



Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 11:37 am BST

PDB ID : 3KHX
Title : Crystal structure of Staphylococcus aureus metallopeptidase (Sapep/DapE) in the apo-form
Authors : Girish, T.S.; Gopal, B.
Deposited on : 2009-10-31
Resolution : 2.30 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.11
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

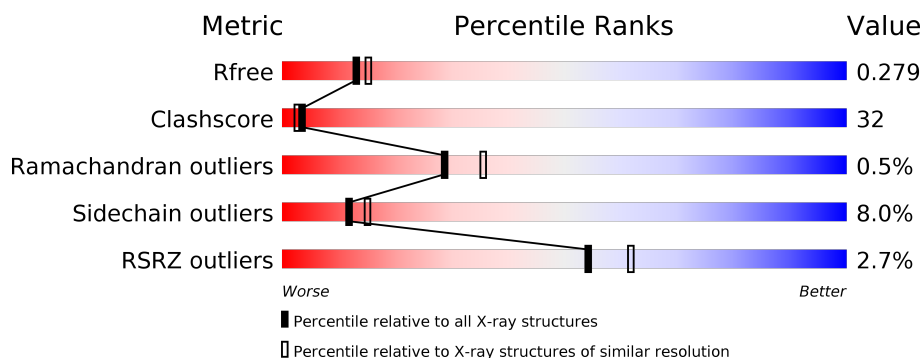
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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 on 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	492	<div> <div>3%</div> <div>54%</div> <div>27%</div> <div>•</div> <div>14%</div> </div>
1	B	492	<div> <div>2%</div> <div>58%</div> <div>23%</div> <div>5%</div> <div>14%</div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6716 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 Putative dipeptidase SACOL1801.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	422	Total	C	N	O	S	1	0	0
			3232	2052	531	635	14			
1	B	425	Total	C	N	O	S	0	0	0
			3302	2099	546	642	15			

There are 46 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-22	MET	-	EXPRESSION TAG	UNP Q5HF23
A	-21	GLY	-	EXPRESSION TAG	UNP Q5HF23
A	-20	SER	-	EXPRESSION TAG	UNP Q5HF23
A	-19	SER	-	EXPRESSION TAG	UNP Q5HF23
A	-18	HIS	-	EXPRESSION TAG	UNP Q5HF23
A	-17	HIS	-	EXPRESSION TAG	UNP Q5HF23
A	-16	HIS	-	EXPRESSION TAG	UNP Q5HF23
A	-15	HIS	-	EXPRESSION TAG	UNP Q5HF23
A	-14	HIS	-	EXPRESSION TAG	UNP Q5HF23
A	-13	HIS	-	EXPRESSION TAG	UNP Q5HF23
A	-12	SER	-	EXPRESSION TAG	UNP Q5HF23
A	-11	SER	-	EXPRESSION TAG	UNP Q5HF23
A	-10	GLY	-	EXPRESSION TAG	UNP Q5HF23
A	-9	LEU	-	EXPRESSION TAG	UNP Q5HF23
A	-8	VAL	-	EXPRESSION TAG	UNP Q5HF23
A	-7	PRO	-	EXPRESSION TAG	UNP Q5HF23
A	-6	ARG	-	EXPRESSION TAG	UNP Q5HF23
A	-5	GLY	-	EXPRESSION TAG	UNP Q5HF23
A	-4	SER	-	EXPRESSION TAG	UNP Q5HF23
A	-3	HIS	-	EXPRESSION TAG	UNP Q5HF23
A	-2	MET	-	EXPRESSION TAG	UNP Q5HF23
A	-1	ALA	-	EXPRESSION TAG	UNP Q5HF23
A	0	SER	-	EXPRESSION TAG	UNP Q5HF23
B	-22	MET	-	EXPRESSION TAG	UNP Q5HF23
B	-21	GLY	-	EXPRESSION TAG	UNP Q5HF23

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-20	SER	-	EXPRESSION TAG	UNP Q5HF23
B	-19	SER	-	EXPRESSION TAG	UNP Q5HF23
B	-18	HIS	-	EXPRESSION TAG	UNP Q5HF23
B	-17	HIS	-	EXPRESSION TAG	UNP Q5HF23
B	-16	HIS	-	EXPRESSION TAG	UNP Q5HF23
B	-15	HIS	-	EXPRESSION TAG	UNP Q5HF23
B	-14	HIS	-	EXPRESSION TAG	UNP Q5HF23
B	-13	HIS	-	EXPRESSION TAG	UNP Q5HF23
B	-12	SER	-	EXPRESSION TAG	UNP Q5HF23
B	-11	SER	-	EXPRESSION TAG	UNP Q5HF23
B	-10	GLY	-	EXPRESSION TAG	UNP Q5HF23
B	-9	LEU	-	EXPRESSION TAG	UNP Q5HF23
B	-8	VAL	-	EXPRESSION TAG	UNP Q5HF23
B	-7	PRO	-	EXPRESSION TAG	UNP Q5HF23
B	-6	ARG	-	EXPRESSION TAG	UNP Q5HF23
B	-5	GLY	-	EXPRESSION TAG	UNP Q5HF23
B	-4	SER	-	EXPRESSION TAG	UNP Q5HF23
B	-3	HIS	-	EXPRESSION TAG	UNP Q5HF23
B	-2	MET	-	EXPRESSION TAG	UNP Q5HF23
B	-1	ALA	-	EXPRESSION TAG	UNP Q5HF23
B	0	SER	-	EXPRESSION TAG	UNP Q5HF23

