



# Full wwPDB X-ray Structure Validation Report ⓘ

May 27, 2020 – 03:07 am BST

PDB ID : 5HV6  
Title : The ATP binding domain of rifampin phosphotransferase from *Listeria monocytogenes*  
Authors : Zhang, P.; Qi, X.  
Deposited on : 2016-01-28  
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](mailto: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.

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

MolProbity : 4.02b-467  
Xtrriage (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

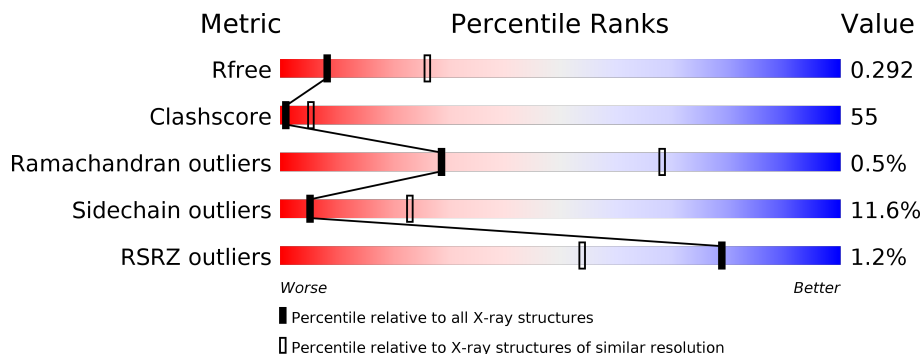
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 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))
$R_{free}$	130704	2092 (3.00-3.00)
Clashscore	141614	2416 (3.00-3.00)
Ramachandran outliers	138981	2333 (3.00-3.00)
Sidechain outliers	138945	2336 (3.00-3.00)
RSRZ outliers	127900	1990 (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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	327	
1	B	327	

## 2 Entry composition i

There are 2 unique types of molecules in this entry. The entry contains 4678 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 Phosphoenolpyruvate synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	308	Total	C	N	O	S	0	0	0
			2400	1518	410	462	10			
1	B	291	Total	C	N	O	S	0	0	0
			2277	1442	391	434	10			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-11	MET	-	expression tag	UNP A0A0S2YLC8
A	-10	ARG	-	expression tag	UNP A0A0S2YLC8
A	-9	GLY	-	expression tag	UNP A0A0S2YLC8
A	-8	SER	-	expression tag	UNP A0A0S2YLC8
A	-7	HIS	-	expression tag	UNP A0A0S2YLC8
A	-6	HIS	-	expression tag	UNP A0A0S2YLC8
A	-5	HIS	-	expression tag	UNP A0A0S2YLC8
A	-4	HIS	-	expression tag	UNP A0A0S2YLC8
A	-3	HIS	-	expression tag	UNP A0A0S2YLC8
A	-2	HIS	-	expression tag	UNP A0A0S2YLC8
A	-1	GLY	-	expression tag	UNP A0A0S2YLC8
A	0	SER	-	expression tag	UNP A0A0S2YLC8
B	-11	MET	-	expression tag	UNP A0A0S2YLC8
B	-10	ARG	-	expression tag	UNP A0A0S2YLC8
B	-9	GLY	-	expression tag	UNP A0A0S2YLC8
B	-8	SER	-	expression tag	UNP A0A0S2YLC8
B	-7	HIS	-	expression tag	UNP A0A0S2YLC8
B	-6	HIS	-	expression tag	UNP A0A0S2YLC8
B	-5	HIS	-	expression tag	UNP A0A0S2YLC8
B	-4	HIS	-	expression tag	UNP A0A0S2YLC8
B	-3	HIS	-	expression tag	UNP A0A0S2YLC8
B	-2	HIS	-	expression tag	UNP A0A0S2YLC8
B	-1	GLY	-	expression tag	UNP A0A0S2YLC8
B	0	SER	-	expression tag	UNP A0A0S2YLC8

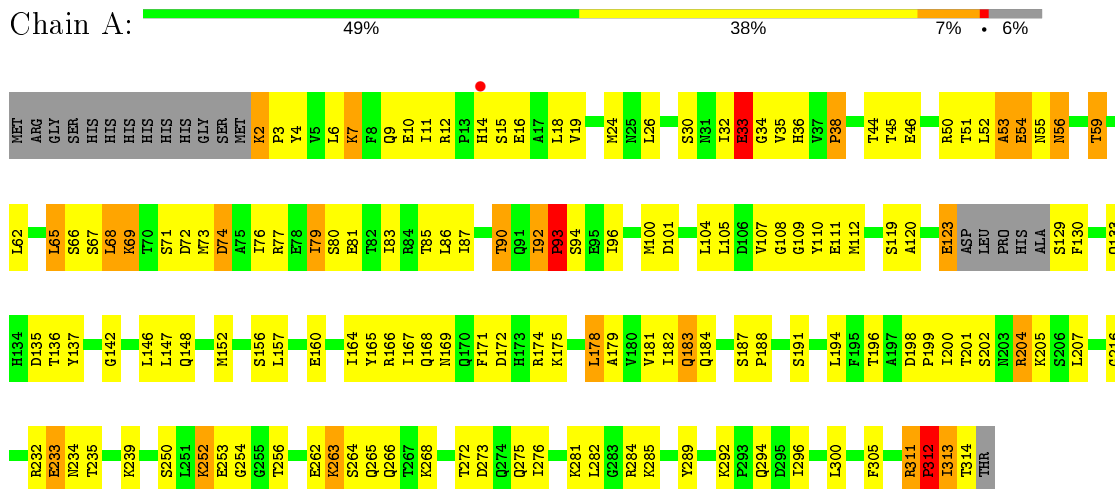
- Molecule 2 is water.

<b>Mol</b>	<b>Chain</b>	<b>Residues</b>	<b>Atoms</b>		<b>ZeroOcc</b>	<b>AltConf</b>
2	A	1	Total	O	0	0
			1	1		

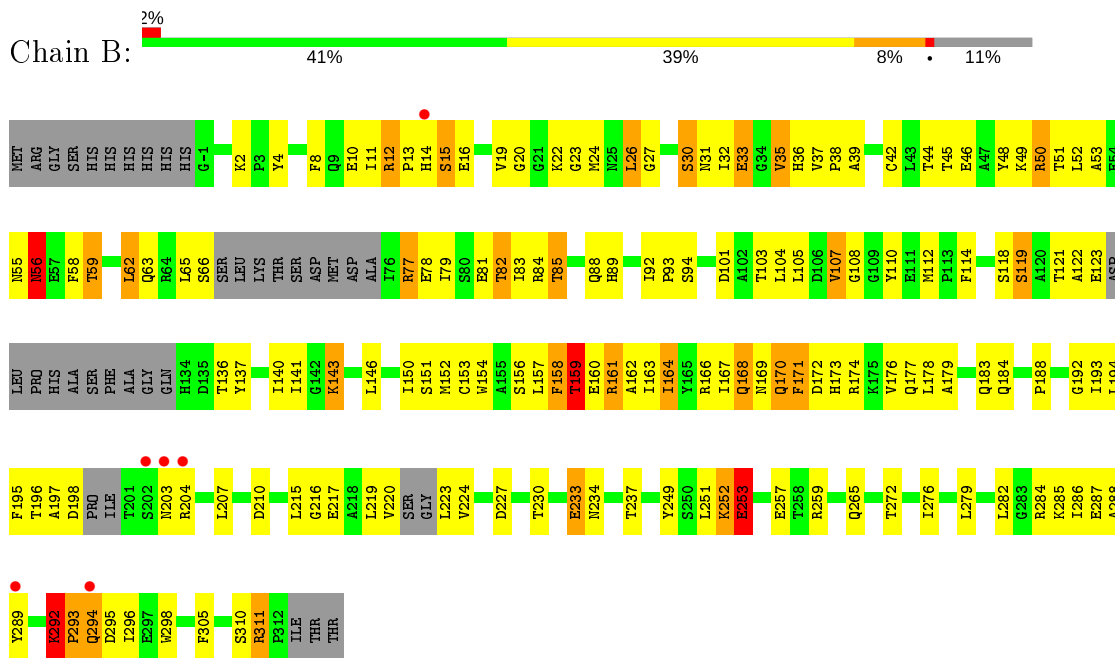
### 3 Residue-property plots

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: Phosphoenolpyruvate synthase



