:LYS:N	2.60	0.64
2:H:126:PRO:HG3	2:H:212:LYS:O	1.98	0.64
3:L:21:ILE:HD12	3:L:21:ILE:N	2.12	0.64
1:A:394:LYS:HD3	4:A:415:HOH:O	1.97	0.63
3:L:136:VAL:HG12	3:L:152:TRP:CH2	2.33	0.63
2:H:157:TRP:HZ3	2:H:213:ILE:CD1	2.11	0.63
1:A:317:HIS:NE2	4:A:427:HOH:O	2.30	0.63
2:H:213:ILE:O	2:H:213:ILE:HG22	1.99	0.63
2:H:4:LEU:HD21	2:H:27:PHE:CZ	2.33	0.63
2:H:186:VAL:HG23	2:H:187:THR:O	1.99	0.63
3:L:134:ALA:HB3	3:L:185:LEU:HG	1.79	0.63
2:H:111:LEU:HD23	2:H:112:VAL:N	2.14	0.63
2:H:2:VAL:HG13	2:H:27:PHE:HD2	1.62	0.63
1:A:344:LYS:C	1:A:346:HIS:N	2.54	0.62
2:H:45:LEU:CD2	3:L:48:PRO:HG2	2.28	0.62
2:H:71:THR:CG2	2:H:72:ALA:N	2.60	0.62
3:L:84:ALA:HA	3:L:110:ILE:CD1	2.28	0.62
1:A:382:VAL:HG12	1:A:383:GLU:H	1.64	0.62
2:H:97:ALA:CB	2:H:103:PHE:HD1	2.12	0.62
3:L:188:ASP:OD1	3:L:188:ASP:C	2.42	0.62
2:H:166:VAL:CG1	2:H:167:HIS:H	2.13	0.62
2:H:20:LEU:N	2:H:20:LEU:HD23	2.15	0.61
3:L:138:CYS:O	3:L:138:CYS:SG	2.58	0.61
2:H:97:ALA:HB3	2:H:103:PHE:CD1	2.35	0.61
3:L:12:ALA:HB1	3:L:111:LYS:CE	2.29	0.61
3:L:169:ASP:OD1	3:L:169:ASP:N	2.32	0.61
3:L:164:LEU:HD23	3:L:164:LEU:N	2.15	0.61
2:H:67:LYS:HZ2	2:H:86:LEU:HA	1.65	0.61
3:L:140:LEU:H	3:L:140:LEU:CD1	2.11	0.61
3:L:38:HIS:CD2	3:L:53:TYR:O	2.54	0.61
2:H:166:VAL:CG1	2:H:167:HIS:N	2.63	0.60
3:L:7:SER:O	3:L:8:PRO:C	2.35	0.60
2:H:40:ARG:HG2	2:H:91:THR:O	2.00	0.60
1:A:346:HIS:O	1:A:347:VAL:HG23	2.01	0.60
3:L:59:GLU:OE2	3:L:60:SER:N	2.27	0.60
3:L:94:GLN:HG2	3:L:95:THR:N	2.15	0.60
1:A:315:THR:HB	1:A:319:THR:HG23	1.83	0.60
1:A:340:MET:CE	1:A:346:HIS:HA	2.32	0.60
2:H:126:PRO:HA	2:H:143:CYS:HB2	1.83	0.60
2:H:108:GLN:O	2:H:109:GLY:C	2.45	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:142:ASN:HA	3:L:176:THR:OG1	2.02	0.60
1:A:378:ILE:HG22	1:A:379:ILE:H	1.67	0.59
2:H:143:CYS:N	2:H:182:SER:OG	2.35	0.59
3:L:144:TYR:O	3:L:145:PRO:C	2.43	0.59
2:H:36:TRP:NE1	2:H:80:TYR:O	2.34	0.59
3:L:114:ASP:N	3:L:114:ASP:OD1	2.35	0.59
1:A:321:VAL:HG12	1:A:366:ASN:HD21	1.68	0.59
3:L:11:LEU:HD23	3:L:11:LEU:N	2.16	0.59
2:H:91:THR:HG23	2:H:113:THR:HA	1.85	0.58
2:H:157:TRP:CZ3	2:H:213:ILE:HD11	2.38	0.58
3:L:5:THR:HG23	3:L:24:ARG:HB2	1.85	0.58
3:L:4:LEU:HD22	3:L:103:GLY:HA2	1.85	0.58
2:H:59:LYS:HD3	3:L:98:ASP:CG	2.28	0.58
2:H:89:GLU:C	2:H:91:THR:N	2.60	0.58
2:H:169:PHE:CD2	3:L:168:THR:HG23	2.38	0.58
3:L:73:THR:HG22	3:L:73:THR:O	2.02	0.58
3:L:150:VAL:HG12	3:L:151:LYS:O	2.03	0.58
3:L:19:ALA:HB3	3:L:79:ILE:HB	1.85	0.58
1:A:375:ASP:HB3	1:A:392:PHE:HB2	1.85	0.58
3:L:79:ILE:HD11	3:L:90:TYR:HE2	1.66	0.58
2:H:166:VAL:HG12	2:H:167:HIS:H	1.69	0.58
2:H:191:TRP:CB	2:H:192:PRO:HD3	2.30	0.58
1:A:359:THR:O	1:A:359:THR:HG22	2.04	0.57
1:A:312:ILE:HG23	1:A:322:ILE:HG12	1.85	0.57
2:H:52:ASP:OD2	2:H:54:ALA:HB3	2.05	0.57
2:H:103:PHE:N	2:H:103:PHE:HD2	2.03	0.57
2:H:153:VAL:HA	2:H:201:ALA:O	2.04	0.57