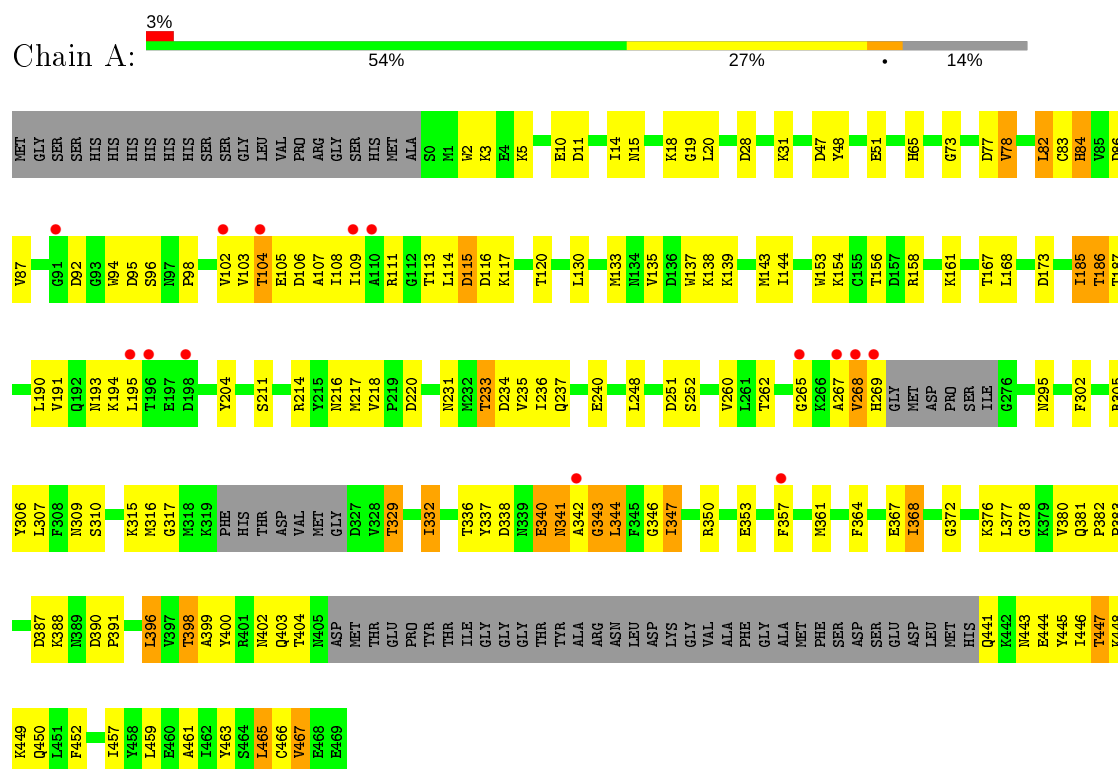
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	92	Total O 92 92	0	0
2	B	90	Total O 90 90	0	0

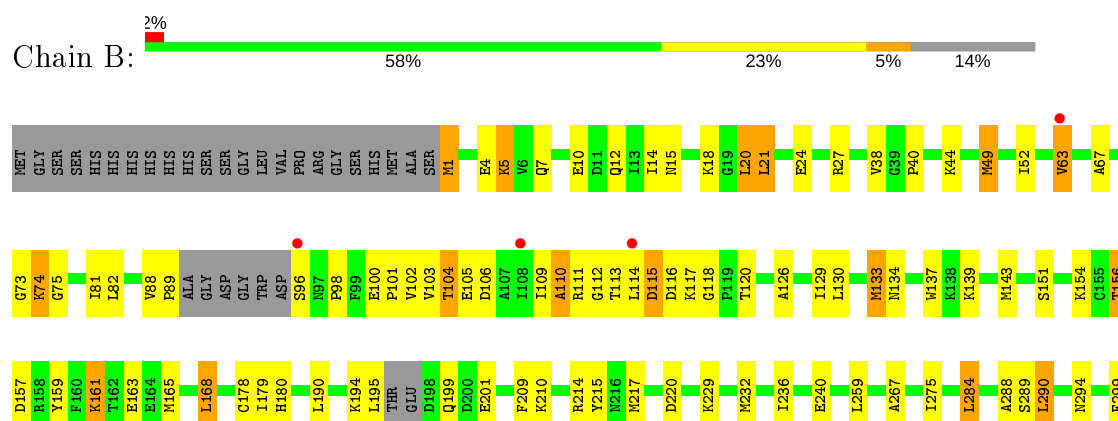
3 Residue-property plots [i](#)

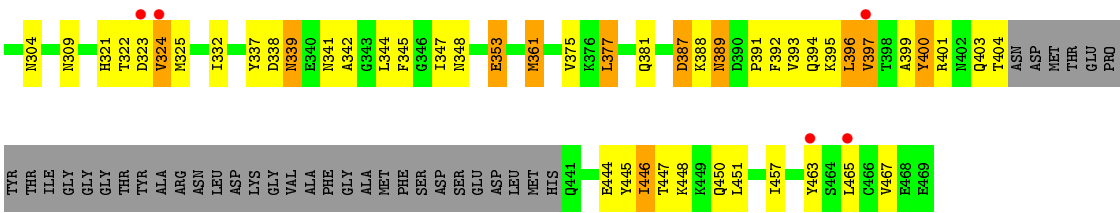
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are 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: Putative dipeptidase SACOL1801



• Molecule 1: Putative dipeptidase SACOL1801





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	64.82Å 134.17Å 68.28Å 90.00° 94.52° 90.00°	Depositor
Resolution (Å)	60.71 – 2.30 60.70 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.3 (60.71-2.30) 99.3 (60.70-2.30)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.27 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.5.0066	Depositor
R, R_{free}	0.246 , 0.278 0.256 , 0.279	Depositor DCC
R_{free} test set	2605 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	44.5	Xtriage
Anisotropy	0.134	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 67.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.018 for l,-k,h	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	6716	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.62% 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 [i](#)

5.1 Standard geometry [i](#)

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	0.56	0/3298	0.69	2/4478 (0.0%)
1	B	0.49	0/3372	0.64	0/4574
All	All	0.53	0/6670	0.66	2/9052 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	96	SER	CB-CA-C	5.84	121.19	110.10
1	A	211	SER	CB-CA-C	5.60	120.74	110.10

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	3232	0	3017	208	0
1	B	3302	0	3113	197	0
2	A	92	0	0	8	0
2	B	90	0	0	4	0
All	All	6716	0	6130	405	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 32.