- Molecule 1: Phosphoenolpyruvate synthase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	85.01Å 86.18Å 100.30Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.27 – 3.00 29.27 – 3.01	Depositor EDS
% Data completeness (in resolution range)	99.9 (29.27-3.00) 84.7 (29.27-3.01)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.24 (at 3.00Å)	Xtrriage
Refinement program	PHENIX 1.9_1692	Depositor
R, $R_{free}$	0.241 , 0.295 0.246 , 0.292	Depositor DCC
$R_{free}$ test set	1326 reflections (9.95%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	48.2	Xtrriage
Anisotropy	0.060	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.29 , 49.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.023 for k,h,-l	Xtrriage
$F_o, F_c$ correlation	0.87	EDS
Total number of atoms	4678	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	44.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.32% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup> Intensities estimated from amplitudes.

<sup>2</sup> Theoretical 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	0.95	2/2438 (0.1%)	1.12	18/3295 (0.5%)
1	B	0.89	1/2310 (0.0%)	1.16	17/3116 (0.5%)
All	All	0.92	3/4748 (0.1%)	1.14	35/6411 (0.5%)

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
1	A	0	3
1	B	0	5
All	All	0	8

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	182	ILE	C-O	-6.02	1.11	1.23
1	B	293	PRO	N-CD	5.12	1.55	1.47
1	A	93	PRO	N-CD	5.09	1.54	1.47

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	233	GLU	CB-CA-C	-19.27	71.85	110.40
1	B	56	ASN	CB-CA-C	-16.87	76.66	110.40
1	B	253	GLU	CB-CA-C	-15.66	79.07	110.40
1	B	253	GLU	C-N-CA	13.82	151.32	122.30
1	B	56	ASN	C-N-CA	11.91	151.47	121.70
1	A	54	GLU	N-CA-CB	11.49	131.28	110.60
1	B	159	THR	O-C-N	-11.03	105.05	122.70
1	B	56	ASN	N-CA-C	10.98	140.65	111.00
1	A	53	ALA	CB-CA-C	-10.88	93.78	110.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	53	ALA	N-CA-C	-10.39	82.95	111.00
1	A	69	LYS	CB-CA-C	9.89	130.18	110.40
1	B	15	SER	CB-CA-C	-9.88	91.34	110.10
1	A	33	GLU	N-CA-C	9.46	136.55	111.00
1	B	253	GLU	O-C-N	-8.81	108.22	123.20
1	A	73	MET	N-CA-C	8.64	134.33	111.00
1	B	252	LYS	C-N-CA	8.37	142.61	121.70
1	B	253	GLU	CA-C-N	8.30	132.81	116.20
1	B	158	PHE	C-N-CA	7.88	141.39	121.70
1	A	264	SER	N-CA-C	-7.33	91.21	111.00
1	A	73	MET	CB-CA-C	-7.28	95.84	110.40
1	B	159	THR	CA-C-N	6.71	131.97	117.20
1	A	234	ASN	N-CA-CB	6.38	122.08	110.60
1	A	311	ARG	C-N-CD	5.98	140.97	128.40
1	A	312	PRO	CA-N-CD	-5.92	103.21	111.50
1	B	12	ARG	C-N-CD	5.85	140.68	128.40
1	A	92	ILE	C-N-CD	5.80	140.57	128.40
1	A	136	THR	N-CA-C	-5.78	95.41	111.00
1	B	292	LYS	C-N-CD	5.68	140.33	128.40
1	B	107	VAL	N-CA-C	5.57	126.04	111.00
1	A	33	GLU	CB-CA-C	-5.51	99.37	110.40
1	B	33	GLU	N-CA-C	-5.37	96.50	111.00
1	A	234	ASN	N-CA-C	5.27	125.23	111.00
1	A	72	ASP	CB-CG-OD2	5.21	122.99	118.30
1	B	198	ASP	CB-CG-OD2	5.21	122.99	118.30
1	A	74	ASP	CB-CG-OD2	5.20	122.98	118.30

There are no chirality outliers.

All (8) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	204	ARG	Mainchain
1	A	205	LYS	Mainchain
1	A	33	GLU	Peptide
1	B	159	THR	Mainchain
1	B	216	GLY	Peptide
1	B	252	LYS	Peptide
1	B	253	GLU	Peptide
1	B	56	ASN	Peptide



## 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	2400	0	2423	190	0
1	B	2277	0	2300	334	0
2	A	1	0	0	0	0
All	All	4678	0	4723	513	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 55.

All (513) 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:220:VAL:HG21	1:B:223:LEU:CB	1.23	1.66
1:B:110:TYR:CE1	1:B:143:LYS:HB2	1.15	1.65
1:B:110:TYR:CD1	1:B:143:LYS:CB	1.82	1.62
1:A:263:LYS:CA	1:A:266:GLN:HG3	1.16	1.61
1:B:13:PRO:HB3	1:B:14:HIS:CE1	1.09	1.61
1:B:65:LEU:HD21	1:B:79:ILE:CG2	1.29	1.61
1:B:13:PRO:HB2	1:B:14:HIS:CG	1.36	1.60
1:A:263:LYS:HA	1:A:266:GLN:CG	1.24	1.59
1:B:220:VAL:HG11	1:B:223:LEU:CD1	1.30	1.55
1:A:6:LEU:CB	1:A:11:ILE:HD11	1.35	1.54
1:B:110:TYR:CD1	1:B:143:LYS:HB3	1.40	1.52
1:B:13:PRO:CB	1:B:14:HIS:ND1	1.73	1.52
1:B:196:THR:HG21	1:B:287:GLU:CD	1.34	1.48
1:A:6:LEU:HB3	1:A:11:ILE:CD1	1.39	1.47
1:B:249:TYR:CD2	1:B:259:ARG:HD2	1.48	1.46
1:B:13:PRO:CB	1:B:14:HIS:CE1	2.00	1.44
1:B:220:VAL:CG2	1:B:223:LEU:HB2	1.48	1.43
1:B:110:TYR:CE1	1:B:143:LYS:CB	1.90	1.41
1:B:48:TYR:CE1	1:B:52:LEU:HD12	1.56	1.40
1:B:286:ILE:CG2	1:B:294:GLN:OE1	1.67	1.39
1:B:65:LEU:CD2	1:B:79:ILE:HG21	1.54	1.36
1:B:55:ASN:CA	1:B:56:ASN:HB2	1.01	1.36
1:B:13:PRO:CB	1:B:14:HIS:CG	2.06	1.33
1:B:13:PRO:HB3	1:B:14:HIS:ND1	1.33	1.33

