



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:39 AM UTC

PDB ID : 4NZS / pdb_00004nzs
Title : Crystal structure of beta-ketothiolase BktB B from Ralstonia eutropha H16
Authors : Kim, E.J.; Son, H.; Kim, S.; Kim, K.J.
Deposited on : 2013-12-12
Resolution : 2.29 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

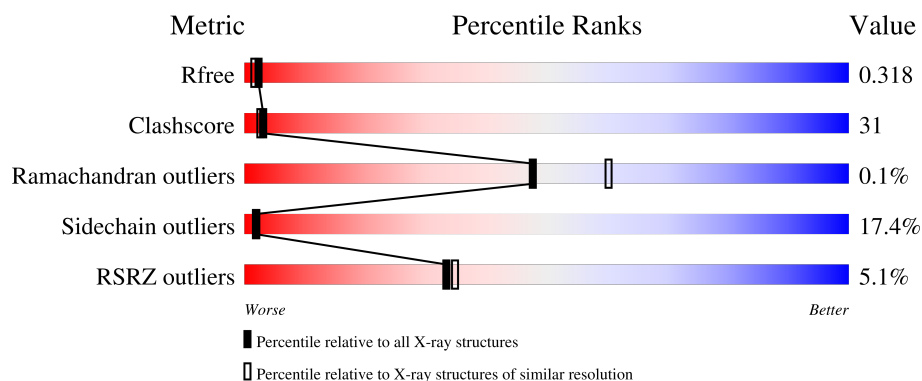
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.29 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (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 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	399	
1	B	399	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5732 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 Beta-ketothiolase BktB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	393	Total	C	N	O	S	0	0	0
			2859	1777	526	542	14			
1	B	393	Total	C	N	O	S	0	0	0
			2859	1777	526	542	14			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	HIS	-	expression tag	UNP Q0KBP1
A	-3	HIS	-	expression tag	UNP Q0KBP1
A	-2	HIS	-	expression tag	UNP Q0KBP1
A	-1	HIS	-	expression tag	UNP Q0KBP1
A	0	HIS	-	expression tag	UNP Q0KBP1
A	1	HIS	-	expression tag	UNP Q0KBP1
B	-4	HIS	-	expression tag	UNP Q0KBP1
B	-3	HIS	-	expression tag	UNP Q0KBP1
B	-2	HIS	-	expression tag	UNP Q0KBP1
B	-1	HIS	-	expression tag	UNP Q0KBP1
B	0	HIS	-	expression tag	UNP Q0KBP1
B	1	HIS	-	expression tag	UNP Q0KBP1

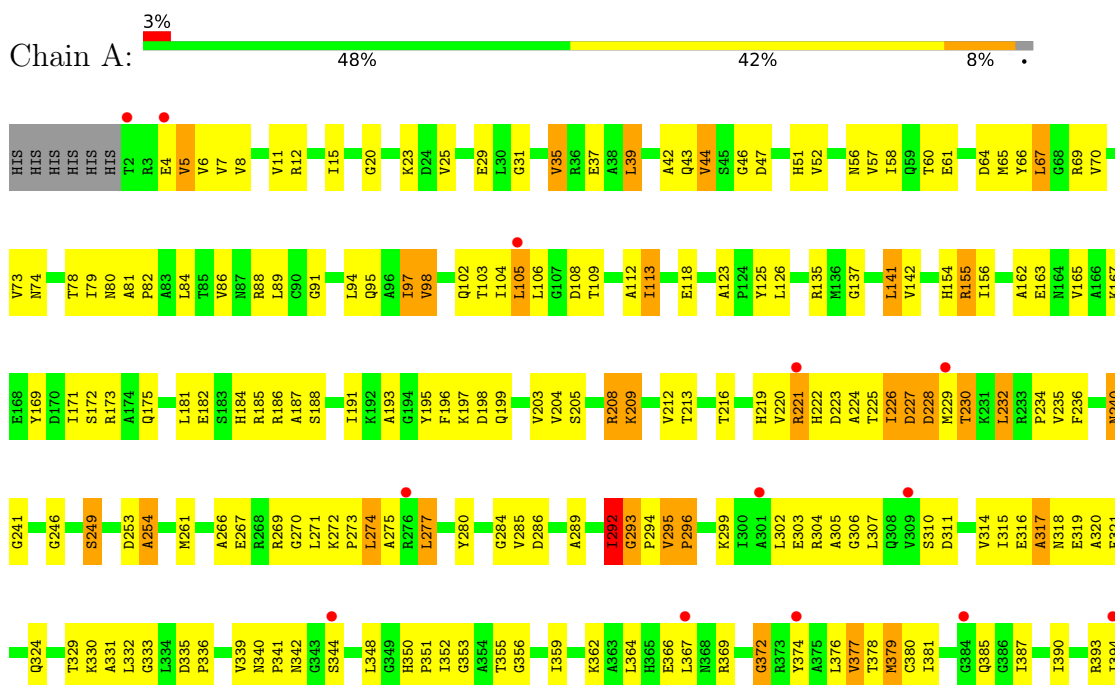
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	3	Total	O	0	0
			3	3		
2	B	11	Total	O	0	0
			11	11		