All (405) close contacts within the same asymmetric unit are listed below, sorted by their clash

magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:185:ILE:HB	1:A:350:ARG:NH1	1.28	1.43
1:A:130:LEU:HD21	1:A:459:LEU:CD1	1.55	1.34
1:A:186:THR:CG2	1:A:380:VAL:HG22	1.60	1.32
1:B:5:LYS:N	1:B:5:LYS:HE3	1.50	1.27
1:A:114:LEU:HD11	1:A:445:TYR:O	1.35	1.25
1:B:74:LYS:HD3	1:B:75:GLY:N	1.51	1.23
1:A:130:LEU:CD2	1:A:459:LEU:HD11	1.68	1.22
1:B:12:GLN:HG3	2:B:529:HOH:O	1.35	1.22
1:A:104:THR:CG2	1:A:107:ALA:H	1.57	1.18
1:A:28:ASP:OD2	1:A:31:LYS:HG3	1.44	1.17
1:A:447:THR:CG2	1:A:449:LYS:H	1.61	1.14
1:A:186:THR:HG21	1:A:380:VAL:HG22	1.20	1.13
1:B:113:THR:HB	1:B:446:ILE:HG22	1.30	1.12
1:A:185:ILE:CB	1:A:350:ARG:HH11	1.64	1.10
1:A:104:THR:HG22	1:A:107:ALA:N	1.66	1.10
1:B:49:MET:HE1	1:B:52:ILE:HD11	1.15	1.07
1:A:310:SER:HB3	2:A:533:HOH:O	1.54	1.07
1:B:392:PHE:CZ	1:B:396:LEU:HD11	1.88	1.07
1:A:130:LEU:HD21	1:A:459:LEU:HD12	1.38	1.05
1:A:130:LEU:CD2	1:A:459:LEU:CD1	2.30	1.05
1:A:377:LEU:HD23	1:A:378:GLY:O	1.57	1.04
1:B:392:PHE:CE2	1:B:396:LEU:CD1	2.41	1.03
1:A:186:THR:CG2	1:A:380:VAL:CG2	2.37	1.03
1:A:387:ASP:OD1	1:A:388:LYS:N	1.92	1.02
1:B:400:TYR:HD2	1:B:401:ARG:N	1.57	1.01
1:A:130:LEU:HD21	1:A:459:LEU:HD11	1.18	1.01
1:A:186:THR:HG23	1:A:380:VAL:HG22	1.41	1.01
1:A:447:THR:HB	1:A:450:GLN:HG3	1.43	1.00
1:A:186:THR:HG21	1:A:380:VAL:CG2	1.91	1.00
1:B:104:THR:CG2	1:B:105:GLU:N	2.21	1.00
1:A:104:THR:HG23	1:A:106:ASP:H	1.24	0.99
1:A:447:THR:HG22	1:A:449:LYS:N	1.78	0.99
1:A:231:ASN:HD21	1:A:234:ASP:CB	1.75	0.98
1:A:83:CYS:HB3	2:A:510:HOH:O	1.63	0.98
1:A:231:ASN:HD21	1:A:234:ASP:HB2	1.24	0.98
1:A:104:THR:HG22	1:A:107:ALA:H	0.82	0.97
1:B:49:MET:CE	1:B:52:ILE:HD11	1.94	0.96
1:A:102:VAL:HG22	1:A:109:ILE:HB	1.48	0.96
1:A:191:VAL:HG21	1:A:376:LYS:HE2	1.46	0.96
1:A:377:LEU:CD2	1:A:378:GLY:O	2.14	0.95
1:B:156:THR:O	1:B:161:LYS:HE2	1.67	0.95