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:65:LEU:CD2	1:B:79:ILE:CG2	2.06	1.31
1:B:16:GLU:O	1:B:121:THR:CG2	1.77	1.30
1:B:220:VAL:CG1	1:B:223:LEU:CD1	2.08	1.30
1:B:55:ASN:HA	1:B:56:ASN:CB	1.17	1.29
1:B:157:LEU:O	1:B:166:ARG:NH1	1.65	1.27
1:B:22:LYS:HB2	1:B:119:SER:OG	1.16	1.27
1:B:196:THR:CG2	1:B:287:GLU:CD	2.05	1.24
1:A:68:LEU:O	1:A:69:LYS:HG3	1.34	1.24
1:B:286:ILE:HG21	1:B:294:GLN:OE1	1.18	1.22
1:B:220:VAL:HG11	1:B:223:LEU:CG	1.70	1.21
1:A:174:ARG:NH2	1:B:217:GLU:OE1	1.74	1.19
1:B:159:THR:HG22	1:B:162:ALA:N	1.57	1.19
1:B:16:GLU:O	1:B:121:THR:HG21	1.03	1.18
1:B:13:PRO:HB2	1:B:14:HIS:CD2	1.78	1.17
1:B:220:VAL:HG21	1:B:223:LEU:CG	1.73	1.17
1:B:249:TYR:CE2	1:B:259:ARG:HD2	1.79	1.16
1:A:68:LEU:CD1	1:A:167:ILE:HD11	1.74	1.16
1:A:6:LEU:HD22	1:A:11:ILE:CD1	1.77	1.15
1:B:249:TYR:CD2	1:B:259:ARG:CD	2.30	1.15
1:B:220:VAL:CG1	1:B:223:LEU:HD12	1.75	1.15
1:A:6:LEU:CD2	1:A:11:ILE:HD13	1.78	1.13
1:A:68:LEU:HD12	1:A:167:ILE:HD11	1.15	1.13
1:A:294:GLN:OE1	1:A:311:ARG:O	1.67	1.12
1:A:263:LYS:HG2	1:A:263:LYS:O	1.49	1.11
1:B:14:HIS:HA	1:B:15:SER:HB2	1.19	1.11
1:A:200:ILE:HD12	1:A:201:THR:H	1.14	1.09
1:B:65:LEU:HD23	1:B:79:ILE:HG21	1.26	1.08
1:B:220:VAL:HG11	1:B:223:LEU:HD11	1.20	1.08
1:B:48:TYR:HE1	1:B:52:LEU:CD1	1.66	1.08
1:B:107:VAL:HB	1:B:184:GLN:NE2	1.68	1.07
1:A:198:ASP:OD1	1:A:200:ILE:CD1	2.01	1.07
1:B:196:THR:HG21	1:B:287:GLU:OE2	1.52	1.07
1:B:220:VAL:CG2	1:B:223:LEU:CB	2.18	1.06
1:B:159:THR:CG2	1:B:162:ALA:H	1.66	1.06
1:A:44:THR:HG22	1:A:46:GLU:H	1.20	1.06
1:B:56:ASN:N	1:B:59:THR:OG1	1.87	1.05
1:B:16:GLU:N	1:B:24:MET:HE3	1.71	1.05
1:B:104:LEU:HD23	1:B:110:TYR:CE2	1.91	1.05
1:A:65:LEU:O	1:A:68:LEU:HB3	1.54	1.05
1:A:68:LEU:CD1	1:A:167:ILE:CD1	2.34	1.04
1:B:65:LEU:HD21	1:B:79:ILE:HG23	1.10	1.04

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:15:SER:O	1:A:19:VAL:HG22	1.58	1.03
1:A:198:ASP:OD1	1:A:200:ILE:HD11	1.55	1.03
1:B:110:TYR:HD1	1:B:143:LYS:CA	1.72	1.03
1:B:55:ASN:CB	1:B:56:ASN:HB2	1.89	1.03
1:A:175:LYS:HA	1:B:220:VAL:CG1	1.89	1.03
1:B:65:LEU:CD1	1:B:83:ILE:HD11	1.88	1.02
1:A:6:LEU:CB	1:A:11:ILE:CD1	2.13	1.02
1:A:313:ILE:HG22	1:A:314:THR:HA	1.41	1.02
1:A:200:ILE:HD12	1:A:201:THR:N	1.74	1.02
1:B:32:ILE:HD11	1:B:289:TYR:CE1	1.94	1.01
1:A:6:LEU:HD13	1:A:11:ILE:HD12	1.43	1.00
1:B:110:TYR:HD1	1:B:143:LYS:CB	1.42	0.99
1:B:249:TYR:HD2	1:B:259:ARG:CD	1.69	0.99
1:A:263:LYS:HA	1:A:266:GLN:CB	1.92	0.99
1:B:22:LYS:CB	1:B:119:SER:OG	2.11	0.98
1:A:175:LYS:HA	1:B:220:VAL:HG12	1.46	0.98
1:B:32:ILE:HD11	1:B:289:TYR:CZ	1.99	0.97
1:B:83:ILE:HG21	1:B:158:PHE:CD2	1.99	0.97
1:B:159:THR:HG22	1:B:162:ALA:H	0.82	0.97
1:B:48:TYR:HE1	1:B:52:LEU:HD12	0.88	0.97
1:B:22:LYS:O	1:B:26:LEU:HB2	1.64	0.97
1:B:58:PHE:CZ	1:B:62:LEU:HD11	2.01	0.96
1:A:52:LEU:HD13	1:A:87:ILE:HD13	1.43	0.96
1:B:13:PRO:CA	1:B:14:HIS:ND1	2.29	0.95
1:A:32:ILE:HD12	1:A:35:VAL:HG11	1.49	0.95
1:B:22:LYS:HD2	1:B:118:SER:O	1.66	0.95
1:B:110:TYR:CD1	1:B:143:LYS:CA	2.49	0.95
1:A:204:ARG:HG3	1:A:204:ARG:HH11	1.29	0.94
1:B:220:VAL:CB	1:B:223:LEU:HG	1.98	0.94
1:A:6:LEU:CG	1:A:11:ILE:CD1	2.45	0.94
1:B:14:HIS:HA	1:B:15:SER:CB	1.98	0.94
1:A:11:ILE:HG22	1:A:15:SER:OG	1.66	0.93
1:B:81:GLU:O	1:B:85:THR:OG1	1.85	0.93
1:B:13:PRO:HB3	1:B:14:HIS:NE2	1.83	0.93
1:A:6:LEU:CD2	1:A:11:ILE:CD1	2.42	0.92
1:B:196:THR:CG2	1:B:287:GLU:OE1	2.15	0.92
1:B:220:VAL:HG11	1:B:223:LEU:HG	1.51	0.92
1:B:22:LYS:HB2	1:B:119:SER:HG	1.28	0.92
1:B:65:LEU:HD11	1:B:83:ILE:CD1	1.99	0.92
1:B:13:PRO:CB	1:B:14:HIS:CD2	2.45	0.91
1:B:196:THR:CG2	1:B:287:GLU:OE2	2.13	0.91