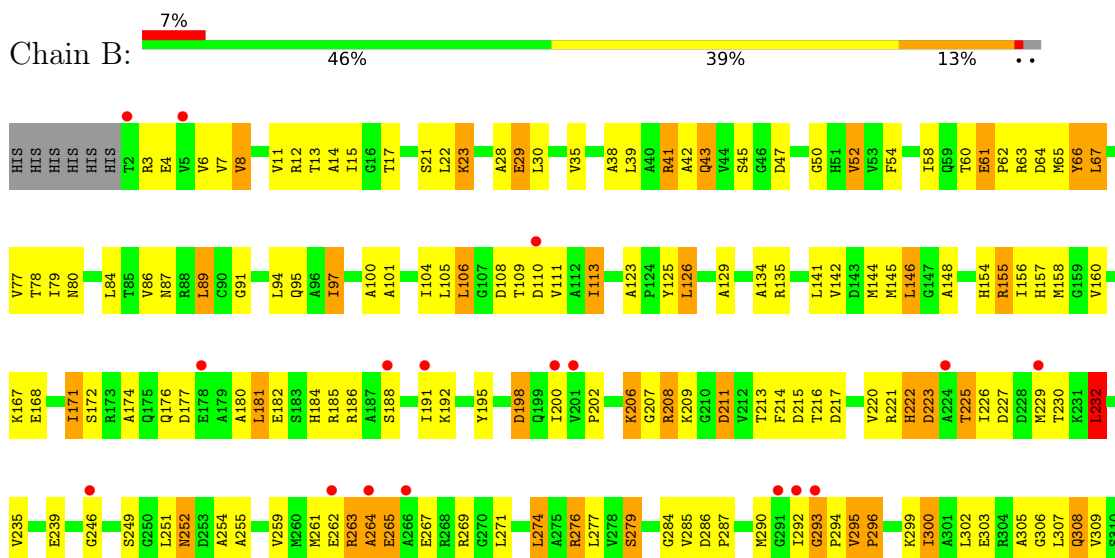
3 Residue-property plots

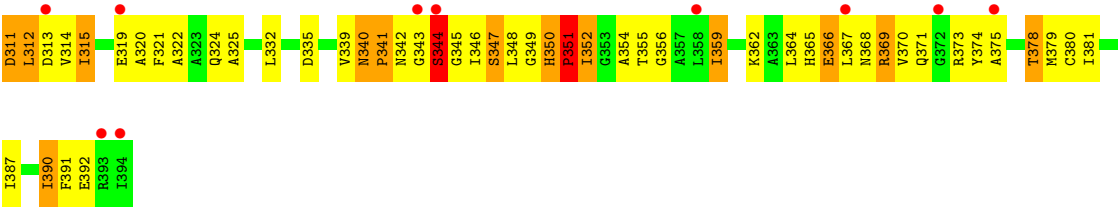
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density.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



• Molecule 1: Beta-ketothiolase BktB





4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	106.95Å 107.24Å 144.14Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	35.74 – 2.29 35.74 – 2.29	Depositor EDS
% Data completeness (in resolution range)	85.4 (35.74-2.29) 85.5 (35.74-2.29)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.66 (at 2.29Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.253 , 0.319 0.251 , 0.318	Depositor DCC
R_{free} test set	1574 reflections (4.21%)	wwPDB-VP
Wilson B-factor (Å ²)	41.7	Xtriage
Anisotropy	0.247	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 46.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	0.042 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5732	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.02% 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	1.06	2/2899 (0.1%)	1.31	30/3930 (0.8%)
1	B	1.04	4/2899 (0.1%)	1.25	25/3930 (0.6%)
All	All	1.05	6/5798 (0.1%)	1.28	55/7860 (0.7%)

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	2
1	B	0	1
All	All	0	3

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	295	VAL	C-N	6.31	1.42	1.34
1	A	295	VAL	C-N	5.81	1.41	1.34
1	B	341	PRO	N-CD	5.36	1.55	1.47
1	A	293	GLY	C-N	5.13	1.41	1.34
1	B	351	PRO	N-CD	5.12	1.54	1.47
1	B	126	LEU	C-O	-5.06	1.17	1.23

All (55) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	293	GLY	O-C-N	11.39	125.17	121.07
1	B	174	ALA	N-CA-C	9.37	121.57	111.36
1	B	225	THR	N-CA-C	9.29	123.27	109.24
1	A	295	VAL	CA-C-N	-8.37	109.77	119.05
1	A	295	VAL	C-N-CA	-8.37	109.77	119.05
1	A	57	VAL	N-CA-C	8.26	118.15	111.62

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	295	VAL	CA-C-N	-8.11	110.05	119.05
1	B	295	VAL	C-N-CA	-8.11	110.05	119.05
1	A	123	ALA	N-CA-C	-7.86	99.23	110.08
1	B	295	VAL	O-C-N	7.59	125.28	120.42
1	B	313	ASP	N-CA-C	7.59	119.63	111.36
1	B	232	LEU	N-CA-C	7.54	120.60	110.35
1	A	295	VAL	O-C-N	7.44	125.18	120.42
1	B	222	HIS	N-CA-C	-7.21	101.54	110.41
1	B	293	GLY	N-CA-C	-7.13	106.16	115.22
1	A	292	ILE	N-CA-C	6.95	118.83	112.29
1	A	67	LEU	N-CA-C	6.95	119.73	111.33
1	A	135	ARG	N-CA-C	-6.72	103.95	111.28
1	A	5	VAL	N-CA-C	6.68	118.41	108.46
1	B	296	PRO	CA-N-CD	-6.66	102.68	112.00
1	A	254	ALA	N-CA-C	6.57	118.12	108.14
1	A	340	ASN	CA-C-N	6.21	125.90	119.82
1	A	340	ASN	C-N-CA	6.21	125.90	119.82
1	A	296	PRO	CA-N-CD	-6.17	103.36	112.00
1	A	372	GLY	N-CA-C	-6.16	102.94	112.85
1	B	200	ILE	CB-CA-C	-6.11	103.93	111.08
1	A	311	ASP	N-CA-C	-6.10	105.42	112.92
1	A	56	ASN	CA-C-N	-6.03	116.77	123.16
1	A	56	ASN	C-N-CA	-6.03	116.77	123.16
1	B	293	GLY	CA-C-N	-5.97	112.42	119.05
1	B	293	GLY	C-N-CA	-5.97	112.42	119.05
1	A	81	ALA	N-CA-C	5.96	118.10	109.04
1	A	137	GLY	N-CA-C	-5.88	106.37	111.95
1	B	264	ALA	N-CA-C	5.86	117.47	111.14
1	B	61	GLU	N-CA-C	-5.85	102.86	108.13
1	A	286	ASP	CA-C-N	5.85	125.05	118.97
1	A	286	ASP	C-N-CA	5.85	125.05	118.97
1	A	293	GLY	CA-C-N	-5.78	112.64	119.05
1	A	293	GLY	C-N-CA	-5.78	112.64	119.05
1	B	52	VAL	N-CA-C	5.75	116.40	108.12
1	B	315	ILE	N-CA-C	5.63	116.16	107.78
1	A	219	HIS	N-CA-C	5.37	117.22	111.36
1	A	35	VAL	N-CA-C	-5.37	105.48	110.53
1	A	98	VAL	CB-CA-C	-5.33	105.05	111.88
1	B	66	TYR	N-CA-C	-5.29	105.27	112.26
1	A	7	VAL	N-CA-C	5.28	115.18	107.37
1	A	318	ASN	N-CA-C	5.25	118.50	110.20
1	B	340	ASN	CA-C-N	-5.25	113.22	119.05