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:104:THR:HG22	1:B:106:ASP:N	1.80	0.95
1:A:329:THR:HG22	1:A:350:ARG:HB2	1.49	0.95
1:A:2:TRP:HZ3	1:A:5:LYS:HZ1	1.09	0.95
1:A:185:ILE:CB	1:A:350:ARG:NH1	2.25	0.94
1:B:49:MET:HE1	1:B:52:ILE:CD1	1.96	0.94
1:A:447:THR:HG22	1:A:450:GLN:H	1.31	0.94
1:B:400:TYR:CD2	1:B:401:ARG:N	2.35	0.93
1:A:447:THR:HG23	1:A:449:LYS:H	1.30	0.93
1:A:231:ASN:ND2	1:A:234:ASP:HB2	1.83	0.93
1:B:88:VAL:HG13	1:B:89:PRO:HD2	1.48	0.93
1:B:104:THR:HG23	1:B:105:GLU:H	1.34	0.93
1:A:104:THR:CG2	1:A:107:ALA:N	2.27	0.92
1:B:104:THR:HG22	1:B:106:ASP:H	1.32	0.92
1:A:2:TRP:CZ3	1:A:5:LYS:NZ	2.38	0.92
1:B:114:LEU:HD11	1:B:445:TYR:O	1.70	0.91
1:B:114:LEU:HD12	1:B:114:LEU:N	1.84	0.91
1:A:447:THR:HG22	1:A:449:LYS:H	1.34	0.91
1:B:5:LYS:H	1:B:5:LYS:HE3	1.23	0.90
1:B:49:MET:HA	1:B:49:MET:CE	2.00	0.90
1:A:447:THR:CG2	1:A:449:LYS:N	2.33	0.90
1:B:74:LYS:HD3	1:B:75:GLY:H	1.37	0.89
1:B:392:PHE:CE2	1:B:396:LEU:HD11	2.06	0.89
1:B:338:ASP:O	1:B:342:ALA:HB1	1.73	0.89
1:B:395:LYS:HD2	1:B:465:LEU:HD23	1.54	0.89
1:A:2:TRP:HZ3	1:A:5:LYS:NZ	1.71	0.88
1:B:209:PHE:CE1	1:B:284:LEU:HD13	2.09	0.88
1:A:204:TYR:CE2	1:A:235:VAL:HG21	2.10	0.87
1:B:104:THR:HG23	1:B:105:GLU:N	1.88	0.87
1:A:104:THR:HG23	1:A:106:ASP:N	1.91	0.86
1:A:231:ASN:HD21	1:A:234:ASP:CG	1.79	0.86
1:A:104:THR:HG23	1:A:105:GLU:N	1.91	0.85
1:B:236:ILE:O	1:B:240:GLU:HG2	1.77	0.84
1:B:400:TYR:CD2	1:B:400:TYR:C	2.51	0.84
1:A:84:HIS:HD2	1:A:86:ASP:OD1	1.60	0.84
1:B:103:VAL:HG13	1:B:103:VAL:O	1.79	0.83
1:B:21:LEU:HD21	1:B:113:THR:CG2	2.08	0.83
1:B:1:MET:O	1:B:4:GLU:CB	2.27	0.83
1:B:113:THR:HB	1:B:446:ILE:CG2	2.09	0.82
1:A:20:LEU:HG	1:A:117:LYS:HD3	1.60	0.82
1:B:63:VAL:CG2	1:B:67:ALA:O	2.27	0.82
1:B:199:GLN:HE22	1:B:294:ASN:H	1.23	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:321:HIS:CE1	1:B:323:ASP:CB	2.62	0.82
1:B:104:THR:HG22	1:B:105:GLU:N	1.95	0.81
1:A:396:LEU:HD22	1:A:461:ALA:CB	2.09	0.81
1:A:84:HIS:CD2	1:A:86:ASP:OD1	2.33	0.81
1:B:154:LYS:HB3	1:B:156:THR:HG22	1.61	0.81
1:A:186:THR:HG23	1:A:380:VAL:HA	1.62	0.81
1:B:63:VAL:HG23	1:B:67:ALA:O	1.80	0.81
1:B:104:THR:CG2	1:B:106:ASP:H	1.95	0.80
1:A:357:PHE:CE1	1:A:361:MET:HE3	2.16	0.80
1:B:290:LEU:N	1:B:290:LEU:HD12	1.94	0.80
1:B:74:LYS:CD	1:B:75:GLY:N	2.41	0.80
1:A:130:LEU:HD23	1:A:459:LEU:HD11	1.62	0.80
1:A:104:THR:CG2	1:A:106:ASP:H	1.95	0.80
1:A:329:THR:CG2	1:A:350:ARG:HB2	2.11	0.80
1:B:49:MET:HA	1:B:49:MET:HE2	1.64	0.79
1:B:156:THR:C	1:B:161:LYS:HE2	2.03	0.79
1:A:168:LEU:HD11	1:A:465:LEU:HD21	1.63	0.79
1:A:5:LYS:HE3	1:A:459:LEU:HD21	1.65	0.78
1:B:321:HIS:CE1	1:B:323:ASP:HB2	2.18	0.78
1:A:251:ASP:OD1	1:A:262:THR:CG2	2.31	0.78
1:A:316:MET:HG2	1:A:364:PHE:CE1	2.18	0.77
1:A:114:LEU:HD13	1:A:444:GLU:HG2	1.66	0.77
1:B:395:LYS:CD	1:B:465:LEU:HD23	2.14	0.77
1:B:5:LYS:N	1:B:5:LYS:CE	2.41	0.77
1:B:392:PHE:CZ	1:B:396:LEU:CD1	2.66	0.77
1:B:21:LEU:HD21	1:B:113:THR:HG22	1.65	0.77
1:B:74:LYS:HD3	1:B:74:LYS:C	2.00	0.77
1:A:204:TYR:OH	1:A:235:VAL:CG2	2.33	0.76
1:A:251:ASP:O	1:A:262:THR:HG22	1.85	0.76
1:B:139:LYS:HD2	1:B:465:LEU:O	1.85	0.75
1:A:447:THR:HG22	1:A:450:GLN:N	2.01	0.75
1:A:168:LEU:HD11	1:A:465:LEU:CD2	2.17	0.74
1:B:73:GLY:HA2	1:B:137:TRP:CG	2.22	0.74
1:B:323:ASP:N	1:B:324:VAL:CB	2.50	0.74
1:B:392:PHE:HA	1:B:465:LEU:HD21	1.68	0.74
1:A:396:LEU:CD2	1:A:461:ALA:CB	2.66	0.74
1:A:343:GLY:HA2	2:A:477:HOH:O	1.88	0.74
1:A:310:SER:OG	1:A:315:LYS:HB2	1.88	0.74
1:A:317:GLY:HA3	2:A:493:HOH:O	1.86	0.74
1:B:338:ASP:O	1:B:342:ALA:CB	2.36	0.74
1:A:448:LYS:O	1:A:452:PHE:CD2	2.42	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:114:LEU:HD11	1:B:445:TYR:C	2.07	0.73
1:B:114:LEU:CD1	1:B:114:LEU:N	2.52	0.72
1:A:28:ASP:CG	1:A:31:LYS:HG3	2.10	0.72
1:B:190:LEU:CD1	1:B:347:ILE:HD11	2.20	0.72
1:B:321:HIS:CE1	1:B:323:ASP:HB3	2.25	0.72
1:A:347:ILE:N	1:A:347:ILE:CD1	2.51	0.71
1:B:15:ASN:O	1:B:18:LYS:HG2	1.90	0.71
1:B:49:MET:HA	1:B:49:MET:HE3	1.72	0.71
1:A:251:ASP:OD1	1:A:262:THR:HG21	1.90	0.71