2:H:211:LYS:NZ	2:H:212:LYS:HE3	2.18	0.57
3:L:42:GLN:O	3:L:88:ALA:HB1	2.04	0.57
2:H:149:PHE:O	2:H:150:PRO:C	2.46	0.57
3:L:28:SER:HB3	3:L:72:ARG:HG2	1.87	0.57
1:A:331:SER:N	1:A:332:PRO:HD3	2.19	0.57
3:L:12:ALA:CB	3:L:111:LYS:HE3	2.31	0.57
2:H:124:VAL:CG1	2:H:209:VAL:HG21	2.34	0.57
3:L:153:LYS:HB2	3:L:197:THR:HB	1.85	0.57
2:H:102:GLY:N	3:L:38:HIS:CE1	2.73	0.57
2:H:140:THR:HA	2:H:185:THR:HA	1.86	0.57
3:L:19:ALA:HB2	3:L:82:VAL:CG2	2.35	0.57
3:L:159:ARG:HH22	3:L:189:GLU:HG2	1.70	0.57
1:A:348:LEU:HD22	1:A:372:PRO:HG3	1.87	0.57
2:H:126:PRO:O	2:H:127:LEU:HD23	2.05	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:157:TRP:HZ2	2:H:182:SER:HB2	1.69	0.57
1:A:319:THR:OG1	1:A:368:GLU:HG3	2.05	0.56
1:A:320:ILE:HG13	1:A:320:ILE:O	2.05	0.56
1:A:357:ILE:HD13	1:A:357:ILE:H	1.69	0.56
3:L:108:LEU:HD12	3:L:109:GLU:N	2.20	0.56
3:L:164:LEU:HG	3:L:182:THR:OG1	2.04	0.56
2:H:50:ARG:HG3	2:H:50:ARG:NH1	2.16	0.56
3:L:80:ASN:HB3	3:L:81:PRO:CD	2.29	0.56
2:H:143:CYS:C	2:H:182:SER:HG	2.14	0.56
1:A:327:GLU:HG3	4:A:404:HOH:O	2.05	0.56
2:H:36:TRP:CE3	2:H:95:PHE:O	2.59	0.56
2:H:216:ARG:HD2	3:L:126:SER:OG	2.05	0.56
3:L:115:ALA:CB	3:L:143:PHE:CA	2.83	0.56
2:H:196:ILE:HG22	2:H:197:THR:N	2.21	0.56
1:A:337:PHE:C	1:A:338:GLU:CG	2.78	0.56
1:A:359:THR:O	1:A:360:GLU:HG3	2.06	0.56
1:A:387:LEU:HD21	3:L:54:ARG:HH22	1.71	0.56
2:H:144:LEU:HD23	2:H:145:VAL:N	2.21	0.56
2:H:126:PRO:HB2	2:H:213:ILE:HA	1.88	0.56
3:L:65:ARG:HB2	3:L:81:PRO:HD2	1.88	0.56
3:L:4:LEU:CG	3:L:103:GLY:HA2	2.36	0.55
2:H:187:THR:CB	2:H:190:THR:OG1	2.50	0.55
2:H:191:TRP:HZ2	2:H:213:ILE:O	1.88	0.55
1:A:355:ASN:H	1:A:356:PRO:HD2	1.65	0.55
2:H:60:TYR:N	2:H:60:TYR:CD1	2.73	0.55
2:H:129:PRO:CG	3:L:122:PHE:HE2	2.14	0.55
2:H:157:TRP:O	2:H:158:ASN:HB2	2.06	0.55
1:A:301:MET:HB2	1:A:336:PRO:HG3	1.88	0.55
1:A:392:PHE:HD1	1:A:392:PHE:O	1.89	0.55
1:A:310:LYS:HE2	2:H:52:ASP:OD1	2.07	0.55
3:L:95:THR:HG23	3:L:100:TRP:CZ3	2.41	0.55
3:L:165:ASN:HB3	3:L:179:MET:HE3	1.88	0.55
1:A:333:CYS:O	1:A:357:ILE:HB	2.06	0.55
3:L:37:MET:CE	3:L:93:GLN:O	2.55	0.55
3:L:70:GLY:HA3	3:L:75:PHE:HD1	1.72	0.55
3:L:6:GLN:HA	3:L:22:SER:O	2.06	0.55
3:L:120:SER:O	3:L:138:CYS:HA	2.06	0.55
3:L:38:HIS:CD2	3:L:38:HIS:N	2.72	0.55
1:A:340:MET:HE3	1:A:346:HIS:CA	2.33	0.55
2:H:12:VAL:HG21	2:H:18:VAL:CG1	2.37	0.55
3:L:112:ARG:CZ	3:L:113:ALA:HB3	2.37	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:133:GLY:HA3	3:L:185:LEU:O	2.07	0.54
1:A:339:ILE:HG23	1:A:377:TYR:O	2.07	0.54
2:H:6:GLN:HA	2:H:21:SER:O	2.06	0.54
1:A:387:LEU:CD2	3:L:54:ARG:NH2	2.71	0.54
3:L:41:GLN:CB	3:L:51:LEU:HD21	2.32	0.54
1:A:333:CYS:HB3	4:A:433:HOH:O	2.07	0.54
3:L:209:PHE:CD1	3:L:209:PHE:C	2.86	0.54
1:A:378:ILE:HG22	1:A:379:ILE:N	2.22	0.54
2:H:103:PHE:N	2:H:103:PHE:CD2	2.74	0.54
2:H:144:LEU:HD21	2:H:146:LYS:HB3	1.88	0.54