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:65:LEU:HD12	1:B:83:ILE:HD11	1.52	0.91
1:A:252:LYS:NZ	1:A:252:LYS:HB2	1.86	0.90
1:A:68:LEU:O	1:A:69:LYS:CG	2.18	0.90
1:A:6:LEU:HD22	1:A:11:ILE:HD13	0.91	0.90
1:B:65:LEU:CD1	1:B:83:ILE:CD1	2.50	0.90
1:A:263:LYS:N	1:A:266:GLN:HG3	1.87	0.90
1:A:86:LEU:O	1:A:90:THR:OG1	1.90	0.90
1:B:159:THR:HB	1:B:162:ALA:HB3	1.53	0.90
1:B:159:THR:CG2	1:B:162:ALA:N	2.31	0.89
1:B:8:PHE:O	1:B:11:ILE:HG22	1.73	0.89
1:B:286:ILE:HG22	1:B:294:GLN:OE1	1.72	0.89
1:A:16:GLU:OE2	1:A:123:GLU:HB3	1.71	0.89
1:B:48:TYR:CE2	1:B:177:GLN:HA	2.08	0.88
1:B:220:VAL:CG1	1:B:223:LEU:HG	2.02	0.88
1:B:58:PHE:CE1	1:B:62:LEU:HD11	2.09	0.88
1:B:107:VAL:CG2	1:B:108:GLY:HA2	2.03	0.88
1:A:55:ASN:OD1	1:A:56:ASN:N	2.06	0.87
1:B:159:THR:HB	1:B:162:ALA:CB	2.04	0.87
1:B:16:GLU:CA	1:B:24:MET:HE3	2.04	0.86
1:B:220:VAL:CG2	1:B:223:LEU:CG	2.49	0.86
1:B:233:GLU:N	1:B:284:ARG:HH22	1.73	0.86
1:A:272:THR:H	1:A:275:GLN:HE21	1.21	0.86
1:B:23:GLY:O	1:B:27:GLY:N	2.08	0.86
1:B:197:ALA:HB2	1:B:204:ARG:HA	1.56	0.85
1:B:294:GLN:HG2	1:B:296:ILE:HG13	1.58	0.85
1:A:81:GLU:O	1:A:85:THR:HG23	1.77	0.85
1:B:197:ALA:CB	1:B:204:ARG:HA	2.07	0.85
1:B:220:VAL:HG13	1:B:223:LEU:HD12	1.56	0.85
1:B:107:VAL:HG22	1:B:108:GLY:HA2	1.59	0.84
1:B:233:GLU:H	1:B:284:ARG:HH22	1.24	0.84
1:A:201:THR:O	1:A:202:SER:OG	1.95	0.84
1:A:18:LEU:O	1:A:45:THR:HG23	1.76	0.84
1:B:104:LEU:O	1:B:107:VAL:HG22	1.77	0.84
1:B:101:ASP:OD1	1:B:110:TYR:OH	1.96	0.83
1:A:52:LEU:HD13	1:A:87:ILE:CD1	2.07	0.83
1:B:249:TYR:HD2	1:B:259:ARG:HD2	1.18	0.83
1:A:119:SER:HB2	1:A:181:VAL:HG23	1.59	0.83
1:A:68:LEU:HD11	1:A:167:ILE:CD1	2.08	0.83
1:A:11:ILE:CG2	1:A:15:SER:OG	2.27	0.82
1:B:13:PRO:CB	1:B:14:HIS:NE2	2.40	0.82
1:B:16:GLU:HA	1:B:24:MET:CE	2.10	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:101:ASP:O	1:A:105:LEU:HG	1.79	0.82
1:A:79:ILE:O	1:A:83:ILE:HD12	1.81	0.81
1:B:110:TYR:HE1	1:B:143:LYS:CB	1.58	0.81
1:A:263:LYS:CB	1:A:266:GLN:HG3	2.09	0.81
1:B:294:GLN:HG2	1:B:296:ILE:CG1	2.11	0.81
1:A:165:TYR:OH	1:B:161:ARG:HD2	1.81	0.81
1:B:55:ASN:HA	1:B:56:ASN:CG	2.00	0.80
1:A:200:ILE:CD1	1:A:201:THR:N	2.45	0.80
1:B:77:ARG:H	1:B:77:ARG:HD2	1.45	0.80
1:B:16:GLU:HA	1:B:24:MET:HE1	1.62	0.80
1:A:68:LEU:HG	1:A:167:ILE:HD13	1.64	0.79
1:A:313:ILE:HG22	1:A:314:THR:CA	2.12	0.79
1:A:168:GLN:HE22	1:B:160:GLU:CD	1.86	0.79
1:B:65:LEU:HD11	1:B:83:ILE:HD11	1.56	0.79
1:B:294:GLN:HG3	1:B:295:ASP:H	1.48	0.79
1:B:33:GLU:HA	1:B:35:VAL:H	1.47	0.79
1:B:107:VAL:HB	1:B:184:GLN:HE21	1.45	0.78
1:B:110:TYR:CD1	1:B:143:LYS:HB2	1.69	0.78
1:B:16:GLU:CA	1:B:24:MET:CE	2.61	0.78
1:B:32:ILE:HG22	1:B:35:VAL:CG1	2.14	0.78
1:A:169:ASN:OD1	1:B:159:THR:HG23	1.83	0.78
1:A:79:ILE:H	1:A:79:ILE:HD12	1.46	0.77
1:B:32:ILE:O	1:B:33:GLU:HG3	1.84	0.77
1:A:6:LEU:CD1	1:A:11:ILE:HD12	2.15	0.77
1:A:55:ASN:O	1:A:59:THR:OG1	2.03	0.77
1:A:135:ASP:OD2	1:A:156:SER:OG	1.99	0.77
1:B:251:LEU:CB	1:B:253:GLU:O	2.33	0.77
1:B:251:LEU:HB3	1:B:253:GLU:O	1.85	0.76
1:B:65:LEU:O	1:B:66:SER:OG	2.03	0.76
1:A:200:ILE:HD13	1:A:201:THR:HB	1.66	0.76
1:B:32:ILE:CD1	1:B:289:TYR:CZ	2.68	0.75
1:B:119:SER:O	1:B:178:LEU:HD12	1.87	0.75
1:A:6:LEU:HD13	1:A:11:ILE:CD1	2.16	0.75
1:B:249:TYR:HD2	1:B:259:ARG:CG	1.99	0.75
1:B:15:SER:HB3	1:B:24:MET:SD	2.27	0.74
1:B:141:ILE:O	1:B:141:ILE:HD12	1.86	0.74
1:B:287:GLU:HG3	1:B:292:LYS:O	1.88	0.74
1:B:16:GLU:N	1:B:24:MET:CE	2.50	0.73
1:A:232:ARG:O	1:A:233:GLU:HB2	1.86	0.73
1:B:55:ASN:CA	1:B:56:ASN:CB	1.95	0.73
1:B:146:LEU:O	1:B:146:LEU:HD12	1.87	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:104:LEU:HD11	1:B:114:PHE:CZ	2.24	0.72
1:B:233:GLU:H	1:B:284:ARG:NH2	1.87	0.72
1:B:55:ASN:CB	1:B:56:ASN:CB	2.60	0.72
1:B:48:TYR:CE1	1:B:52:LEU:CD1	2.49	0.72
1:B:196:THR:HG21	1:B:287:GLU:CG	2.19	0.72
1:B:220:VAL:HG21	1:B:223:LEU:HB2	0.73	0.72
1:B:48:TYR:CD1	1:B:52:LEU:HD12	2.22	0.72
1:B:65:LEU:HD11	1:B:83:ILE:HD12	1.72	0.72
1:B:15:SER:O	1:B:24:MET:SD	2.47	0.72
1:A:204:ARG:HG3	1:A:204:ARG:NH1	2.04	0.72
1:B:220:VAL:CB	1:B:223:LEU:CG	2.67	0.72
1:B:14:HIS:CA	1:B:15:SER:HB2	2.10	0.71
1:B:11:ILE:HD11	1:B:19:VAL:HG21	1.73	0.71
1:A:110:TYR:HB3	1:A:142:GLY:HA2	1.73	0.70
1:B:107:VAL:CB	1:B:184:GLN:NE2	2.52	0.70
1:B:15:SER:CB	1:B:24:MET:SD	2.80	0.70
1:A:16:GLU:HB2	1:A:24:MET:HG3	1.74	0.70
1:B:110:TYR:HD1	1:B:143:LYS:N	1.89	0.70
1:B:79:ILE:O	1:B:82:THR:OG1	2.10	0.69
1:B:16:GLU:HG2	1:B:24:MET:CE	2.23	0.69