1:B:395:LYS:CE	1:B:465:LEU:HD23	2.19	0.71
1:B:290:LEU:CD1	1:B:290:LEU:N	2.52	0.71
1:B:126:ALA:O	1:B:130:LEU:HG	1.91	0.70
1:B:102:VAL:HG22	1:B:109:ILE:CB	2.22	0.70
1:B:74:LYS:HE2	1:B:74:LYS:HA	1.71	0.70
1:A:186:THR:HG23	1:A:380:VAL:CG2	2.13	0.70
1:B:49:MET:CE	1:B:52:ILE:CD1	2.61	0.70
1:B:403:GLN:O	1:B:404:THR:OG1	2.10	0.69
1:B:387:ASP:OD2	1:B:387:ASP:N	2.22	0.69
1:B:391:PRO:O	1:B:395:LYS:HE2	1.93	0.69
1:B:209:PHE:CE1	1:B:284:LEU:CD1	2.75	0.69
1:B:74:LYS:CD	1:B:75:GLY:H	2.02	0.69
1:B:1:MET:O	1:B:4:GLU:N	2.25	0.69
1:A:138:LYS:HG2	1:A:466:CYS:O	1.93	0.69
1:A:204:TYR:CZ	1:A:235:VAL:HG21	2.27	0.68
1:A:231:ASN:ND2	1:A:234:ASP:OD2	2.27	0.68
1:B:102:VAL:CG2	1:B:109:ILE:CB	2.71	0.68
1:B:322:THR:C	1:B:324:VAL:CB	2.62	0.68
1:A:194:LYS:HE3	1:A:342:ALA:HB2	1.76	0.68
1:A:347:ILE:HD13	1:A:347:ILE:N	2.09	0.67
1:B:323:ASP:HA	1:B:324:VAL:C	2.14	0.67
1:A:316:MET:HG2	1:A:364:PHE:CZ	2.29	0.67
1:B:214:ARG:NH1	1:B:217:MET:HG3	2.09	0.67
1:B:275:ILE:HG22	1:B:275:ILE:O	1.94	0.67
1:A:233:THR:O	1:A:237:GLN:HG3	1.95	0.66
1:B:5:LYS:CA	1:B:5:LYS:HE3	2.26	0.66
1:B:103:VAL:CG1	1:B:103:VAL:O	2.43	0.66
1:B:109:ILE:O	1:B:110:ALA:HB2	1.95	0.65
1:B:63:VAL:HG22	1:B:67:ALA:O	1.96	0.65
1:B:27:ARG:HG3	1:B:38:VAL:HB	1.77	0.65
1:A:3:LYS:NZ	2:A:527:HOH:O	2.28	0.65
1:B:209:PHE:CZ	1:B:284:LEU:HD13	2.32	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:392:PHE:CE2	1:B:396:LEU:HD13	2.31	0.64
1:B:7:GLN:O	1:B:10:GLU:HG3	1.97	0.64
1:A:185:ILE:HB	1:A:350:ARG:HH12	1.49	0.64
1:A:396:LEU:CD2	1:A:461:ALA:HB3	2.26	0.64
1:A:447:THR:CG2	1:A:450:GLN:H	2.07	0.64
1:B:113:THR:O	1:B:118:GLY:HA3	1.98	0.64
1:A:377:LEU:HG	1:A:378:GLY:N	2.13	0.64
1:B:209:PHE:O	1:B:210:LYS:HD3	1.98	0.64
1:A:104:THR:CG2	1:A:106:ASP:N	2.57	0.63
1:A:120:THR:HG23	1:A:143:MET:CE	2.29	0.63
1:A:47:ASP:O	1:A:51:GLU:HG3	1.99	0.62
1:B:154:LYS:CB	1:B:156:THR:CG2	2.78	0.62
1:B:381:GLN:NE2	2:B:496:HOH:O	2.24	0.62
1:A:104:THR:CG2	1:A:105:GLU:N	2.63	0.62
1:B:400:TYR:HA	1:B:457:ILE:HD13	1.82	0.62
1:A:220:ASP:HB2	1:A:265:GLY:O	2.00	0.62
1:A:377:LEU:HD21	1:A:378:GLY:O	2.00	0.62
1:B:20:LEU:HD13	1:B:117:LYS:HB3	1.82	0.62
1:B:5:LYS:H	1:B:5:LYS:CE	2.07	0.62
1:B:154:LYS:CB	1:B:156:THR:HG22	2.29	0.61
1:B:154:LYS:HB2	1:B:156:THR:HG23	1.82	0.61
1:A:194:LYS:HE2	1:A:295:ASN:OD1	2.01	0.61
1:A:5:LYS:HD2	1:A:133:MET:HE3	1.82	0.61
1:A:114:LEU:CD1	1:A:445:TYR:O	2.29	0.61
1:A:396:LEU:HD22	1:A:461:ALA:HB2	1.82	0.61
1:A:217:MET:HA	1:A:268:VAL:O	2.01	0.61
1:B:154:LYS:HB3	1:B:156:THR:CG2	2.31	0.61
1:A:447:THR:CB	1:A:450:GLN:HG3	2.26	0.60
1:A:193:ASN:ND2	1:A:372:GLY:HA2	2.17	0.60
1:B:5:LYS:CA	1:B:5:LYS:CE	2.80	0.60
1:A:252:SER:HA	1:A:260:VAL:O	2.02	0.59
1:B:394:GLN:O	1:B:397:VAL:HG13	2.03	0.59
1:B:392:PHE:HA	1:B:465:LEU:CD2	2.31	0.59
1:B:88:VAL:CG1	1:B:89:PRO:HD2	2.30	0.59
1:A:332:ILE:HD12	1:A:332:ILE:N	2.17	0.59
1:A:78:VAL:H	1:A:167:THR:HG22	1.67	0.59
1:B:289:SER:C	1:B:290:LEU:HD12	2.23	0.58
1:B:341:ASN:CB	1:B:342:ALA:HA	2.33	0.58
1:A:204:TYR:OH	1:A:235:VAL:HG21	2.03	0.58
1:B:18:LYS:NZ	2:B:483:HOH:O	2.32	0.58
1:A:216:ASN:O	1:A:269:HIS:HA	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:10:GLU:HG2	1:A:452:PHE:CE1	2.39	0.57
1:A:113:THR:HB	1:A:446:ILE:HG22	1.85	0.57
1:B:88:VAL:HG13	1:B:89:PRO:CD	2.29	0.57
1:A:14:ILE:O	1:A:18:LYS:HG3	2.04	0.57
1:A:329:THR:HG22	1:A:350:ARG:CB	2.30	0.57
1:A:377:LEU:HD21	1:A:380:VAL:HG23	1.86	0.57
1:A:133:MET:HG3	1:A:135:VAL:HG13	1.85	0.56
1:A:185:ILE:HB	1:A:350:ARG:HH11	0.75	0.56
1:B:321:HIS:NE2	1:B:323:ASP:HB3	2.21	0.56
1:A:154:LYS:HB3	1:A:156:THR:HG22	1.87	0.56
1:B:74:LYS:CE	1:B:75:GLY:H	2.17	0.56
1:A:353:GLU:OE1	1:A:383:PRO:HB2	2.06	0.56
1:A:463:TYR:O	1:A:467:VAL:HG23	2.06	0.56
1:B:10:GLU:O	1:B:14:ILE:HG12	2.05	0.56
1:A:336:THR:O	1:A:343:GLY:HA2	2.06	0.55
1:A:193:ASN:HD22	1:A:372:GLY:HA2	1.70	0.55
1:A:396:LEU:HD22	1:A:461:ALA:HB3	1.84	0.55
1:B:165:MET:HE3	1:B:353:GLU:HG3	1.88	0.55
1:A:367:GLU:CG	1:A:367:GLU:O	2.55	0.55
1:A:83:CYS:CB	2:A:510:HOH:O	2.36	0.55