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	340	ASN	C-N-CA	-5.25	113.22	119.05
1	B	252	ASN	N-CA-C	5.22	116.74	108.96
1	A	46	GLY	N-CA-C	-5.10	106.21	113.86
1	A	223	ASP	N-CA-C	-5.07	107.59	112.97
1	B	350	HIS	CA-C-N	-5.04	113.54	119.84
1	B	350	HIS	C-N-CA	-5.04	113.54	119.84
1	B	344	SER	N-CA-C	5.03	115.17	108.23

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	310	SER	Peptide
1	A	317	ALA	Peptide
1	B	351	PRO	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	2859	0	2896	159	0
1	B	2859	0	2896	225	0
2	A	3	0	0	0	0
2	B	11	0	0	2	0
All	All	5732	0	5792	362	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 31.

All (362) 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:181:LEU:CD2	1:B:226:ILE:HG13	1.74	1.18
1:B:181:LEU:HD23	1:B:226:ILE:HG13	1.24	1.09
1:B:308:GLN:OE1	1:B:309:VAL:N	1.90	1.04

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:314:VAL:CG1	1:B:367:LEU:HD13	1.91	1.00
1:B:312:LEU:HD23	1:B:374:TYR:C	1.87	1.00
1:A:91:GLY:HA2	1:A:379:MET:HE3	1.46	0.97
1:B:295:VAL:HG13	1:B:332:LEU:HD21	1.42	0.97
1:B:176:GLN:HB2	1:B:322:ALA:CB	1.97	0.95
1:A:319:GLU:CD	1:A:344:SER:HB3	1.90	0.95
1:B:320:ALA:C	1:B:321:PHE:CD1	2.45	0.93
1:B:263:ARG:HD3	1:B:267:GLU:OE1	1.70	0.92
1:B:314:VAL:HG12	1:B:367:LEU:CD1	2.01	0.90
1:B:314:VAL:CG1	1:B:367:LEU:CD1	2.51	0.89
1:A:42:ALA:HB3	1:A:44:VAL:HG12	1.55	0.88
1:B:176:GLN:HB2	1:B:322:ALA:HB3	1.55	0.87
1:B:227:ASP:O	1:B:230:THR:HG22	1.74	0.87
1:B:356:GLY:HA3	1:B:379:MET:HE1	1.57	0.87
1:A:91:GLY:HA2	1:A:379:MET:CE	2.04	0.87
1:A:292:ILE:O	1:A:296:PRO:HD2	1.73	0.86
1:B:319:GLU:HG2	1:B:325:ALA:CB	2.07	0.84
1:B:319:GLU:OE2	1:B:346:ILE:HD12	1.77	0.84
1:B:292:ILE:O	1:B:296:PRO:HD2	1.77	0.84
1:B:206:LYS:HA	1:B:211:ASP:HB2	1.58	0.83
1:B:261:MET:HE2	1:B:265:GLU:HB3	1.60	0.83
1:B:226:ILE:H	1:B:226:ILE:HD12	1.43	0.82
1:A:285:VAL:CG1	1:A:296:PRO:HG3	2.09	0.82
1:B:306:GLY:O	1:B:307:LEU:HD22	1.79	0.82
1:A:316:GLU:O	1:A:377:VAL:HA	1.81	0.81
1:B:263:ARG:O	1:B:267:GLU:HG2	1.81	0.81
1:B:373:ARG:O	1:B:392:GLU:HA	1.81	0.81
1:B:314:VAL:HG11	1:B:367:LEU:HD13	1.62	0.80
1:B:181:LEU:CD2	1:B:226:ILE:CG1	2.57	0.80
1:B:314:VAL:HG12	1:B:367:LEU:HD13	1.63	0.79
1:B:312:LEU:HD23	1:B:375:ALA:N	1.98	0.79
1:A:42:ALA:HB3	1:A:44:VAL:CG1	2.13	0.79
1:A:369:ARG:HG2	1:A:369:ARG:HH21	1.48	0.78
1:B:188:SER:HB2	1:B:222:HIS:HD2	1.49	0.77
1:B:184:HIS:CD2	1:B:221:ARG:H	2.03	0.77
1:A:66:TYR:CE2	1:B:89:LEU:HD22	2.20	0.76
1:A:261:MET:HE1	1:A:269:ARG:HD2	1.66	0.76
1:A:369:ARG:HG2	1:A:369:ARG:NH2	2.01	0.76
1:B:181:LEU:HD23	1:B:226:ILE:CG1	2.10	0.75
1:B:176:GLN:CB	1:B:322:ALA:HB1	2.17	0.74
1:A:182:GLU:OE2	1:A:186:ARG:NE	2.20	0.74