1:A:350:ARG:NH1	1:A:370:GLU:OE1	2.41	0.54
3:L:212:ASN:O	3:L:213:GLU:O	2.26	0.54
3:L:18:ARG:HG2	3:L:80:ASN:ND2	2.23	0.53
1:A:299:TYR:CD2	1:A:299:TYR:N	2.76	0.53
2:H:127:LEU:HG	2:H:143:CYS:HA	1.91	0.53
3:L:203:SER:OG	3:L:204:PRO:HD2	2.08	0.53
1:A:309:VAL:HG23	1:A:309:VAL:O	2.09	0.53
3:L:124:PRO:HG3	3:L:135:SER:H	1.73	0.53
2:H:45:LEU:HD12	2:H:45:LEU:H	1.72	0.53
3:L:168:THR:OG1	3:L:178:SER:O	2.25	0.53
3:L:192:ARG:HH11	3:L:192:ARG:CG	2.17	0.53
1:A:331:SER:H	1:A:332:PRO:HD3	1.74	0.53
3:L:58:LEU:HD21	4:L:245:HOH:O	2.07	0.53
3:L:150:VAL:CG1	3:L:151:LYS:N	2.72	0.53
3:L:144:TYR:HB3	3:L:145:PRO:HD3	1.90	0.53
3:L:159:ARG:NH1	3:L:185:LEU:HD22	2.23	0.53
2:H:13:LYS:HA	2:H:115:SER:O	2.09	0.53
3:L:163:VAL:HA	3:L:182:THR:O	2.09	0.53
3:L:75:PHE:C	3:L:76:THR:HG22	2.32	0.53
3:L:79:ILE:HD11	3:L:90:TYR:CE2	2.43	0.53
3:L:136:VAL:CG1	3:L:152:TRP:CH2	2.91	0.53
2:H:201:ALA:O	2:H:203:PRO:HD3	2.08	0.52
2:H:191:TRP:CH2	2:H:215:PRO:HA	2.43	0.52
1:A:379:ILE:O	1:A:380:ILE:HG12	2.08	0.52
2:H:17:SER:HB2	2:H:84:SER:CB	2.39	0.52
2:H:36:TRP:HE3	2:H:95:PHE:O	1.91	0.52
2:H:162:LEU:C	2:H:164:SER:N	2.67	0.52
2:H:197:THR:HG23	2:H:198:CYS:N	2.25	0.52
3:L:4:LEU:HD11	4:L:219:HOH:O	2.08	0.52
3:L:185:LEU:HD12	3:L:190:TYR:HB2	1.91	0.52
2:H:157:TRP:CZ3	2:H:213:ILE:CD1	2.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:8:PRO:HB2	3:L:11:LEU:HD22	1.92	0.52
3:L:18:ARG:CG	3:L:80:ASN:ND2	2.73	0.52
3:L:63:PRO:HG2	3:L:66:PHE:HE2	1.75	0.52
3:L:154:ILE:O	3:L:195:SER:O	2.28	0.52
3:L:18:ARG:CZ	3:L:18:ARG:CB	2.87	0.52
2:H:20:LEU:HB2	2:H:81:LEU:HB3	1.91	0.52
3:L:37:MET:HG2	3:L:75:PHE:CD2	2.43	0.52
3:L:93:GLN:HG2	3:L:94:GLN:N	2.25	0.52
2:H:52:ASP:O	2:H:56:GLY:N	2.43	0.52
3:L:20:THR:C	3:L:21:ILE:HD12	2.35	0.52
2:H:33:TYR:CE2	2:H:52:ASP:OD1	2.63	0.52
1:A:360:GLU:HB3	1:A:362:ASP:H	1.74	0.52
2:H:144:LEU:CD2	2:H:146:LYS:HB3	2.40	0.52
3:L:168:THR:OG1	3:L:178:SER:C	2.53	0.52
3:L:52:ILE:HD11	3:L:67:SER:HA	1.92	0.51
3:L:30:VAL:HA	3:L:35:SER:HA	1.91	0.51
2:H:62:PRO:C	2:H:64:PHE:H	2.18	0.51
2:H:155:LEU:HD21	2:H:157:TRP:HE1	1.76	0.51
3:L:151:LYS:HZ2	3:L:199:GLU:CD	2.19	0.51
2:H:71:THR:CG2	2:H:72:ALA:H	2.22	0.51
2:H:2:VAL:O	2:H:3:GLN:HG3	2.11	0.51
3:L:4:LEU:HD22	3:L:103:GLY:N	2.24	0.51
2:H:174:GLN:CD	3:L:164:LEU:HD22	2.35	0.51
2:H:200:VAL:CB	2:H:209:VAL:HG13	2.34	0.51
1:A:376:SER:H	1:A:392:PHE:CA	2.23	0.51
1:A:376:SER:N	1:A:392:PHE:HB3	2.27	0.51
3:L:4:LEU:CD2	3:L:103:GLY:HA2	2.40	0.51
2:H:135:THR:HG21	2:H:139:VAL:HA	1.92	0.50
2:H:141:LEU:HA	3:L:122:PHE:HZ	1.76	0.50
2:H:157:TRP:CD1	2:H:166:VAL:HG21	2.46	0.50
2:H:152:PRO:O	2:H:203:PRO:HD2	2.11	0.50
3:L:59:GLU:CD	3:L:60:SER:H	2.15	0.50
3:L:12:ALA:HA	3:L:109:GLU:O	2.11	0.50
3:L:136:VAL:CG1	3:L:152:TRP:CZ3	2.94	0.50
3:L:146:LYS:HD2	3:L:167:TRP:CZ3	2.47	0.50
2:H:28:ASN:ND2	2:H:30:LYS:HB2	2.27	0.50