1:B:165:MET:CE	1:B:353:GLU:HG3	2.37	0.55
1:B:288:ALA:HB2	1:B:304:ASN:HD21	1.72	0.54
1:B:395:LYS:HD2	1:B:465:LEU:CD2	2.34	0.54
1:A:447:THR:HG22	1:A:449:LYS:CA	2.37	0.54
1:A:338:ASP:OD2	1:A:340:GLU:N	2.40	0.54
1:B:154:LYS:HB2	1:B:156:THR:CG2	2.37	0.54
1:B:322:THR:O	1:B:325:MET:N	2.40	0.54
1:A:14:ILE:CG2	1:A:15:ASN:N	2.71	0.54
1:B:1:MET:C	1:B:4:GLU:H	2.11	0.54
1:A:111:ARG:NH1	1:A:441:GLN:CB	2.71	0.54
1:A:194:LYS:HE2	1:A:295:ASN:CG	2.28	0.54
1:A:95:ASP:H	1:A:443:ASN:HD21	1.55	0.53
1:A:14:ILE:HD11	1:A:108:ILE:HG12	1.90	0.53
1:B:129:ILE:O	1:B:133:MET:HG2	2.08	0.53
1:A:204:TYR:HE2	1:A:235:VAL:HG21	1.68	0.53
1:A:65:HIS:HD2	2:A:518:HOH:O	1.90	0.53
1:A:357:PHE:CZ	1:A:361:MET:HE3	2.44	0.53
1:B:109:ILE:O	1:B:110:ALA:CB	2.58	0.52
1:B:190:LEU:HD11	1:B:347:ILE:CD1	2.39	0.52
1:A:104:THR:HG21	1:A:107:ALA:N	2.17	0.52
1:A:14:ILE:HD11	1:A:108:ILE:CG1	2.40	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:377:LEU:HG	1:A:378:GLY:H	1.73	0.52
1:A:102:VAL:CG2	1:A:109:ILE:HB	2.30	0.52
1:A:82:LEU:N	1:A:82:LEU:HD23	2.24	0.52
1:B:168:LEU:H	1:B:168:LEU:HD23	1.75	0.52
1:B:361:MET:HG2	1:B:377:LEU:HD11	1.92	0.52
1:A:357:PHE:CE1	1:A:361:MET:CE	2.93	0.52
1:A:111:ARG:HH12	1:A:441:GLN:CB	2.23	0.52
1:A:448:LYS:O	1:A:452:PHE:HD2	1.93	0.52
1:A:447:THR:CG2	1:A:449:LYS:CB	2.88	0.52
1:B:105:GLU:O	1:B:448:LYS:NZ	2.32	0.52
1:A:11:ASP:O	1:A:14:ILE:HG22	2.10	0.52
1:A:344:LEU:HD12	1:A:344:LEU:C	2.29	0.52
1:B:113:THR:C	1:B:114:LEU:HD12	2.29	0.52
1:B:82:LEU:N	1:B:82:LEU:HD12	2.25	0.52
1:A:332:ILE:N	1:A:332:ILE:CD1	2.73	0.51
1:A:77:ASP:HB2	1:A:139:LYS:HG2	1.91	0.51
1:B:404:THR:HG22	1:B:404:THR:O	2.09	0.51
1:A:302:PHE:HE1	1:A:368:ILE:HG13	1.75	0.51
1:A:396:LEU:HD13	1:A:457:ILE:HG22	1.92	0.51
1:B:1:MET:O	1:B:4:GLU:CA	2.58	0.51
1:A:28:ASP:OD2	1:A:31:LYS:CG	2.37	0.51
1:A:135:VAL:HB	1:A:467:VAL:HG11	1.93	0.51
1:B:21:LEU:HD21	1:B:113:THR:HG23	1.90	0.51
1:B:463:TYR:CD1	1:B:467:VAL:HG21	2.45	0.51
1:A:102:VAL:CG2	1:A:102:VAL:O	2.59	0.51
1:A:2:TRP:CE3	1:A:5:LYS:NZ	2.69	0.51
1:A:186:THR:HG23	1:A:380:VAL:CA	2.37	0.51
1:B:190:LEU:CD1	1:B:347:ILE:CD1	2.88	0.51
1:A:114:LEU:HD12	1:A:114:LEU:N	2.25	0.50
1:B:168:LEU:HD12	1:B:391:PRO:HG2	1.94	0.50
1:B:168:LEU:N	1:B:168:LEU:HD23	2.27	0.50
1:A:338:ASP:OD2	1:A:340:GLU:HB3	2.12	0.50
1:A:73:GLY:HA2	1:A:137:TRP:CG	2.46	0.50
1:B:115:ASP:N	1:B:115:ASP:OD2	2.44	0.50
1:B:392:PHE:CA	1:B:465:LEU:HD21	2.37	0.50
1:A:399:ALA:O	1:A:403:GLN:HG3	2.11	0.50
1:A:78:VAL:O	1:A:167:THR:HG22	2.12	0.50
1:B:82:LEU:CD1	1:B:82:LEU:N	2.74	0.50
1:A:402:ASN:OD1	1:A:402:ASN:O	2.30	0.50
1:A:367:GLU:O	1:A:367:GLU:HG3	2.11	0.50
1:B:102:VAL:HG23	1:B:102:VAL:O	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:201:GLU:CD	1:B:229:LYS:HD3	2.32	0.49
1:A:102:VAL:CG2	1:A:109:ILE:HD12	2.42	0.49
1:B:275:ILE:CG2	1:B:275:ILE:O	2.60	0.49
1:B:446:ILE:HD11	1:B:451:LEU:HB2	1.93	0.49
1:B:446:ILE:HG12	1:B:447:THR:O	2.13	0.49
1:A:310:SER:OG	1:A:315:LYS:CB	2.60	0.49
1:A:194:LYS:CE	1:A:342:ALA:HB2	2.41	0.49
1:B:113:THR:CB	1:B:446:ILE:HG22	2.21	0.49
1:A:251:ASP:H	1:A:262:THR:HG23	1.76	0.48
1:A:28:ASP:OD2	1:A:31:LYS:HE3	2.13	0.48
1:B:159:TYR:CE2	1:B:163:GLU:HB2	2.48	0.48
1:B:168:LEU:CD1	1:B:391:PRO:HG2	2.43	0.48
1:B:112:GLY:H	1:B:444:GLU:CD	2.17	0.48
1:B:21:LEU:HA	1:B:21:LEU:HD12	1.73	0.48
1:B:339:ASN:O	2:B:530:HOH:O	2.20	0.48
1:B:190:LEU:HD12	1:B:347:ILE:HD11	1.92	0.48
1:A:158:ARG:O	1:A:161:LYS:HB2	2.14	0.48
1:A:337:TYR:HA	1:A:343:GLY:CA	2.42	0.48
1:B:114:LEU:CD1	1:B:445:TYR:O	2.52	0.48
1:B:232:MET:HE2	1:B:259:LEU:HB2	1.95	0.48
1:A:103:VAL:HG12	1:A:103:VAL:O	2.13	0.48
1:B:111:ARG:HD3	1:B:444:GLU:OE2	2.14	0.48
1:A:218:VAL:HB	1:A:267:ALA:HA	1.95	0.48
1:A:447:THR:CG2	1:A:449:LYS:CA	2.91	0.48
1:B:323:ASP:CA	1:B:324:VAL:CB	2.91	0.48
1:B:375:VAL:HG12	1:B:377:LEU:HD13	1.96	0.48
1:A:10:GLU:HG2	1:A:452:PHE:CZ	2.49	0.47
1:B:321:HIS:NE2	1:B:323:ASP:CB	2.77	0.47
1:B:400:TYR:HD2	1:B:401:ARG:CA	2.24	0.47
1:B:447:THR:OG1	1:B:450:GLN:HB3	2.15	0.47