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:176:GLN:CB	1:B:322:ALA:CB	2.65	0.74
1:B:294:PRO:HB3	1:B:378:THR:OG1	1.87	0.73
1:B:4:GLU:HB3	1:B:263:ARG:HB2	1.71	0.71
1:B:171:ILE:HG22	1:B:176:GLN:HE21	1.55	0.70
1:B:312:LEU:HG	1:B:374:TYR:HB3	1.73	0.70
1:B:8:VAL:HG11	1:B:271:LEU:HD13	1.73	0.70
1:A:84:LEU:HD11	1:B:86:VAL:HG13	1.72	0.69
1:A:106:LEU:CD2	1:B:105:LEU:HD23	2.22	0.69
1:A:266:ALA:O	1:A:270:GLY:N	2.26	0.69
1:B:320:ALA:C	1:B:321:PHE:HD1	1.99	0.69
1:B:320:ALA:HB1	1:B:321:PHE:CE1	2.28	0.69
1:A:285:VAL:HG11	1:A:296:PRO:HG3	1.75	0.69
1:B:186:ARG:NH1	1:B:340:ASN:O	2.26	0.69
1:B:320:ALA:CB	1:B:321:PHE:HD1	2.06	0.68
1:B:41:ARG:NH1	1:B:198:ASP:O	2.25	0.68
1:B:184:HIS:HD2	1:B:221:ARG:H	1.39	0.68
1:B:312:LEU:HD23	1:B:374:TYR:O	1.94	0.67
1:B:306:GLY:C	1:B:307:LEU:HD22	2.19	0.67
1:A:381:ILE:HB	1:A:385:GLN:HB2	1.76	0.67
1:B:314:VAL:HG11	1:B:367:LEU:CD1	2.21	0.66
1:B:305:ALA:O	1:B:307:LEU:HD23	1.96	0.66
1:B:320:ALA:O	1:B:321:PHE:CD1	2.49	0.66
1:A:319:GLU:CG	1:A:344:SER:HB3	2.25	0.66
1:B:308:GLN:CD	1:B:309:VAL:H	2.01	0.66
1:A:94:LEU:HD23	1:A:387:ILE:HG23	1.78	0.65
1:B:356:GLY:CA	1:B:379:MET:HE1	2.27	0.65
1:B:206:LYS:CA	1:B:211:ASP:HB2	2.26	0.65
1:B:292:ILE:C	1:B:292:ILE:HD12	2.22	0.65
1:B:312:LEU:CD2	1:B:375:ALA:N	2.59	0.65
1:A:261:MET:HE1	1:A:269:ARG:CD	2.27	0.65
1:A:226:ILE:O	1:A:230:THR:HG22	1.96	0.64
1:B:60:THR:HB	1:B:64:ASP:OD2	1.96	0.64
1:B:320:ALA:CB	1:B:321:PHE:CD1	2.80	0.64
1:B:342:ASN:HD21	1:B:366:GLU:HG3	1.63	0.64
1:A:195:TYR:CE2	1:A:369:ARG:NE	2.66	0.63
1:B:320:ALA:HB1	1:B:321:PHE:CD1	2.34	0.63
1:A:58:ILE:HG12	1:B:65:MET:HE1	1.81	0.63
1:B:319:GLU:HG2	1:B:325:ALA:HB1	1.78	0.62
1:B:311:ASP:N	1:B:311:ASP:OD1	2.30	0.62
1:B:350:HIS:CE1	1:B:352:ILE:HG22	2.35	0.62
1:A:191:ILE:CD1	1:A:220:VAL:HG11	2.29	0.62

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:182:GLU:O	1:B:186:ARG:HG3	2.00	0.62
1:B:302:LEU:O	1:B:306:GLY:N	2.33	0.62
1:B:319:GLU:HB2	1:B:344:SER:HB3	1.82	0.62
1:B:279:SER:OG	1:B:390:ILE:HD11	1.99	0.62
1:B:64:ASP:O	1:B:67:LEU:HB2	2.00	0.62
1:B:186:ARG:NH1	1:B:341:PRO:HA	2.14	0.61
1:B:91:GLY:HA2	1:B:379:MET:HE3	1.81	0.61
1:B:186:ARG:HH11	1:B:341:PRO:HA	1.66	0.61
1:B:335:ASP:O	1:B:339:VAL:HG23	2.00	0.60
1:A:193:ALA:HB1	1:A:195:TYR:CE1	2.37	0.60
1:B:11:VAL:HG12	1:B:38:ALA:HB2	1.84	0.60
1:A:188:SER:HB2	1:A:222:HIS:ND1	2.17	0.59
1:A:42:ALA:CB	1:A:44:VAL:HG12	2.32	0.59
1:A:225:THR:HB	1:A:228:ASP:HB2	1.84	0.59
1:B:345:GLY:HA2	1:B:349:GLY:O	2.02	0.59
1:A:306:GLY:C	1:A:307:LEU:HD23	2.28	0.59
1:A:319:GLU:CD	1:A:344:SER:CB	2.74	0.58
1:A:249:SER:HB2	1:A:320:ALA:O	2.03	0.58
1:B:154:HIS:O	1:B:155:ARG:HG2	2.04	0.57
1:B:249:SER:HB2	1:B:320:ALA:O	2.03	0.57
1:B:207:GLY:HA3	2:B:411:HOH:O	2.04	0.57
1:B:312:LEU:HD23	1:B:375:ALA:CA	2.35	0.57
1:A:5:VAL:HG23	1:A:277:LEU:HB2	1.85	0.57
1:A:285:VAL:HG12	1:A:296:PRO:HG3	1.86	0.56
1:B:176:GLN:CB	1:B:322:ALA:HB3	2.32	0.56
1:B:184:HIS:HA	1:B:347:SER:OG	2.05	0.56
1:B:350:HIS:HD2	1:B:355:THR:OG1	1.88	0.56
1:A:94:LEU:HB2	1:A:379:MET:HE1	1.88	0.56
1:B:302:LEU:HB3	1:B:307:LEU:O	2.05	0.56
1:A:305:ALA:HB2	1:A:390:ILE:HD11	1.85	0.56
1:B:7:VAL:HG21	1:B:364:LEU:HD21	1.88	0.56
1:B:191:ILE:HD11	1:B:220:VAL:HG21	1.88	0.56
1:A:284:GLY:O	1:B:80:ASN:HA	2.05	0.55
1:B:188:SER:CB	1:B:222:HIS:HD2	2.19	0.55
1:B:22:LEU:O	1:B:23:LYS:C	2.49	0.55
1:B:321:PHE:CD1	1:B:321:PHE:N	2.74	0.55
1:B:321:PHE:H	1:B:324:GLN:HE21	1.53	0.55
1:A:369:ARG:HH21	1:A:369:ARG:CG	2.14	0.55
1:B:274:LEU:H	1:B:274:LEU:HD12	1.72	0.55
1:B:94:LEU:HB2	1:B:379:MET:CE	2.37	0.55
1:B:172:SER:O	1:B:176:GLN:HG2	2.07	0.55