3:L:27:GLU:O	3:L:73:THR:HG23	2.11	0.50
3:L:95:THR:HA	3:L:100:TRP:CG	2.47	0.50
3:L:194:ASN:HA	3:L:211:ARG:HD3	1.93	0.50
2:H:143:CYS:O	2:H:182:SER:OG	2.24	0.50
1:A:382:VAL:HG12	1:A:383:GLU:N	2.26	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:12:VAL:HG11	2:H:18:VAL:HG12	1.93	0.50
2:H:106:TRP:HZ2	3:L:40:TYR:CE2	2.29	0.50
1:A:353:THR:HG22	1:A:354:VAL:N	2.27	0.49
3:L:29:VAL:HG22	3:L:72:ARG:O	2.12	0.49
3:L:138:CYS:HB3	3:L:152:TRP:HZ2	1.77	0.49
1:A:304:GLY:HA3	1:A:328:GLY:HA3	1.94	0.49
3:L:12:ALA:HB3	3:L:111:LYS:NZ	2.26	0.49
3:L:40:TYR:HA	3:L:49:LYS:O	2.11	0.49
3:L:83:GLU:C	3:L:85:ASP:H	2.20	0.49
1:A:367:ILE:HD12	1:A:367:ILE:N	2.27	0.49
1:A:380:ILE:HB	1:A:387:LEU:HB2	1.93	0.49
2:H:148:TYR:HE1	2:H:151:GLU:HA	1.78	0.49
2:H:162:LEU:HG	2:H:163:SER:H	1.77	0.49
3:L:18:ARG:HB3	3:L:18:ARG:CZ	2.40	0.49
3:L:86:ASP:O	3:L:108:LEU:HD23	2.13	0.49
3:L:111:LYS:HG2	3:L:144:TYR:OH	2.13	0.49
3:L:210:ASN:O	3:L:212:ASN:N	2.45	0.49
2:H:93:VAL:HA	2:H:111:LEU:HA	1.94	0.49
3:L:13:VAL:CG1	3:L:14:SER:N	2.74	0.49
3:L:49:LYS:HB2	3:L:49:LYS:HZ2	1.77	0.49
1:A:353:THR:CG2	1:A:356:PRO:CG	2.75	0.49
3:L:109:GLU:OE2	3:L:177:TYR:OH	2.31	0.49
2:H:35:HIS:CD2	2:H:35:HIS:N	2.81	0.49
1:A:303:THR:HG23	1:A:329:ASP:HB2	1.95	0.49
2:H:40:ARG:O	2:H:41:PRO:C	2.56	0.49
3:L:30:VAL:HB	3:L:72:ARG:HG3	1.94	0.49
3:L:54:ARG:O	3:L:56:SER:N	2.46	0.49
3:L:81:PRO:O	3:L:83:GLU:HG3	2.13	0.49
3:L:144:TYR:CD1	3:L:144:TYR:C	2.91	0.49
3:L:151:LYS:HB2	3:L:199:GLU:HB3	1.95	0.48
3:L:191:GLU:O	3:L:193:HIS:N	2.45	0.48
3:L:1:ASP:HA	4:L:268:HOH:O	2.12	0.48
3:L:146:LYS:HD2	3:L:167:TRP:CH2	2.48	0.48
1:A:337:PHE:CE2	1:A:339:ILE:HG13	2.49	0.48
2:H:124:VAL:HA	2:H:144:LEU:O	2.14	0.48
2:H:101:GLU:HA	3:L:38:HIS:CE1	2.49	0.48
3:L:151:LYS:HZ1	3:L:199:GLU:HG2	1.78	0.48
1:A:302:CYS:HB2	1:A:336:PRO:HD3	1.95	0.48
1:A:387:LEU:HD22	3:L:54:ARG:NH2	2.29	0.48
1:A:365:VAL:HG12	1:A:367:ILE:HD11	1.96	0.48
2:H:158:ASN:C	2:H:160:GLY:H	2.20	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:5:THR:CG2	3:L:24:ARG:O	2.61	0.48
2:H:3:GLN:O	2:H:25:SER:HB3	2.13	0.48
2:H:163:SER:C	2:H:165:GLY:H	2.22	0.48
1:A:302:CYS:SG	1:A:326:TYR:CZ	3.07	0.48
2:H:14:PRO:HG3	2:H:114:VAL:CG1	2.40	0.48
3:L:18:ARG:NH2	4:L:218:HOH:O	2.47	0.48
3:L:43:LYS:HB3	3:L:44:PRO:HD2	1.96	0.48
3:L:80:ASN:CB	3:L:81:PRO:HD3	2.33	0.48
3:L:192:ARG:CG	3:L:192:ARG:NH1	2.77	0.48
1:A:303:THR:HA	1:A:382:VAL:HG21	1.95	0.48
1:A:357:ILE:HD13	1:A:357:ILE:N	2.29	0.48
2:H:157:TRP:HZ3	2:H:213:ILE:HD11	1.76	0.48
3:L:37:MET:C	3:L:38:HIS:HD2	2.21	0.48
2:H:36:TRP:HA	2:H:95:PHE:O	2.13	0.47
2:H:61:ASP:O	2:H:62:PRO:C	2.56	0.47
2:H:83:LEU:HD11	2:H:94:TYR:CD2	2.49	0.47
2:H:108:GLN:O	2:H:108:GLN:OE1	2.32	0.47
3:L:4:LEU:HD22	3:L:103:GLY:CA	2.43	0.47
3:L:195:SER:HB2	3:L:210:ASN:OD1	2.14	0.47
3:L:140:LEU:HD21	3:L:150:VAL:HG21	1.95	0.47