1:A:305:ARG:HD3	1:A:306:TYR:CZ	2.50	0.47
1:A:168:LEU:CD1	1:A:465:LEU:HD21	2.40	0.47
1:B:190:LEU:HD11	1:B:347:ILE:HD11	1.93	0.46
1:B:88:VAL:CG1	1:B:89:PRO:CD	2.93	0.46
1:A:114:LEU:HD11	1:A:445:TYR:C	2.23	0.46
1:A:341:ASN:O	1:A:342:ALA:HB3	2.15	0.46
1:A:204:TYR:CZ	1:A:235:VAL:CG2	2.95	0.46
1:A:195:LEU:H	1:A:295:ASN:HD21	1.62	0.46
1:B:10:GLU:O	1:B:14:ILE:CG1	2.63	0.46
1:B:194:LYS:O	1:B:195:LEU:C	2.54	0.46
1:B:157:ASP:O	1:B:161:LYS:HD3	2.14	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:346:GLY:C	1:A:347:ILE:HD12	2.35	0.46
1:B:446:ILE:CD1	1:B:451:LEU:HB2	2.45	0.46
1:A:346:GLY:C	1:A:347:ILE:CD1	2.84	0.45
1:B:393:VAL:O	1:B:397:VAL:HG12	2.17	0.45
1:B:332:ILE:N	1:B:332:ILE:HD13	2.32	0.45
1:B:400:TYR:CD2	1:B:401:ARG:CA	2.99	0.45
1:B:88:VAL:CG1	1:B:89:PRO:N	2.80	0.45
1:A:336:THR:O	1:A:343:GLY:CA	2.64	0.45
1:B:399:ALA:HB3	1:B:457:ILE:HG23	1.98	0.45
1:A:114:LEU:HD13	1:A:444:GLU:CG	2.41	0.45
1:B:63:VAL:HG21	1:B:159:TYR:HD1	1.82	0.45
1:A:102:VAL:HG23	1:A:102:VAL:O	2.16	0.45
1:A:185:ILE:N	1:A:381:GLN:O	2.42	0.45
1:A:94:TRP:CD2	1:A:98:PRO:HG3	2.53	0.44
1:B:101:PRO:HA	1:B:109:ILE:O	2.18	0.44
1:B:337:TYR:C	1:B:337:TYR:CD2	2.90	0.44
1:A:186:THR:CG2	1:A:380:VAL:HG23	2.42	0.44
1:B:387:ASP:C	1:B:389:ASN:H	2.21	0.44
1:A:130:LEU:HD23	1:A:459:LEU:CD1	2.32	0.44
1:B:387:ASP:C	1:B:389:ASN:N	2.69	0.44
1:B:395:LYS:HE2	1:B:465:LEU:HD23	1.97	0.44
1:A:347:ILE:N	1:A:347:ILE:HD12	2.33	0.44
1:A:114:LEU:CD1	1:A:114:LEU:N	2.80	0.44
1:A:168:LEU:HD11	1:A:465:LEU:HD22	1.96	0.43
1:B:288:ALA:HB2	1:B:304:ASN:ND2	2.32	0.43
1:B:96:SER:O	1:B:98:PRO:HD3	2.17	0.43
1:B:344:LEU:C	1:B:344:LEU:HD23	2.39	0.43
1:A:390:ASP:HA	1:A:391:PRO:HD3	1.88	0.43
1:A:400:TYR:O	1:A:404:THR:N	2.30	0.43
1:B:199:GLN:NE2	1:B:294:ASN:CG	2.72	0.43
1:B:447:THR:OG1	1:B:450:GLN:CB	2.66	0.43
1:B:392:PHE:CB	1:B:465:LEU:HD21	2.48	0.43
1:B:394:GLN:HA	1:B:397:VAL:CG1	2.48	0.43
1:A:236:ILE:O	1:A:240:GLU:HG3	2.19	0.43
1:A:343:GLY:CA	2:A:477:HOH:O	2.57	0.42
1:A:168:LEU:N	1:A:168:LEU:HD23	2.33	0.42
1:B:165:MET:HG3	1:B:178:CYS:SG	2.59	0.42
1:B:40:PRO:HB2	1:B:44:LYS:NZ	2.34	0.42
1:A:185:ILE:HG12	1:A:186:THR:N	2.33	0.42
1:A:400:TYR:O	1:A:404:THR:HG22	2.20	0.42
1:A:463:TYR:CD1	1:A:467:VAL:HG21	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:179:ILE:O	1:B:180:HIS:C	2.55	0.42
1:B:190:LEU:HG	1:B:347:ILE:HD12	2.02	0.42
1:B:154:LYS:C	1:B:156:THR:N	2.73	0.42
1:B:392:PHE:HB2	1:B:465:LEU:HD21	2.00	0.42
1:A:218:VAL:N	1:A:268:VAL:O	2.39	0.42
1:A:78:VAL:H	1:A:167:THR:CG2	2.32	0.42
1:A:144:ILE:HG22	1:A:153:TRP:CH2	2.55	0.41
1:A:403:GLN:HB3	1:A:403:GLN:HE21	1.64	0.41
1:A:138:LYS:CG	1:A:466:CYS:O	2.67	0.41
1:A:5:LYS:HD2	1:A:133:MET:CE	2.49	0.41
1:B:114:LEU:C	1:B:115:ASP:OD2	2.58	0.41
1:B:49:MET:CA	1:B:49:MET:CE	2.82	0.41
1:A:120:THR:CG2	1:A:143:MET:CE	2.98	0.41
1:A:193:ASN:HD22	1:A:372:GLY:CA	2.33	0.41
1:A:19:GLY:HA3	1:A:48:TYR:OH	2.20	0.41
1:A:447:THR:HG21	1:A:449:LYS:CB	2.50	0.41
1:B:104:THR:HG22	1:B:105:GLU:C	2.39	0.41
1:B:299:PHE:CE2	1:B:345:PHE:CZ	3.08	0.41
1:B:397:VAL:CG2	1:B:397:VAL:O	2.69	0.41
1:B:24:GLU:CG	1:B:24:GLU:O	2.68	0.41
1:A:187:THR:HA	1:A:347:ILE:O	2.20	0.41
1:B:120:THR:HG23	1:B:143:MET:CE	2.50	0.41
1:A:381:GLN:HA	1:A:382:PRO:HD3	1.73	0.41
1:B:220:ASP:HB3	1:B:267:ALA:HB2	2.03	0.41
1:A:3:LYS:HE2	1:A:3:LYS:HB2	1.84	0.41
1:A:398:THR:O	1:A:402:ASN:HB2	2.21	0.41
1:A:173:ASP:OD2	1:A:173:ASP:N	2.48	0.41
1:A:115:ASP:HA	1:A:116:ASP:HA	1.51	0.41
1:A:190:LEU:N	1:A:190:LEU:HD23	2.36	0.41
1:B:100:GLU:HA	1:B:101:PRO:HD2	1.91	0.41
1:B:81:ILE:C	1:B:82:LEU:HD12	2.41	0.40
1:B:115:ASP:HA	1:B:116:ASP:HA	1.64	0.40
1:B:88:VAL:HG12	1:B:89:PRO:N	2.34	0.40
1:B:73:GLY:HA2	1:B:137:TRP:CD1	2.57	0.40
1:B:388:LYS:HG2	1:B:388:LYS:O	2.22	0.40
1:B:215:TYR:O	1:B:348:ASN:HB2	2.21	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	414/492 (84%)	397 (96%)	15 (4%)	2 (0%)	29	35
1	B	417/492 (85%)	399 (96%)	16 (4%)	2 (0%)	29	35
All	All	831/984 (84%)	796 (96%)	31 (4%)	4 (0%)	29	35