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:262:GLU:O	1:B:263:ARG:C	2.49	0.55
1:B:312:LEU:CD2	1:B:375:ALA:CA	2.84	0.55
1:B:181:LEU:HD21	1:B:226:ILE:N	2.21	0.55
1:B:319:GLU:HG2	1:B:325:ALA:HB2	1.90	0.54
1:A:294:PRO:HD3	1:A:380:CYS:HB3	1.90	0.54
1:B:14:ALA:HB3	1:B:214:PHE:CE2	2.43	0.54
1:B:345:GLY:HA2	1:B:349:GLY:C	2.32	0.54
1:A:321:PHE:O	1:A:324:GLN:HG3	2.07	0.54
1:B:202:PRO:HB3	1:B:215:ASP:HB3	1.90	0.54
1:A:66:TYR:CE2	1:B:89:LEU:CD2	2.91	0.54
1:B:104:ILE:HA	1:B:109:THR:O	2.08	0.53
1:A:126:LEU:HD23	1:B:126:LEU:HD22	1.88	0.53
1:A:69:ARG:O	1:A:73:VAL:HG23	2.08	0.53
1:B:232:LEU:HD13	1:B:246:GLY:HA3	1.89	0.53
1:A:296:PRO:HA	1:A:299:LYS:HE3	1.91	0.53
1:B:171:ILE:CG2	1:B:176:GLN:HE21	2.20	0.53
1:B:254:ALA:HB3	1:B:351:PRO:HG3	1.89	0.53
1:A:209:LYS:HD2	1:A:209:LYS:O	2.09	0.53
1:A:186:ARG:NH1	1:A:339:VAL:O	2.42	0.53
1:A:66:TYR:CZ	1:B:89:LEU:HD22	2.44	0.52
1:A:86:VAL:HG13	1:B:84:LEU:HD11	1.90	0.52
1:A:88:ARG:O	1:A:381:ILE:HG23	2.10	0.52
1:A:163:GLU:HA	1:A:163:GLU:OE1	2.08	0.52
1:B:186:ARG:HD2	1:B:341:PRO:O	2.10	0.52
1:B:249:SER:OG	1:B:350:HIS:HB2	2.09	0.52
1:B:350:HIS:NE2	1:B:352:ILE:HG22	2.24	0.52
1:A:118:GLU:OE2	1:A:353:GLY:N	2.42	0.52
1:B:351:PRO:O	1:B:351:PRO:HG2	2.10	0.52
1:A:47:ASP:HA	1:A:78:THR:HG23	1.92	0.52
1:B:12:ARG:NH1	1:B:13:THR:O	2.43	0.52
1:A:293:GLY:N	1:A:294:PRO:CD	2.73	0.52
1:B:181:LEU:CD2	1:B:226:ILE:N	2.73	0.52
1:A:52:VAL:HG13	1:A:113:ILE:HG22	1.92	0.51
1:B:195:TYR:CE1	1:B:369:ARG:HD2	2.44	0.51
1:B:295:VAL:HG12	1:B:299:LYS:HE2	1.92	0.51
1:A:104:ILE:HD11	1:A:112:ALA:HB3	1.90	0.51
1:A:267:GLU:O	1:A:270:GLY:N	2.43	0.51
1:B:15:ILE:CD1	1:B:348:LEU:HB3	2.41	0.51
1:B:217:ASP:O	1:B:220:VAL:HG23	2.10	0.51
1:B:232:LEU:CD1	1:B:246:GLY:HA3	2.40	0.51
1:A:227:ASP:HA	1:A:230:THR:HG23	1.93	0.51