1:A:263:LYS:CA	1:A:266:GLN:CG	2.11	0.69
1:A:252:LYS:HZ1	1:A:252:LYS:HB2	1.57	0.69
1:B:159:THR:O	1:B:163:ILE:HG13	1.93	0.69
1:B:32:ILE:HG22	1:B:35:VAL:HG12	1.74	0.69
1:B:13:PRO:HA	1:B:14:HIS:ND1	2.06	0.69
1:B:16:GLU:CG	1:B:24:MET:HE3	2.22	0.69
1:B:220:VAL:HB	1:B:223:LEU:HG	1.74	0.68
1:B:104:LEU:CD2	1:B:110:TYR:CE2	2.74	0.68
1:B:44:THR:HG22	1:B:46:GLU:H	1.59	0.68
1:A:68:LEU:HD11	1:A:167:ILE:HG12	1.74	0.68
1:B:122:ALA:O	1:B:123:GLU:HB2	1.93	0.68
1:B:164:ILE:O	1:B:168:GLN:HG2	1.93	0.68
1:A:6:LEU:CD1	1:A:11:ILE:CD1	2.71	0.68
1:A:135:ASP:HB3	1:A:137:TYR:HE1	1.59	0.68
1:A:198:ASP:OD1	1:A:200:ILE:HD12	1.91	0.68
1:B:14:HIS:CA	1:B:15:SER:CB	2.72	0.67
1:A:119:SER:HB3	1:A:179:ALA:O	1.94	0.67
1:B:104:LEU:HD23	1:B:110:TYR:CD2	2.29	0.67
1:B:11:ILE:C	1:B:12:ARG:HG3	2.15	0.67
1:A:120:ALA:HB2	1:A:133:GLN:HB3	1.77	0.67
1:B:15:SER:C	1:B:24:MET:CE	2.63	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:282:LEU:O	1:B:286:ILE:HD13	1.96	0.66
1:A:157:LEU:HD11	1:A:178:LEU:HB2	1.77	0.66
1:A:68:LEU:HD23	1:A:68:LEU:C	2.15	0.66
1:A:34:GLY:HA2	1:A:35:VAL:HB	1.78	0.65
1:A:123:GLU:N	1:A:123:GLU:OE2	2.28	0.65
1:A:68:LEU:HD11	1:A:167:ILE:CG1	2.26	0.65
1:B:58:PHE:O	1:B:62:LEU:HD12	1.96	0.65
1:A:200:ILE:CD1	1:A:201:THR:HB	2.26	0.65
1:B:16:GLU:HG2	1:B:24:MET:HE3	1.79	0.65
1:B:15:SER:O	1:B:24:MET:CE	2.45	0.65
1:B:220:VAL:CG1	1:B:223:LEU:HD11	1.99	0.64
1:A:15:SER:HB3	1:A:19:VAL:CG2	2.27	0.64
1:A:38:PRO:HB2	1:A:183:GLN:HG2	1.80	0.64
1:A:194:LEU:HD22	1:A:296:ILE:HD12	1.78	0.64
1:B:48:TYR:HE2	1:B:177:GLN:HA	1.58	0.64
1:A:15:SER:HB3	1:A:19:VAL:HG22	1.80	0.63
1:A:168:GLN:HG3	1:B:84:ARG:NH2	2.13	0.63
1:A:168:GLN:NE2	1:B:160:GLU:CD	2.51	0.63
1:B:15:SER:C	1:B:24:MET:HE3	2.19	0.63
1:B:110:TYR:HD1	1:B:143:LYS:HB3	1.05	0.63
1:B:107:VAL:HG23	1:B:108:GLY:HA2	1.80	0.62
1:B:16:GLU:O	1:B:121:THR:HG22	1.93	0.62
1:B:107:VAL:CB	1:B:184:GLN:HE21	2.11	0.62
1:B:203:ASN:O	1:B:204:ARG:HB2	2.00	0.62
1:B:215:LEU:N	1:B:215:LEU:HD12	2.15	0.62
1:A:239:LYS:NZ	1:A:273:ASP:OD1	2.30	0.62
1:A:34:GLY:HA3	1:A:36:HIS:CD2	2.35	0.61
1:A:92:ILE:HD12	1:A:147:LEU:HD23	1.81	0.61
1:B:146:LEU:O	1:B:150:ILE:HG13	2.00	0.61
1:B:220:VAL:CG2	1:B:223:LEU:HG	2.25	0.61
1:B:251:LEU:HB2	1:B:253:GLU:O	2.00	0.61
1:B:77:ARG:HH11	1:B:77:ARG:HG2	1.63	0.61
1:B:27:GLY:O	1:B:30:SER:N	2.33	0.61
1:A:201:THR:C	1:A:202:SER:HG	1.98	0.61
1:B:233:GLU:O	1:B:234:ASN:HB2	2.00	0.61
1:A:135:ASP:HB3	1:A:137:TYR:CE1	2.35	0.61
1:B:19:VAL:HG12	1:B:42:CYS:SG	2.41	0.61
1:B:136:THR:CG2	1:B:217:GLU:HG2	2.31	0.61
1:A:15:SER:O	1:A:19:VAL:CG2	2.43	0.60
1:A:68:LEU:CG	1:A:167:ILE:HD13	2.30	0.60
1:A:272:THR:HG23	1:A:275:GLN:NE2	2.17	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:48:TYR:CD2	1:B:177:GLN:HA	2.35	0.60
1:A:265:GLN:HA	1:A:268:LYS:HB2	1.84	0.60
1:B:65:LEU:HD21	1:B:79:ILE:HG22	1.63	0.60
1:A:263:LYS:HB2	1:A:266:GLN:CD	2.22	0.60
1:A:165:TYR:OH	1:B:161:ARG:CD	2.47	0.59
1:A:198:ASP:CG	1:A:200:ILE:HD11	2.22	0.59
1:A:313:ILE:HA	1:A:314:THR:C	2.23	0.59
1:B:11:ILE:O	1:B:12:ARG:HG3	2.02	0.59
1:A:101:ASP:OD1	1:A:110:TYR:OH	2.18	0.59
1:A:74:ASP:OD1	1:A:77:ARG:NH2	2.36	0.59
1:A:314:THR:O	1:A:314:THR:HG22	2.01	0.59
1:A:2:LYS:O	1:A:2:LYS:HD3	2.03	0.59
1:B:194:LEU:CD2	1:B:207:LEU:HD11	2.33	0.58
1:B:294:GLN:HG3	1:B:295:ASP:N	2.16	0.58
1:B:58:PHE:CE2	1:B:62:LEU:CD1	2.86	0.58
1:B:153:CYS:O	1:B:156:SER:HB3	2.04	0.58
1:B:22:LYS:CB	1:B:119:SER:HG	2.05	0.58
1:B:167:ILE:O	1:B:170:GLN:CD	2.42	0.58
1:B:65:LEU:HD12	1:B:83:ILE:CD1	2.23	0.58
1:B:169:ASN:N	1:B:170:GLN:HA	2.19	0.58
1:B:194:LEU:HD21	1:B:207:LEU:HD11	1.86	0.58
1:B:11:ILE:CG2	1:B:27:GLY:HA3	2.34	0.57
1:A:7:LYS:O	1:A:11:ILE:HG12	2.04	0.57
1:A:52:LEU:C	1:A:53:ALA:O	2.28	0.57
1:A:52:LEU:CD1	1:A:87:ILE:CD1	2.80	0.57
1:B:210:ASP:OD2	1:B:219:LEU:HD11	2.04	0.57
1:B:58:PHE:CE2	1:B:62:LEU:HD11	2.38	0.57
1:A:2:LYS:CA	1:A:46:GLU:OE1	2.53	0.57
1:B:171:PHE:N	1:B:171:PHE:CD1	2.73	0.57
1:A:265:GLN:HG3	1:A:268:LYS:HD3	1.86	0.56
1:B:20:GLY:HA2	1:B:121:THR:HG23	1.85	0.56
1:A:262:GLU:O	1:A:263:LYS:HB3	2.05	0.56
1:B:220:VAL:CG2	1:B:223:LEU:HD12	2.34	0.56
1:B:15:SER:HB2	1:B:24:MET:SD	2.46	0.55
1:B:169:ASN:HB3	1:B:171:PHE:HE1	1.71	0.55
1:B:49:LYS:O	1:B:53:ALA:HB3	2.07	0.55
1:B:51:THR:HG23	1:B:93:PRO:HD3	1.87	0.55
1:A:68:LEU:CG	1:A:167:ILE:CD1	2.84	0.55
1:A:32:ILE:HD12	1:A:35:VAL:CG1	2.32	0.55
1:B:286:ILE:HG22	1:B:294:GLN:CD	2.27	0.55