2:H:9:ALA:O	2:H:204:ALA:HA	2.14	0.47
2:H:158:ASN:C	2:H:160:GLY:N	2.70	0.47
3:L:89:THR:HA	3:L:106:THR:O	2.14	0.47
2:H:141:LEU:N	2:H:141:LEU:HD23	2.30	0.47
1:A:306:PHE:CZ	1:A:335:ILE:HG13	2.49	0.47
2:H:72:ALA:HA	2:H:79:ALA:HA	1.95	0.47
2:H:89:GLU:O	2:H:91:THR:N	2.35	0.47
2:H:157:TRP:CZ2	2:H:182:SER:HB2	2.50	0.47
2:H:130:GLY:HA3	4:L:216:HOH:O	2.14	0.47
3:L:32:TYR:C	3:L:34:ASN:H	2.23	0.47
3:L:95:THR:HG23	3:L:100:TRP:CD2	2.50	0.47
3:L:43:LYS:HB3	3:L:44:PRO:CD	2.45	0.47
3:L:63:PRO:HG2	3:L:66:PHE:CE2	2.49	0.47
3:L:66:PHE:CD1	3:L:79:ILE:HG12	2.50	0.46
3:L:119:VAL:CB	3:L:207:LYS:HD3	2.39	0.46
3:L:123:PRO:HB3	3:L:209:PHE:CZ	2.50	0.46
2:H:184:VAL:O	2:H:184:VAL:CG1	2.60	0.46
2:H:188:SER:C	2:H:190:THR:H	2.22	0.46
3:L:98:ASP:C	3:L:100:TRP:N	2.74	0.46
3:L:191:GLU:HG3	3:L:192:ARG:HD3	1.97	0.46
1:A:321:VAL:HA	1:A:367:ILE:O	2.14	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:337:PHE:O	1:A:338:GLU:HG2	2.16	0.46
2:H:42:GLU:O	2:H:42:GLU:HG2	2.14	0.46
3:L:13:VAL:HG12	3:L:14:SER:H	1.79	0.46
3:L:187:LYS:HE2	3:L:187:LYS:HB3	1.64	0.46
1:A:360:GLU:HB2	1:A:363:SER:HB2	1.98	0.46
2:H:51:ILE:HG12	2:H:52:ASP:N	2.30	0.46
2:H:135:THR:OG1	2:H:138:SER:O	2.33	0.46
3:L:115:ALA:HB1	3:L:143:PHE:CB	2.46	0.46
1:A:326:TYR:HE2	1:A:330:GLY:O	1.99	0.46
2:H:29:ILE:O	2:H:53:PRO:HG2	2.15	0.46
2:H:124:VAL:HG13	2:H:209:VAL:HG21	1.98	0.46
2:H:151:GLU:O	2:H:151:GLU:HG2	2.16	0.46
2:H:12:VAL:O	2:H:114:VAL:HA	2.15	0.46
2:H:21:SER:HA	2:H:80:TYR:HA	1.97	0.46
2:H:124:VAL:O	2:H:211:LYS:HE3	2.16	0.46
2:H:143:CYS:H	2:H:182:SER:HG	1.61	0.46
1:A:306:PHE:HE2	1:A:335:ILE:HD11	1.80	0.46
2:H:216:ARG:NH2	4:H:269:HOH:O	2.48	0.46
1:A:353:THR:CG2	1:A:354:VAL:N	2.79	0.46
2:H:214:GLU:HA	2:H:215:PRO:HD3	1.52	0.46
1:A:386:GLN:O	1:A:388:LYS:HG3	2.17	0.45
2:H:28:ASN:HD21	2:H:30:LYS:HD2	1.80	0.45
2:H:74:THR:O	2:H:75:SER:C	2.60	0.45
3:L:24:ARG:HG2	3:L:74:ASP:CG	2.41	0.45
3:L:188:ASP:O	3:L:192:ARG:HD3	2.17	0.45
1:A:348:LEU:CD2	1:A:372:PRO:HG3	2.45	0.45
2:H:115:SER:HG	2:H:149:PHE:HE2	1.62	0.45
3:L:12:ALA:CB	3:L:111:LYS:NZ	2.79	0.45
3:L:12:ALA:C	3:L:111:LYS:HZ2	2.23	0.45
2:H:45:LEU:HD21	3:L:48:PRO:CG	2.42	0.45
3:L:123:PRO:HB3	3:L:209:PHE:CE1	2.51	0.45
3:L:191:GLU:C	3:L:193:HIS:N	2.71	0.45
3:L:95:THR:O	3:L:95:THR:CG2	2.59	0.45
3:L:165:ASN:HA	3:L:180:SER:O	2.16	0.45
1:A:331:SER:N	1:A:332:PRO:CD	2.79	0.45
2:H:22:CYS:O	2:H:78:ALA:HB1	2.16	0.45
2:H:84:SER:OG	2:H:85:SER:N	2.49	0.45
3:L:183:LEU:CG	3:L:185:LEU:HD21	2.43	0.45
4:A:396:HOH:O	3:L:58:LEU:CD2	2.64	0.45
3:L:4:LEU:HD22	3:L:102:PHE:O	2.17	0.45
3:L:146:LYS:HD3	4:L:253:HOH:O	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:194:ASN:OD1	3:L:195:SER:HB3	2.15	0.45
2:H:166:VAL:HG22	2:H:184:VAL:HG23	1.99	0.45
3:L:30:VAL:CG2	3:L:72:ARG:HG3	2.46	0.45