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:312:LEU:CD2	1:B:374:TYR:C	2.73	0.51
1:B:188:SER:HB2	1:B:222:HIS:CD2	2.38	0.51
1:A:8:VAL:O	1:A:274:LEU:HD21	2.11	0.51
1:A:82:PRO:HD3	1:B:284:GLY:N	2.26	0.51
1:B:188:SER:CB	1:B:222:HIS:CD2	2.94	0.51
1:B:355:THR:O	1:B:359:ILE:HG12	2.10	0.51
1:B:296:PRO:O	1:B:300:ILE:HD12	2.11	0.51
1:A:274:LEU:HD23	1:A:274:LEU:N	2.26	0.50
1:A:125:TYR:HA	1:A:142:VAL:O	2.11	0.50
1:A:285:VAL:HG11	1:A:296:PRO:CG	2.41	0.50
1:A:359:ILE:HA	1:A:362:LYS:HD3	1.93	0.50
1:B:217:ASP:CB	1:B:348:LEU:HD23	2.42	0.50
1:A:64:ASP:HA	1:A:67:LEU:HB2	1.94	0.50
1:B:308:GLN:OE1	1:B:309:VAL:HG23	2.12	0.50
1:A:69:ARG:NH2	1:B:285:VAL:O	2.45	0.50
1:B:249:SER:OG	1:B:350:HIS:CB	2.60	0.50
1:A:70:VAL:O	1:A:74:ASN:HB2	2.12	0.49
1:B:42:ALA:O	1:B:43:GLN:HB2	2.12	0.49
1:A:172:SER:O	1:A:173:ARG:C	2.54	0.49
1:B:345:GLY:O	1:B:349:GLY:HA2	2.12	0.49
1:A:169:TYR:HB2	1:A:171:ILE:CD1	2.41	0.49
1:A:204:VAL:O	1:A:204:VAL:HG23	2.12	0.49
1:B:157:HIS:ND1	1:B:158:MET:N	2.60	0.49
1:B:50:GLY:N	1:B:111:VAL:O	2.34	0.49
1:A:78:THR:HB	1:A:80:ASN:H	1.77	0.49
1:B:362:LYS:O	1:B:366:GLU:HB2	2.11	0.49
1:B:7:VAL:O	1:B:274:LEU:HD12	2.11	0.49
1:B:367:LEU:HD13	1:B:375:ALA:HB2	1.94	0.49
1:B:176:GLN:CA	1:B:322:ALA:HB1	2.43	0.49
1:B:290:MET:HB2	1:B:381:ILE:O	2.13	0.49
1:B:123:ALA:CB	1:B:145:MET:HB2	2.44	0.48
1:A:229:MET:HA	1:A:232:LEU:HD22	1.95	0.48
1:B:286:ASP:O	1:B:287:PRO:C	2.55	0.48
1:A:91:GLY:HA2	1:A:379:MET:HE2	1.93	0.48
1:B:306:GLY:C	1:B:307:LEU:CD2	2.87	0.48
1:A:8:VAL:HG11	1:A:271:LEU:HD13	1.96	0.48
1:A:216:THR:HG23	1:A:216:THR:O	2.14	0.48
1:B:158:MET:HE2	1:B:158:MET:HA	1.96	0.48
1:A:66:TYR:OH	1:B:148:ALA:O	2.20	0.48
1:B:181:LEU:HD22	1:B:226:ILE:HA	1.96	0.48
1:A:60:THR:HB	1:A:64:ASP:OD2	2.14	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:186:ARG:HD3	1:A:341:PRO:HA	1.96	0.48
1:B:186:ARG:HH11	1:B:341:PRO:CA	2.26	0.48
1:A:350:HIS:HD2	1:A:355:THR:OG1	1.97	0.48
1:B:6:VAL:HG12	1:B:276:ARG:CB	2.43	0.48
1:B:293:GLY:N	1:B:294:PRO:CD	2.77	0.48
1:B:320:ALA:HB1	1:B:321:PHE:HE1	1.77	0.48
1:A:181:LEU:CD1	1:A:224:ALA:HB1	2.44	0.47
1:A:6:VAL:HG22	1:A:261:MET:O	2.14	0.47
1:A:307:LEU:HD13	1:A:374:TYR:CE2	2.48	0.47
1:A:39:LEU:HD12	1:A:44:VAL:HG13	1.95	0.47
1:B:176:GLN:HB2	1:B:322:ALA:HB1	1.77	0.47
1:B:29:GLU:CD	1:B:63:ARG:HH22	2.22	0.47
1:B:222:HIS:HB3	1:B:223:ASP:OD1	2.14	0.47
1:A:169:TYR:CD1	1:A:169:TYR:N	2.82	0.47
1:A:187:ALA:HA	1:A:342:ASN:O	2.14	0.47
1:A:304:ARG:NH1	1:B:108:ASP:OD1	2.48	0.47
1:B:35:VAL:HG22	1:B:113:ILE:HD11	1.97	0.47
1:B:125:TYR:HA	1:B:142:VAL:O	2.15	0.47
1:B:223:ASP:OD1	1:B:223:ASP:N	2.47	0.47
1:B:312:LEU:CD2	1:B:375:ALA:C	2.87	0.47
1:B:86:VAL:HG11	1:B:95:GLN:HG2	1.97	0.47
1:A:31:GLY:O	1:A:35:VAL:HG23	2.15	0.47
1:A:274:LEU:O	1:A:275:ALA:HB2	2.15	0.47
1:B:156:ILE:HD12	1:B:160:VAL:HB	1.96	0.47
1:A:103:THR:HB	1:A:108:ASP:HB2	1.95	0.46
1:B:181:LEU:CD2	1:B:226:ILE:CD1	2.93	0.46
1:B:351:PRO:HG2	1:B:354:ALA:HB3	1.96	0.46
1:A:88:ARG:O	1:A:381:ILE:CG2	2.64	0.46
1:A:5:VAL:HG13	1:A:105:LEU:CD1	2.46	0.46
1:A:20:GLY:O	1:A:23:LYS:HB2	2.15	0.46
1:B:176:GLN:HA	1:B:322:ALA:HB1	1.98	0.46
1:A:88:ARG:CD	1:A:381:ILE:HD13	2.45	0.46
1:A:195:TYR:CE2	1:A:369:ARG:CD	2.98	0.46
1:A:235:VAL:HG23	1:A:236:PHE:CE2	2.50	0.46
1:A:240:ASN:O	1:A:240:ASN:ND2	2.40	0.46
1:B:110:ASP:O	1:B:261:MET:HA	2.15	0.46
1:B:181:LEU:HD21	1:B:226:ILE:HD12	1.98	0.46
1:B:181:LEU:HD22	1:B:226:ILE:HG13	1.81	0.46
1:A:182:GLU:O	1:A:185:ARG:N	2.48	0.46
1:B:181:LEU:HD21	1:B:226:ILE:CD1	2.45	0.46
1:B:186:ARG:CD	1:B:341:PRO:O	2.64	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:106:LEU:HD11	1:B:106:LEU:HD21	1.97	0.45
1:A:37:GLU:OE1	1:A:37:GLU:HA	2.15	0.45
1:B:206:LYS:HB3	1:B:211:ASP:HB2	1.98	0.45
1:B:339:VAL:O	1:B:340:ASN:HB2	2.16	0.45
1:B:365:HIS:O	1:B:368:ASN:HB2	2.16	0.45
1:A:187:ALA:O	1:A:191:ILE:HG13	2.17	0.45
1:A:302:LEU:O	1:A:303:GLU:C	2.60	0.45
1:B:276:ARG:O	1:B:391:PHE:HB3	2.17	0.45
1:B:180:ALA:HB3	1:B:229:MET:HE3	1.98	0.45
1:B:374:TYR:CD1	1:B:392:GLU:HG3	2.52	0.45
1:A:267:GLU:C	1:A:270:GLY:H	2.24	0.45
1:A:292:ILE:H	1:A:292:ILE:HG13	1.56	0.45
1:B:207:GLY:H	1:B:211:ASP:HA	1.81	0.45
1:A:61:GLU:C	1:B:144:MET:HE1	2.41	0.45
1:B:314:VAL:CG1	1:B:367:LEU:HD12	2.41	0.45
1:B:94:LEU:O	1:B:95:GLN:C	2.58	0.45
1:B:111:VAL:HG23	2:B:401:HOH:O	2.17	0.45