1:A:35:VAL:O	1:A:35:VAL:HG12	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:65:LEU:CD2	1:B:79:ILE:HG23	2.01	0.54
1:B:58:PHE:CZ	1:B:62:LEU:CD1	2.85	0.54
1:A:18:LEU:HG	1:A:44:THR:HG23	1.89	0.54
1:B:196:THR:HG22	1:B:287:GLU:CD	2.21	0.54
1:B:119:SER:O	1:B:178:LEU:CD1	2.54	0.54
1:B:20:GLY:HA2	1:B:121:THR:CG2	2.38	0.54
1:B:286:ILE:CG2	1:B:294:GLN:CD	2.66	0.54
1:B:15:SER:O	1:B:19:VAL:CG2	2.56	0.54
1:B:220:VAL:CG2	1:B:223:LEU:N	2.70	0.54
1:A:204:ARG:NH1	1:A:204:ARG:CG	2.68	0.54
1:A:6:LEU:CB	1:A:11:ILE:HD12	2.30	0.54
1:B:196:THR:HG22	1:B:287:GLU:OE2	2.01	0.54
1:B:294:GLN:HG2	1:B:296:ILE:HG12	1.88	0.54
1:B:32:ILE:CG2	1:B:35:VAL:CG1	2.85	0.54
1:B:159:THR:HB	1:B:162:ALA:HB2	1.85	0.54
1:B:101:ASP:O	1:B:105:LEU:HD13	2.07	0.53
1:B:220:VAL:HG21	1:B:223:LEU:CD1	2.35	0.53
1:B:233:GLU:CA	1:B:284:ARG:HH22	2.20	0.53
1:B:272:THR:O	1:B:276:ILE:HG13	2.08	0.53
1:B:16:GLU:HG2	1:B:24:MET:HE2	1.90	0.53
1:B:136:THR:HG22	1:B:217:GLU:HG2	1.88	0.53
1:B:171:PHE:HD1	1:B:171:PHE:N	2.07	0.53
1:B:4:TYR:OH	1:B:50:ARG:HG3	2.08	0.53
1:B:220:VAL:CG2	1:B:223:LEU:CD1	2.86	0.53
1:B:104:LEU:HD23	1:B:110:TYR:HE2	1.63	0.53
1:B:107:VAL:HG23	1:B:108:GLY:CA	2.38	0.53
1:B:197:ALA:HB1	1:B:204:ARG:HA	1.90	0.53
1:A:252:LYS:HZ2	1:A:252:LYS:HB2	1.70	0.53
1:B:293:PRO:O	1:B:293:PRO:HG2	2.08	0.52
1:A:15:SER:C	1:A:19:VAL:HG22	2.28	0.52
1:B:233:GLU:OE1	1:B:233:GLU:HA	2.10	0.52
1:A:174:ARG:O	1:B:220:VAL:CG1	2.44	0.52
1:B:107:VAL:CG2	1:B:108:GLY:CA	2.85	0.52
1:B:251:LEU:HD21	1:B:257:GLU:HB3	1.92	0.52
1:A:68:LEU:HG	1:A:167:ILE:CD1	2.37	0.52
1:B:194:LEU:HD23	1:B:195:PHE:N	2.25	0.52
1:B:136:THR:O	1:B:217:GLU:OE2	2.28	0.52
1:B:104:LEU:CD2	1:B:110:TYR:CD2	2.93	0.52
1:B:159:THR:HG22	1:B:161:ARG:N	2.24	0.52
1:B:220:VAL:HG21	1:B:223:LEU:CA	2.24	0.52
1:B:59:THR:O	1:B:63:GLN:HB2	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:34:GLY:HA2	1:A:35:VAL:C	2.30	0.51
1:A:263:LYS:CG	1:A:263:LYS:O	2.33	0.51
1:B:13:PRO:HA	1:B:15:SER:HB2	1.91	0.51
1:B:295:ASP:HB3	1:B:311:ARG:O	2.10	0.51
1:A:172:ASP:OD1	1:A:175:LYS:HD2	2.10	0.51
1:A:263:LYS:CB	1:A:266:GLN:CG	2.81	0.51
1:B:26:LEU:HD11	1:B:183:GLN:HE21	1.75	0.51
1:B:193:ILE:HG22	1:B:194:LEU:N	2.26	0.51
1:B:39:ALA:HB3	1:B:184:GLN:OE1	2.10	0.51
1:B:77:ARG:NH1	1:B:77:ARG:HG2	2.26	0.51
1:A:68:LEU:CD1	1:A:167:ILE:HD13	2.31	0.51
1:B:49:LYS:O	1:B:53:ALA:CB	2.58	0.51
1:B:78:GLU:O	1:B:81:GLU:HB3	2.11	0.51
1:B:292:LYS:HG2	1:B:293:PRO:HD2	1.93	0.50
1:A:300:LEU:HD13	1:A:305:PHE:CZ	2.46	0.50
1:B:107:VAL:O	1:B:184:GLN:NE2	2.44	0.50
1:A:51:THR:HG22	1:A:52:LEU:HD23	1.93	0.50
1:A:112:MET:HB2	1:A:184:GLN:NE2	2.26	0.50
1:B:2:LYS:CB	1:B:46:GLU:HG3	2.42	0.50
1:A:2:LYS:C	1:A:2:LYS:HD3	2.31	0.50
1:B:163:ILE:O	1:B:166:ARG:HB3	2.11	0.50
1:B:196:THR:HG23	1:B:287:GLU:OE1	2.06	0.50
1:A:109:GLY:C	1:A:111:GLU:H	2.14	0.50
1:B:31:ASN:O	1:B:32:ILE:HD12	2.11	0.50
1:B:15:SER:C	1:B:24:MET:SD	2.89	0.49
1:B:32:ILE:CD1	1:B:289:TYR:CE2	2.95	0.49
1:A:191:SER:HB3	1:A:216:GLY:HA2	1.93	0.49
1:B:286:ILE:N	1:B:286:ILE:CD1	2.75	0.49
1:A:2:LYS:HA	1:A:46:GLU:OE1	2.12	0.49
1:B:32:ILE:HG22	1:B:33:GLU:N	2.27	0.49
1:A:148:GLN:O	1:A:152:MET:HG3	2.13	0.49
1:B:13:PRO:CA	1:B:14:HIS:CG	2.86	0.49
1:A:2:LYS:HB3	1:A:3:PRO:HD3	1.94	0.49
1:A:253:GLU:HG2	1:A:254:GLY:N	2.27	0.49
1:A:165:TYR:OH	1:B:161:ARG:NH1	2.46	0.49
1:A:52:LEU:CD1	1:A:87:ILE:HD13	2.29	0.48
1:A:196:THR:HG22	1:A:207:LEU:HD12	1.94	0.48
1:B:220:VAL:HG23	1:B:223:LEU:N	2.27	0.48
1:A:198:ASP:HB3	1:A:201:THR:HG22	1.95	0.48
1:B:122:ALA:HB3	1:B:123:GLU:OE1	2.12	0.48
1:A:312:PRO:HD2	1:A:312:PRO:O	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:66:SER:C	1:A:68:LEU:H	2.17	0.48
1:B:172:ASP:OD1	1:B:173:HIS:N	2.47	0.48
1:B:58:PHE:CD2	1:B:62:LEU:HD13	2.49	0.48
1:B:286:ILE:HD12	1:B:286:ILE:N	2.29	0.48
1:A:77:ARG:HG2	1:A:160:GLU:OE1	2.14	0.48
1:B:150:ILE:O	1:B:153:CYS:HB2	2.13	0.48
1:B:173:HIS:O	1:B:176:VAL:HG12	2.14	0.48
1:A:281:LYS:O	1:A:285:LYS:N	2.41	0.48
1:B:285:LYS:O	1:B:288:ALA:HB3	2.14	0.47
1:A:35:VAL:H	1:A:282:LEU:HD11	1.79	0.47
1:A:68:LEU:HG	1:A:69:LYS:N	2.30	0.47
1:B:193:ILE:HG22	1:B:194:LEU:H	1.80	0.47
1:A:100:MET:HE1	1:A:146:LEU:HD11	1.95	0.47
1:B:2:LYS:HB3	1:B:46:GLU:HG3	1.97	0.47
1:A:272:THR:O	1:A:275:GLN:N	2.48	0.47
1:A:76:ILE:O	1:A:80:SER:HB3	2.15	0.47
1:B:151:SER:O	1:B:152:MET:C	2.53	0.47
1:B:15:SER:O	1:B:19:VAL:HG22	2.15	0.47
1:B:192:GLY:HA3	1:B:210:ASP:O	2.15	0.47
1:B:32:ILE:CG2	1:B:33:GLU:N	2.78	0.47
1:B:110:TYR:HE1	1:B:143:LYS:HB2	0.69	0.47