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	268	VAL
1	B	110	ALA
1	A	343	GLY
1	B	324	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	330/418 (79%)	304 (92%)	26 (8%)	12	15
1	B	343/418 (82%)	315 (92%)	28 (8%)	11	14
All	All	673/836 (80%)	619 (92%)	54 (8%)	12	15

All (54) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	78	VAL
1	A	82	LEU

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Mol	Chain	Res	Type
1	A	84	HIS
1	A	87	VAL
1	A	92	ASP
1	A	104	THR
1	A	115	ASP
1	A	185	ILE
1	A	186	THR
1	A	214	ARG
1	A	233	THR
1	A	248	LEU
1	A	307	LEU
1	A	309	ASN
1	A	329	THR
1	A	332	ILE
1	A	340	GLU
1	A	341	ASN
1	A	344	LEU
1	A	347	ILE
1	A	368	ILE
1	A	396	LEU
1	A	398	THR
1	A	447	THR
1	A	465	LEU
1	A	467	VAL
1	B	1	MET
1	B	5	LYS
1	B	20	LEU
1	B	21	LEU
1	B	49	MET
1	B	63	VAL
1	B	74	LYS
1	B	104	THR
1	B	115	ASP
1	B	133	MET
1	B	134	ASN
1	B	151	SER
1	B	156	THR
1	B	161	LYS
1	B	168	LEU
1	B	284	LEU
1	B	290	LEU
1	B	309	ASN

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Mol	Chain	Res	Type
1	B	339	ASN
1	B	353	GLU
1	B	361	MET
1	B	377	LEU
1	B	387	ASP
1	B	389	ASN
1	B	396	LEU
1	B	397	VAL
1	B	400	TYR
1	B	446	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	65	HIS
1	A	76	ASN
1	A	84	HIS
1	A	193	ASN
1	A	231	ASN
1	A	237	GLN
1	A	294	ASN
1	A	295	ASN
1	A	339	ASN
1	A	341	ASN
1	A	370	GLN
1	A	389	ASN
1	A	402	ASN
1	A	403	GLN
1	A	443	ASN
1	B	12	GLN
1	B	76	ASN
1	B	199	GLN
1	B	297	GLN
1	B	309	ASN
1	B	384	HIS
1	B	389	ASN
1	B	403	GLN
1	B	453	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 carbohydrates 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

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	422/492 (85%)	0.37	14 (3%) 46 53	31, 49, 76, 95	1 (0%)
1	B	425/492 (86%)	0.36	9 (2%) 63 70	29, 47, 83, 110	0
All	All	847/984 (86%)	0.37	23 (2%) 54 62	29, 48, 80, 110	1 (0%)

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	268	VAL	4.7
1	A	267	ALA	3.9
1	B	96	SER	3.2
1	A	196	THR	3.2
1	B	397	VAL	3.0
1	B	465	LEU	2.8
1	B	114	LEU	2.6
1	A	91	GLY	2.5
1	B	323	ASP	2.5
1	A	265	GLY	2.4
1	A	357	PHE	2.4
1	A	110	ALA	2.4
1	A	109	ILE	2.3
1	B	108	ILE	2.3
1	B	463	TYR	2.3
1	A	195	LEU	2.2
1	A	198	ASP	2.2
1	A	269	HIS	2.1
1	B	63	VAL	2.1
1	A	102	VAL	2.1
1	A	342	ALA	2.1
1	B	324	VAL	2.1
1	A	104	THR	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates 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.