1:B:269:ARG:HB3	1:B:271:LEU:HD12	1.99	0.45
1:B:344:SER:O	1:B:347:SER:N	2.48	0.45
1:B:66:TYR:O	1:B:67:LEU:C	2.60	0.44
1:A:332:LEU:O	1:A:333:GLY:C	2.58	0.44
1:A:66:TYR:CD2	1:B:89:LEU:HD21	2.52	0.44
1:A:184:HIS:O	1:A:185:ARG:C	2.60	0.44
1:A:235:VAL:HG23	1:A:236:PHE:CD2	2.52	0.44
1:A:102:GLN:O	1:A:106:LEU:HD12	2.17	0.44
1:A:317:ALA:HA	1:A:378:THR:O	2.18	0.44
1:B:208:ARG:HG3	1:B:209:LYS:HG2	1.99	0.44
1:A:65:MET:CE	1:B:58:ILE:HG12	2.48	0.44
1:A:227:ASP:OD1	1:A:227:ASP:N	2.51	0.44
1:B:77:VAL:HG12	1:B:78:THR:N	2.33	0.44
1:B:167:LYS:O	1:B:168:GLU:C	2.61	0.44
1:B:263:ARG:HH11	1:B:263:ARG:CG	2.30	0.44
1:A:285:VAL:CG1	1:A:296:PRO:CG	2.90	0.44
1:B:295:VAL:HG13	1:B:332:LEU:CD2	2.29	0.44
1:B:308:GLN:H	1:B:311:ASP:CG	2.26	0.44
1:B:343:GLY:O	1:B:362:LYS:NZ	2.51	0.44
1:B:181:LEU:HD22	1:B:226:ILE:CA	2.47	0.44
1:A:280:TYR:CD1	1:A:280:TYR:C	2.96	0.43
1:A:305:ALA:CB	1:A:390:ILE:HD11	2.46	0.43
1:A:193:ALA:CB	1:A:195:TYR:CD1	3.01	0.43
1:B:89:LEU:HB2	1:B:381:ILE:HG23	1.99	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:227:ASP:HA	1:B:230:THR:HG22	1.98	0.43
1:B:251:LEU:C	1:B:252:ASN:ND2	2.76	0.43
1:A:165:VAL:CG2	1:A:289:ALA:HB1	2.49	0.43
1:A:221:ARG:HB2	1:A:221:ARG:HH11	1.82	0.43
1:B:262:GLU:C	1:B:264:ALA:N	2.76	0.43
1:A:169:TYR:HB2	1:A:171:ILE:HD12	2.01	0.43
1:B:94:LEU:HB3	1:B:387:ILE:HD13	2.00	0.43
1:A:65:MET:O	1:B:87:ASN:ND2	2.52	0.43
1:A:141:LEU:HD12	1:A:141:LEU:HA	1.82	0.43
1:A:221:ARG:HH11	1:A:221:ARG:CB	2.31	0.43
1:A:253:ASP:O	1:A:254:ALA:HB2	2.19	0.43
1:A:285:VAL:HG11	1:A:293:GLY:HA2	2.01	0.43
1:A:366:GLU:O	1:A:367:LEU:C	2.62	0.43
1:B:262:GLU:O	1:B:264:ALA:N	2.51	0.43
1:A:89:LEU:HB2	1:A:381:ILE:CG2	2.48	0.43
1:B:227:ASP:HA	1:B:230:THR:CG2	2.48	0.43
1:A:225:THR:HG22	1:A:227:ASP:H	1.84	0.43
1:B:362:LYS:HB2	1:B:362:LYS:HE3	1.95	0.42
1:A:292:ILE:C	1:A:294:PRO:CD	2.92	0.42
1:A:66:TYR:CD2	1:B:89:LEU:CD2	3.02	0.42
1:A:393:ARG:O	1:A:394:ILE:O	2.37	0.42
1:B:30:LEU:HB3	1:B:255:ALA:HB2	2.00	0.42
1:A:294:PRO:HG3	1:A:379:MET:C	2.45	0.42
1:B:104:ILE:HG23	1:B:110:ASP:HA	2.00	0.42
1:A:356:GLY:HA3	1:A:379:MET:HE1	2.02	0.42
1:A:126:LEU:HD23	1:B:126:LEU:CD2	2.50	0.42
1:A:269:ARG:HB3	1:A:271:LEU:HG	2.02	0.42
1:B:177:ASP:CG	1:B:229:MET:HB3	2.45	0.42
1:B:252:ASN:C	1:B:351:PRO:HB3	2.45	0.42
1:A:193:ALA:HB3	1:A:195:TYR:CD1	2.54	0.42
1:A:12:ARG:HE	1:A:199:GLN:HE21	1.67	0.42
1:A:97:ILE:HG22	1:A:98:VAL:N	2.34	0.42
1:B:39:LEU:HD12	1:B:259:VAL:HG21	2.02	0.42
1:B:314:VAL:C	1:B:315:ILE:HG13	2.45	0.42
1:A:5:VAL:HG13	1:A:105:LEU:HD12	2.02	0.41
1:A:351:PRO:O	1:A:352:ILE:C	2.62	0.41
1:B:380:CYS:C	1:B:381:ILE:HG13	2.44	0.41
1:A:232:LEU:HD23	1:A:246:GLY:HA3	2.01	0.41
1:A:372:GLY:O	1:A:393:ARG:HD3	2.20	0.41
1:B:8:VAL:C	1:B:274:LEU:HD11	2.45	0.41
1:A:106:LEU:HD11	1:B:106:LEU:CD2	2.50	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:315:ILE:HG12	1:A:376:LEU:HB3	2.02	0.41
1:A:330:LYS:O	1:A:331:ALA:C	2.63	0.41
1:A:154:HIS:C	1:A:155:ARG:CG	2.93	0.41
1:B:100:ALA:O	1:B:101:ALA:C	2.63	0.41
1:B:146:LEU:HD13	1:B:146:LEU:HA	1.90	0.41
1:B:320:ALA:HB3	1:B:321:PHE:HD1	1.81	0.41
1:A:95:GLN:O	1:A:95:GLN:NE2	2.51	0.41
1:B:129:ALA:HB1	1:B:134:ALA:HB2	2.02	0.41
1:A:5:VAL:HG11	1:A:104:ILE:HB	2.02	0.41
1:A:181:LEU:HD12	1:A:224:ALA:HB1	2.02	0.41
1:B:367:LEU:O	1:B:371:GLN:N	2.53	0.41
1:A:193:ALA:CB	1:A:195:TYR:CE1	3.04	0.41
1:A:234:PRO:CB	1:A:241:GLY:HA3	2.51	0.41
1:B:180:ALA:CB	1:B:229:MET:HE3	2.51	0.41
1:B:254:ALA:O	1:B:354:ALA:HB2	2.21	0.41
1:A:162:ALA:O	1:A:163:GLU:C	2.63	0.41
1:A:191:ILE:HD11	1:A:220:VAL:HG11	2.01	0.41
1:A:196:PHE:O	1:A:197:LYS:C	2.62	0.41
1:A:208:ARG:HG2	1:A:209:LYS:N	2.33	0.41
1:A:295:VAL:HG22	1:A:332:LEU:HG	2.04	0.41
1:A:335:ASP:O	1:A:336:PRO:C	2.63	0.41
1:B:206:LYS:CB	1:B:211:ASP:HB2	2.51	0.41
1:A:88:ARG:HD2	1:A:381:ILE:HD13	2.02	0.40
1:B:61:GLU:OE1	1:B:63:ARG:HD3	2.21	0.40
1:B:28:ALA:CB	1:B:63:ARG:NH2	2.84	0.40
1:B:94:LEU:O	1:B:97:ILE:N	2.54	0.40
1:A:15:ILE:HD12	1:A:348:LEU:HB3	2.02	0.40
1:B:47:ASP:HA	1:B:78:THR:HB	2.04	0.40
1:A:5:VAL:HG22	1:A:105:LEU:HD11	2.02	0.40
1:A:126:LEU:CD2	1:B:126:LEU:CD2	3.00	0.40
1:B:292:ILE:C	1:B:292:ILE:CD1	2.91	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	391/399 (98%)	356 (91%)	34 (9%)	1 (0%)	36	46
1	B	391/399 (98%)	363 (93%)	28 (7%)	0	100	100
All	All	782/798 (98%)	719 (92%)	62 (8%)	1 (0%)	48	60