2:H:45:LEU:HD22	3:L:102:PHE:HE1	1.81	0.45
2:H:91:THR:HG23	2:H:112:VAL:O	2.16	0.45
2:H:180:LEU:HD23	2:H:181:SER:N	2.32	0.45
1:A:326:TYR:CE2	1:A:330:GLY:O	2.70	0.45
2:H:11:LEU:HD22	2:H:119:THR:HG22	1.99	0.45
2:H:2:VAL:O	2:H:3:GLN:CG	2.65	0.45
2:H:28:ASN:O	2:H:31:ASP:HB2	2.17	0.45
2:H:34:MET:HE3	2:H:79:ALA:HB2	1.98	0.45
3:L:188:ASP:HA	3:L:191:GLU:HG2	1.98	0.45
2:H:11:LEU:CD2	2:H:119:THR:HG22	2.47	0.44
2:H:36:TRP:CE2	2:H:81:LEU:HB2	2.52	0.44
2:H:2:VAL:O	2:H:3:GLN:CB	2.65	0.44
2:H:32:THR:OG1	2:H:33:TYR:N	2.47	0.44
2:H:61:ASP:OD1	2:H:61:ASP:C	2.60	0.44
2:H:126:PRO:CG	2:H:212:LYS:O	2.64	0.44
2:H:128:ALA:HA	2:H:129:PRO:HD3	1.81	0.44
3:L:12:ALA:CB	3:L:111:LYS:CE	2.94	0.44
3:L:19:ALA:HB2	3:L:82:VAL:HG21	1.99	0.44
2:H:30:LYS:O	2:H:31:ASP:C	2.58	0.44
3:L:37:MET:C	3:L:38:HIS:CD2	2.95	0.44
2:H:167:HIS:NE2	3:L:142:ASN:OD1	2.49	0.44
2:H:169:PHE:HA	2:H:170:PRO:HD3	1.87	0.44
3:L:49:LYS:NZ	3:L:49:LYS:CB	2.79	0.44
3:L:87:VAL:HG11	3:L:170:GLN:HB2	1.98	0.44
3:L:151:LYS:NZ	3:L:199:GLU:CD	2.76	0.44
1:A:333:CYS:N	4:A:433:HOH:O	2.51	0.44
1:A:394:LYS:CD	4:A:415:HOH:O	2.61	0.44
1:A:337:PHE:HD1	1:A:380:ILE:CD1	2.31	0.44
2:H:2:VAL:HG13	2:H:27:PHE:CE2	2.53	0.44
2:H:100:TYR:CD1	2:H:100:TYR:C	2.96	0.44
2:H:176:ASP:O	2:H:177:LEU:HD23	2.18	0.44
1:A:312:ILE:HG23	1:A:322:ILE:CG1	2.48	0.44
2:H:108:GLN:O	2:H:109:GLY:O	2.36	0.44
3:L:39:TRP:CG	3:L:77:LEU:HD22	2.53	0.44
1:A:306:PHE:CE2	1:A:335:ILE:HD11	2.53	0.43
2:H:25:SER:OG	2:H:26:GLY:N	2.51	0.43
3:L:49:LYS:HB2	3:L:49:LYS:HZ3	1.81	0.43
1:A:389:LEU:N	1:A:389:LEU:HD23	2.33	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:359:THR:C	1:A:360:GLU:HG3	2.44	0.43
2:H:2:VAL:CA	2:H:25:SER:OG	2.63	0.43
2:H:162:LEU:CG	2:H:163:SER:H	2.30	0.43
3:L:108:LEU:HD12	3:L:109:GLU:H	1.81	0.43
2:H:167:HIS:CE1	3:L:172:SER:HB2	2.54	0.43
2:H:216:ARG:CZ	4:H:269:HOH:O	2.66	0.43
2:H:37:VAL:HG11	2:H:45:LEU:HB3	2.01	0.43
1:A:364:PRO:CD	2:H:31:ASP:OD1	2.67	0.43
2:H:149:PHE:HB3	2:H:150:PRO:CD	2.44	0.43
1:A:337:PHE:HD1	1:A:380:ILE:HD11	1.84	0.42
3:L:12:ALA:O	3:L:111:LYS:NZ	2.46	0.42
1:A:341:ASP:OD1	1:A:343:GLU:O	2.37	0.42
1:A:376:SER:H	1:A:392:PHE:HB3	1.84	0.42
1:A:387:LEU:CD2	3:L:54:ARG:HH22	2.32	0.42
2:H:197:THR:CG2	2:H:198:CYS:N	2.79	0.42
3:L:130:THR:O	3:L:130:THR:HG22	2.18	0.42
1:A:387:LEU:HD21	3:L:54:ARG:NH2	2.32	0.42
3:L:129:LEU:C	3:L:131:SER:H	2.26	0.42
1:A:337:PHE:CD2	1:A:351:LEU:HD13	2.55	0.42
2:H:42:GLU:O	2:H:42:GLU:CG	2.67	0.42
2:H:147:GLY:N	2:H:178:TYR:O	2.47	0.42
3:L:150:VAL:HG12	3:L:151:LYS:H	1.81	0.42
1:A:390:ASN:C	1:A:392:PHE:H	2.28	0.42
1:A:392:PHE:O	1:A:392:PHE:CD1	2.70	0.42
2:H:149:PHE:CD1	2:H:149:PHE:C	2.96	0.42
2:H:167:HIS:O	2:H:183:SER:HB3	2.18	0.42
2:H:124:VAL:HG11	2:H:209:VAL:HG21	2.02	0.42
2:H:169:PHE:CE2	3:L:180:SER:HB2	2.54	0.42
2:H:20:LEU:N	2:H:20:LEU:CD2	2.82	0.42