1:B:230:THR:O	1:B:237:THR:HB	2.16	0.46
1:B:58:PHE:CE2	1:B:62:LEU:HD13	2.49	0.46
1:A:313:ILE:HG22	1:A:314:THR:C	2.36	0.46
1:B:176:VAL:HG22	1:B:177:GLN:N	2.29	0.46
1:B:83:ILE:CD1	1:B:158:PHE:HD2	2.29	0.46
1:B:169:ASN:H	1:B:170:GLN:HA	1.80	0.46
1:B:178:LEU:HD12	1:B:179:ALA:H	1.81	0.46
1:B:45:THR:O	1:B:48:TYR:HB3	2.15	0.46
1:A:263:LYS:HA	1:A:266:GLN:HB2	1.90	0.46
1:B:194:LEU:HD12	1:B:279:LEU:HG	1.98	0.46
1:A:11:ILE:CG2	1:A:15:SER:CB	2.93	0.45
1:A:32:ILE:HG13	1:A:32:ILE:H	1.64	0.45
1:B:22:LYS:O	1:B:26:LEU:CB	2.51	0.45
1:B:32:ILE:HD13	1:B:289:TYR:CE2	2.52	0.45
1:A:2:LYS:C	1:A:2:LYS:CD	2.85	0.45
1:B:294:GLN:CG	1:B:295:ASP:N	2.79	0.45
1:B:83:ILE:HD13	1:B:158:PHE:CD2	2.51	0.45
1:A:313:ILE:CA	1:A:314:THR:C	2.85	0.45
1:B:79:ILE:HD12	1:B:79:ILE:HA	1.66	0.45
1:A:34:GLY:CA	1:A:35:VAL:C	2.85	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:119:SER:CB	1:A:181:VAL:HG23	2.39	0.45
1:A:232:ARG:O	1:A:233:GLU:CB	2.58	0.45
1:A:93:PRO:HD2	1:A:96:ILE:CD1	2.47	0.45
1:A:263:LYS:CB	1:A:266:GLN:CD	2.86	0.44
1:A:45:THR:HG22	1:A:179:ALA:HB2	1.98	0.44
1:B:137:TYR:CD2	1:B:137:TYR:N	2.85	0.44
1:A:15:SER:HB3	1:A:19:VAL:HG21	1.98	0.44
1:B:279:LEU:HB2	1:B:305:PHE:CE1	2.51	0.44
1:A:11:ILE:HG22	1:A:15:SER:CB	2.45	0.44
1:A:273:ASP:HA	1:A:276:ILE:HD12	1.98	0.44
1:B:114:PHE:O	1:B:140:ILE:HG12	2.17	0.44
1:B:36:HIS:C	1:B:37:VAL:HG23	2.38	0.44
1:A:104:LEU:HD12	1:A:104:LEU:HA	1.72	0.44
1:A:68:LEU:CG	1:A:69:LYS:N	2.80	0.44
1:A:166:ARG:HG2	1:A:171:PHE:HB2	2.00	0.44
1:A:2:LYS:C	1:A:46:GLU:OE1	2.55	0.44
1:B:44:THR:HG22	1:B:45:THR:N	2.32	0.44
1:A:26:LEU:HD23	1:A:26:LEU:HA	1.76	0.43
1:B:110:TYR:N	1:B:110:TYR:HD2	2.16	0.43
1:A:204:ARG:O	1:A:284:ARG:NH2	2.51	0.43
1:A:34:GLY:HA2	1:A:35:VAL:CB	2.42	0.43
1:B:58:PHE:CD1	1:B:62:LEU:HD11	2.52	0.43
1:B:110:TYR:CD1	1:B:143:LYS:HA	2.47	0.43
1:B:58:PHE:CD2	1:B:62:LEU:CD1	3.00	0.43
1:B:294:GLN:CD	1:B:310:SER:OG	2.57	0.43
1:A:160:GLU:O	1:A:164:ILE:HG12	2.19	0.43
1:A:100:MET:HE3	1:A:100:MET:HB3	1.69	0.43
1:A:107:VAL:HG23	1:A:108:GLY:HA2	2.00	0.43
1:B:151:SER:O	1:B:154:TRP:N	2.52	0.43
1:A:292:LYS:HA	1:A:292:LYS:HD3	1.84	0.43
1:B:33:GLU:HA	1:B:35:VAL:N	2.24	0.43
1:B:19:VAL:CG1	1:B:42:CYS:SG	3.07	0.43
1:A:112:MET:HB2	1:A:184:GLN:HE21	1.84	0.43
1:B:82:THR:HA	1:B:85:THR:OG1	2.19	0.43
1:B:110:TYR:CE1	1:B:143:LYS:CD	3.02	0.43
1:B:107:VAL:C	1:B:184:GLN:NE2	2.72	0.43
1:A:199:PRO:HD2	1:A:200:ILE:H	1.84	0.42
1:B:110:TYR:N	1:B:110:TYR:CD2	2.86	0.42
1:B:193:ILE:O	1:B:298:TRP:NE1	2.46	0.42
1:B:249:TYR:CD2	1:B:259:ARG:NE	2.84	0.42
1:A:119:SER:CB	1:A:179:ALA:O	2.65	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:52:LEU:O	1:A:53:ALA:C	2.56	0.42
1:B:36:HIS:O	1:B:37:VAL:HG23	2.20	0.42
1:A:66:SER:C	1:A:68:LEU:N	2.73	0.42
1:B:15:SER:O	1:B:24:MET:HE1	2.19	0.42
1:B:168:GLN:HB3	1:B:168:GLN:HE21	1.70	0.42
1:A:4:TYR:O	1:A:44:THR:N	2.40	0.42
1:A:129:SER:OG	1:A:130:PHE:N	2.49	0.42
1:B:31:ASN:O	1:B:32:ILE:HG13	2.20	0.41
1:B:295:ASP:HB3	1:B:311:ARG:HG3	2.02	0.41
1:B:14:HIS:CE1	1:B:24:MET:HG2	2.55	0.41
1:B:79:ILE:HA	1:B:82:THR:OG1	2.19	0.41
1:A:93:PRO:HD2	1:A:96:ILE:HD12	2.02	0.41
1:B:220:VAL:HG13	1:B:223:LEU:CD1	2.21	0.41
1:B:2:LYS:HB2	1:B:46:GLU:HG3	2.03	0.41
1:A:198:ASP:OD1	1:A:200:ILE:CG1	2.66	0.41
1:B:157:LEU:HB2	1:B:158:PHE:HD1	1.85	0.41
1:B:168:GLN:H	1:B:168:GLN:HG2	1.73	0.41
1:B:203:ASN:O	1:B:204:ARG:CB	2.68	0.41
1:B:210:ASP:HA	1:B:227:ASP:O	2.21	0.41
1:B:233:GLU:N	1:B:284:ARG:NH2	2.51	0.41
1:B:31:ASN:O	1:B:32:ILE:CG1	2.69	0.41
1:B:101:ASP:O	1:B:105:LEU:CD1	2.69	0.41
1:B:11:ILE:HG21	1:B:27:GLY:HA3	2.02	0.41
1:A:67:SER:OG	1:A:67:SER:O	2.39	0.41
1:B:13:PRO:HB2	1:B:14:HIS:CB	2.28	0.41
1:A:272:THR:HG23	1:A:275:GLN:HE21	1.85	0.41
1:A:93:PRO:HB2	1:A:96:ILE:HG13	2.02	0.41
1:B:284:ARG:O	1:B:287:GLU:N	2.54	0.40
1:B:37:VAL:HG13	1:B:38:PRO:HD2	2.03	0.40
1:A:123:GLU:CA	1:A:123:GLU:OE2	2.70	0.40
1:A:313:ILE:H	1:A:313:ILE:HG13	1.79	0.40
1:B:83:ILE:HD12	1:B:158:PHE:HD2	1.86	0.40
1:A:289:TYR:C	1:A:289:TYR:CD2	2.94	0.40
1:B:141:ILE:H	1:B:141:ILE:HG13	1.69	0.40
1:B:36:HIS:O	1:B:37:VAL:CG2	2.69	0.40
1:B:92:ILE:HA	1:B:93:PRO:HD3	1.85	0.40
1:A:187:SER:HA	1:A:188:PRO:HD3	1.77	0.40
1:A:196:THR:HG22	1:A:207:LEU:CD1	2.50	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	304/327 (93%)	277 (91%)	26 (9%)	1 (0%)	41	76
1	B	281/327 (86%)	259 (92%)	20 (7%)	2 (1%)	22	60
All	All	585/654 (89%)	536 (92%)	46 (8%)	3 (0%)	29	68