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	273	PRO

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	287/293 (98%)	245 (85%)	42 (15%)	3	3
1	B	287/293 (98%)	229 (80%)	58 (20%)	1	1
All	All	574/586 (98%)	474 (83%)	100 (17%)	2	2

All (100) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	GLU
1	A	11	VAL
1	A	25	VAL
1	A	29	GLU
1	A	39	LEU
1	A	43	GLN
1	A	44	VAL
1	A	51	HIS
1	A	79	ILE
1	A	97	ILE
1	A	105	LEU
1	A	109	THR

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	113	ILE
1	A	141	LEU
1	A	155	ARG
1	A	156	ILE
1	A	167	LYS
1	A	175	GLN
1	A	198	ASP
1	A	203	VAL
1	A	205	SER
1	A	208	ARG
1	A	209	LYS
1	A	212	VAL
1	A	213	THR
1	A	221	ARG
1	A	226	ILE
1	A	227	ASP
1	A	228	ASP
1	A	230	THR
1	A	232	LEU
1	A	240	ASN
1	A	249	SER
1	A	272	LYS
1	A	274	LEU
1	A	277	LEU
1	A	292	ILE
1	A	314	VAL
1	A	329	THR
1	A	364	LEU
1	A	377	VAL
1	A	379	MET
1	B	3	ARG
1	B	8	VAL
1	B	17	THR
1	B	21	SER
1	B	23	LYS
1	B	29	GLU
1	B	41	ARG
1	B	43	GLN
1	B	45	SER
1	B	52	VAL
1	B	54	PHE
1	B	62	PRO

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	67	LEU
1	B	79	ILE
1	B	89	LEU
1	B	97	ILE
1	B	106	LEU
1	B	113	ILE
1	B	135	ARG
1	B	141	LEU
1	B	146	LEU
1	B	155	ARG
1	B	171	ILE
1	B	181	LEU
1	B	185	ARG
1	B	192	LYS
1	B	198	ASP
1	B	206	LYS
1	B	208	ARG
1	B	211	ASP
1	B	213	THR
1	B	216	THR
1	B	223	ASP
1	B	225	THR
1	B	232	LEU
1	B	235	VAL
1	B	239	GLU
1	B	263	ARG
1	B	265	GLU
1	B	274	LEU
1	B	276	ARG
1	B	277	LEU
1	B	279	SER
1	B	300	ILE
1	B	303	GLU
1	B	308	GLN
1	B	311	ASP
1	B	312	LEU
1	B	344	SER
1	B	347	SER
1	B	351	PRO
1	B	352	ILE
1	B	359	ILE
1	B	366	GLU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	369	ARG
1	B	370	VAL
1	B	378	THR
1	B	390	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	51	HIS
1	A	59	GLN
1	A	80	ASN
1	A	102	GLN
1	A	175	GLN
1	A	199	GLN
1	A	252	ASN
1	A	308	GLN
1	A	342	ASN
1	A	350	HIS
1	B	59	GLN
1	B	154	HIS
1	B	176	GLN
1	B	184	HIS
1	B	222	HIS
1	B	252	ASN
1	B	340	ASN
1	B	342	ASN
1	B	350	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	393/399 (98%)	0.38	13 (3%) 49 51	21, 52, 80, 116	0
1	B	393/399 (98%)	0.63	27 (6%) 23 25	20, 62, 93, 110	0
All	All	786/798 (98%)	0.50	40 (5%) 33 35	20, 57, 89, 116	0

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	319	GLU	6.8
1	B	394	ILE	4.5
1	B	344	SER	4.2
1	B	262	GLU	3.8
1	B	343	GLY	3.7
1	B	291	GLY	3.4
1	A	367	LEU	3.4
1	B	229	MET	3.2
1	B	367	LEU	3.1
1	B	5	VAL	3.0
1	A	276	ARG	3.0
1	B	375	ALA	2.9
1	A	2	THR	2.8
1	B	266	ALA	2.8
1	B	224	ALA	2.7
1	B	2	THR	2.6
1	B	393	ARG	2.5
1	A	394	ILE	2.5
1	B	178	GLU	2.4
1	B	358	LEU	2.4
1	B	246	GLY	2.4
1	A	4	GLU	2.4
1	A	344	SER	2.4
1	A	221	ARG	2.4

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	384	GLY	2.3
1	B	201	VAL	2.3
1	A	374	TYR	2.2
1	B	191	ILE	2.2
1	A	309	VAL	2.2
1	B	313	ASP	2.2
1	B	264	ALA	2.2
1	B	372	GLY	2.2
1	A	301	ALA	2.2
1	B	293	GLY	2.1
1	B	200	ILE	2.1
1	B	110	ASP	2.0
1	B	292	ILE	2.0
1	A	229	MET	2.0
1	B	188	SER	2.0
1	A	105	LEU	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 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.