3:L:28:SER:CA	3:L:73:THR:OG1	2.63	0.42
1:A:387:LEU:N	1:A:387:LEU:HD23	2.34	0.42
2:H:135:THR:OG1	2:H:138:SER:N	2.52	0.42
1:A:376:SER:H	1:A:392:PHE:CB	2.32	0.42
1:A:384:PRO:O	1:A:385:GLY:C	2.61	0.41
2:H:6:GLN:HE22	2:H:110:THR:H	1.68	0.41
2:H:111:LEU:HB3	2:H:152:PRO:HG2	2.02	0.41
3:L:30:VAL:HG23	3:L:72:ARG:CA	2.35	0.41
1:A:337:PHE:CD1	1:A:380:ILE:HD11	2.56	0.41
2:H:4:LEU:HA	2:H:23:THR:O	2.20	0.41
3:L:5:THR:HG23	3:L:24:ARG:CA	2.50	0.41
3:L:90:TYR:N	3:L:90:TYR:CD1	2.88	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:94:GLN:NE2	3:L:99:PRO:O	2.54	0.41
1:A:310:LYS:HB3	1:A:311:GLU:H	1.28	0.41
3:L:22:SER:HA	3:L:76:THR:HA	2.01	0.41
3:L:87:VAL:HG22	3:L:110:ILE:HG12	2.01	0.41
1:A:353:THR:HG22	1:A:356:PRO:CD	2.49	0.41
3:L:172:SER:C	4:L:286:HOH:O	2.62	0.41
2:H:196:ILE:CG2	2:H:197:THR:N	2.83	0.41
2:H:212:LYS:H	2:H:212:LYS:HG3	1.41	0.41
3:L:150:VAL:CG1	3:L:151:LYS:H	2.33	0.41
1:A:353:THR:CG2	1:A:354:VAL:H	2.34	0.41
1:A:375:ASP:HA	1:A:392:PHE:CA	2.50	0.41
2:H:34:MET:CE	2:H:79:ALA:HB2	2.50	0.41
2:H:135:THR:HG1	2:H:138:SER:C	2.28	0.41
2:H:163:SER:C	2:H:165:GLY:N	2.79	0.41
3:L:25:ALA:HB3	3:L:73:THR:HG23	2.02	0.41
2:H:52:ASP:HB3	2:H:55:ASN:ND2	2.36	0.41
2:H:97:ALA:CB	2:H:103:PHE:CD1	2.98	0.41
2:H:99:ASP:HB3	2:H:100:TYR:H	1.57	0.41
3:L:136:VAL:HB	3:L:183:LEU:HB3	2.02	0.41
1:A:334:LYS:O	1:A:336:PRO:HD3	2.21	0.41
1:A:364:PRO:HD3	2:H:31:ASP:OD1	2.21	0.41
2:H:29:ILE:HG13	2:H:34:MET:HE2	1.98	0.41
3:L:6:GLN:NE2	3:L:105:GLY:HA2	2.36	0.41
2:H:197:THR:HG23	2:H:210:ASP:OD1	2.21	0.41
3:L:194:ASN:HA	3:L:211:ARG:CD	2.51	0.41
2:H:148:TYR:O	2:H:178:TYR:HB2	2.21	0.40
2:H:191:TRP:CD1	2:H:196:ILE:CG1	3.00	0.40
3:L:179:MET:CG	3:L:180:SER:N	2.80	0.40
1:A:316:GLN:O	1:A:316:GLN:CG	2.69	0.40
2:H:62:PRO:O	2:H:64:PHE:N	2.55	0.40
2:H:151:GLU:HG3	2:H:178:TYR:CD1	2.57	0.40
2:H:158:ASN:O	2:H:159:SER:HB2	2.20	0.40
3:L:97:VAL:O	3:L:99:PRO:N	2.54	0.40
1:A:375:ASP:HA	1:A:392:PHE:HB2	2.03	0.40
2:H:130:GLY:O	2:H:132:ALA:N	2.54	0.40
3:L:59:GLU:O	3:L:60:SER:C	2.64	0.40
1:A:325:GLN:HE21	1:A:325:GLN:HB3	1.58	0.40
2:H:182:SER:O	2:H:183:SER:HB2	2.22	0.40
1:A:375:ASP:CA	1:A:392:PHE:HB2	2.52	0.40
3:L:58:LEU:N	3:L:58:LEU:CD2	2.48	0.40
3:L:207:LYS:NZ	4:L:278:HOH:O	2.48	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	95/97 (98%)	65 (68%)	19 (20%)	11 (12%)	0	1
2	H	214/216 (99%)	157 (73%)	36 (17%)	21 (10%)	0	2
3	L	194/217 (89%)	142 (73%)	37 (19%)	15 (8%)	1	4
All	All	503/530 (95%)	364 (72%)	92 (18%)	47 (9%)	0	2

All (47) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	H	41	PRO
2	H	45	LEU
2	H	99	ASP
2	H	108	GLN
2	H	191	TRP
3	L	55	ALA
3	L	169	ASP
1	A	299	TYR
1	A	373	PHE
1	A	382	VAL
2	H	3	GLN
2	H	44	GLY
2	H	62	PRO
2	H	63	LYS
2	H	109	GLY
2	H	178	TYR
3	L	98	ASP
3	L	192	ARG
3	L	211	ARG
3	L	212	ASN
1	A	329	ASP