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	94	SER
1	A	312	PRO
1	B	188	PRO

### 5.3.2 Protein sidechains [i](#)

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	262/278 (94%)	232 (88%)	30 (12%)	5	24
1	B	247/278 (89%)	218 (88%)	29 (12%)	5	22
All	All	509/556 (92%)	450 (88%)	59 (12%)	5	23

All (59) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	LYS
1	A	7	LYS
1	A	9	GLN

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	A	10	GLU
1	A	12	ARG
1	A	14	HIS
1	A	30	SER
1	A	33	GLU
1	A	38	PRO
1	A	50	ARG
1	A	54	GLU
1	A	56	ASN
1	A	59	THR
1	A	62	LEU
1	A	65	LEU
1	A	68	LEU
1	A	71	SER
1	A	79	ILE
1	A	90	THR
1	A	93	PRO
1	A	94	SER
1	A	123	GLU
1	A	178	LEU
1	A	183	GLN
1	A	235	THR
1	A	250	SER
1	A	252	LYS
1	A	256	THR
1	A	263	LYS
1	A	313	ILE
1	B	10	GLU
1	B	26	LEU
1	B	30	SER
1	B	35	VAL
1	B	50	ARG
1	B	56	ASN
1	B	59	THR
1	B	62	LEU
1	B	77	ARG
1	B	82	THR
1	B	85	THR
1	B	88	GLN
1	B	89	HIS
1	B	103	THR
1	B	112	MET

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Mol	Chain	Res	Type
1	B	119	SER
1	B	143	LYS
1	B	161	ARG
1	B	164	ILE
1	B	168	GLN
1	B	170	GLN
1	B	171	PHE
1	B	174	ARG
1	B	224	VAL
1	B	233	GLU
1	B	265	GLN
1	B	292	LYS
1	B	294	GLN
1	B	311	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	36	HIS
1	A	168	GLN
1	A	184	GLN
1	A	266	GLN
1	A	275	GLN
1	B	36	HIS
1	B	139	ASN
1	B	168	GLN
1	B	184	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 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 [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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	308/327 (94%)	-0.36	1 (0%) 94 84	8, 31, 80, 110	0
1	B	291/327 (88%)	-0.17	6 (2%) 63 34	9, 48, 92, 113	0
All	All	599/654 (91%)	-0.27	7 (1%) 79 54	8, 40, 89, 113	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	14	HIS	4.0
1	B	203	ASN	3.9
1	B	294	GLN	2.9
1	A	14	HIS	2.6
1	B	202	SER	2.2
1	B	289	TYR	2.1
1	B	204	ARG	2.1

### 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

There are no such residues in this entry.