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Mol	Chain	Res	Type
1	A	333	CYS
1	A	354	VAL
1	A	355	ASN
1	A	356	PRO
2	H	9	ALA
2	H	183	SER
3	L	59	GLU
3	L	84	ALA
3	L	162	GLY
1	A	348	LEU
2	H	149	PHE
3	L	14	SER
3	L	18	ARG
1	A	332	PRO
1	A	370	GLU
2	H	163	SER
2	H	170	PRO
2	H	69	THR
2	H	90	ASP
2	H	131	ALA
2	H	102	GLY
2	H	192	PRO
3	L	103	GLY
3	L	144	TYR
3	L	81	PRO
3	L	80	ASN

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	87/87 (100%)	62 (71%)	25 (29%)	0	2
2	H	181/181 (100%)	137 (76%)	44 (24%)	1	4
3	L	179/191 (94%)	135 (75%)	44 (25%)	1	4
All	All	447/459 (97%)	334 (75%)	113 (25%)	0	3

All (113) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	298	SER
1	A	299	TYR
1	A	301	MET
1	A	302	CYS
1	A	303	THR
1	A	308	VAL
1	A	309	VAL
1	A	325	GLN
1	A	329	ASP
1	A	335	ILE
1	A	343	GLU
1	A	348	LEU
1	A	352	ILE
1	A	353	THR
1	A	354	VAL
1	A	355	ASN
1	A	357	ILE
1	A	359	THR
1	A	370	GLU
1	A	371	PRO
1	A	373	PHE
1	A	375	ASP
1	A	380	ILE
1	A	389	LEU
1	A	393	LYS
2	H	1	GLU
2	H	7	SER
2	H	12	VAL
2	H	20	LEU
2	H	23	THR
2	H	35	HIS
2	H	43	GLN
2	H	50	ARG
2	H	55	ASN
2	H	60	TYR
2	H	62	PRO
2	H	74	THR
2	H	77	ASN
2	H	85	SER
2	H	86	LEU
2	H	87	THR
2	H	89	GLU

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Mol	Chain	Res	Type
2	H	90	ASP
2	H	100	TYR
2	H	103	PHE
2	H	108	GLN
2	H	116	SER
2	H	120	THR
2	H	124	VAL
2	H	135	THR
2	H	139	VAL
2	H	140	THR
2	H	141	LEU
2	H	143	CYS
2	H	155	LEU
2	H	156	THR
2	H	158	ASN
2	H	159	SER
2	H	162	LEU
2	H	168	THR
2	H	170	PRO
2	H	175	SER
2	H	179	THR
2	H	189	SER
2	H	194	GLN
2	H	195	THR
2	H	197	THR
2	H	209	VAL
2	H	212	LYS
3	L	4	LEU
3	L	5	THR
3	L	7	SER
3	L	24	ARG
3	L	27	GLU
3	L	38	HIS
3	L	49	LYS
3	L	53	TYR
3	L	58	LEU
3	L	59	GLU
3	L	74	ASP
3	L	76	THR
3	L	77	LEU
3	L	82	VAL
3	L	89	THR

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Mol	Chain	Res	Type
3	L	94	GLN
3	L	96	ASN
3	L	102	PHE
3	L	112	ARG
3	L	118	THR
3	L	125	SER
3	L	135	SER
3	L	139	PHE
3	L	140	LEU
3	L	164	LEU
3	L	165	ASN
3	L	168	THR
3	L	169	ASP
3	L	171	ASP
3	L	176	THR
3	L	177	TYR
3	L	182	THR
3	L	184	THR
3	L	185	LEU
3	L	186	THR
3	L	187	LYS
3	L	192	ARG
3	L	194	ASN
3	L	195	SER
3	L	201	THR
3	L	203	SER
3	L	204	PRO
3	L	206	VAL
3	L	209	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	366	ASN
1	A	390	ASN
2	H	55	ASN
2	H	108	GLN
2	H	158	ASN
3	L	41	GLN
3	L	46	GLN
3	L	96	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

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

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

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	97/97 (100%)	-1.46	0 100 100	7, 43, 60, 66	0
2	H	216/216 (100%)	-1.51	0 100 100	10, 33, 58, 66	0
3	L	204/217 (94%)	-1.43	0 100 100	18, 43, 56, 65	0
All	All	517/530 (97%)	-1.47	0 100 100	7, 40, 59, 66	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.