



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 08:52 PM UTC

PDB ID : 4WYS / pdb_00004wys
Title : Crystal structure of thiolase from Escherichia coli
Authors : Kim, S.; Ha, S.C.; Ahn, J.W.; Kim, E.J.; Lim, J.H.; Kim, K.J.
Deposited on : 2014-11-18
Resolution : 2.10 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

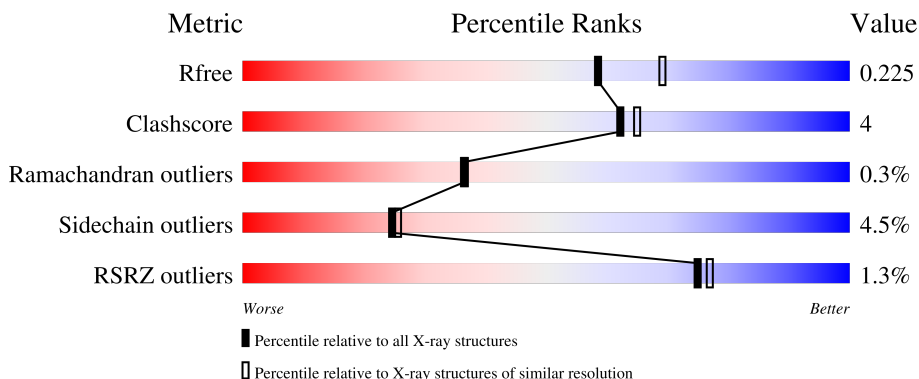
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	405	2% 85% 12% ..
1	B	405	86% 11% ..
1	C	405	% 85% 11% ..
1	D	405	% 86% 10% ..

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 12024 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 Acetyl-CoA acetyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	396	Total	C	N	O	S	0	0	0
			2833	1772	501	542	18			
1	B	396	Total	C	N	O	S	0	0	0
			2833	1772	501	542	18			
1	C	396	Total	C	N	O	S	0	0	0
			2833	1772	501	542	18			
1	D	396	Total	C	N	O	S	0	0	0
			2833	1772	501	542	18			

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	MET	-	initiating methionine	UNP P76461
A	-1	GLY	-	expression tag	UNP P76461
A	0	ALA	-	expression tag	UNP P76461
A	394	ASN	-	expression tag	UNP P76461
A	395	LEU	-	expression tag	UNP P76461
A	396	GLU	-	expression tag	UNP P76461
A	397	HIS	-	expression tag	UNP P76461
A	398	HIS	-	expression tag	UNP P76461
A	399	HIS	-	expression tag	UNP P76461
A	400	HIS	-	expression tag	UNP P76461
A	401	HIS	-	expression tag	UNP P76461
A	402	HIS	-	expression tag	UNP P76461
B	-2	MET	-	initiating methionine	UNP P76461
B	-1	GLY	-	expression tag	UNP P76461
B	0	ALA	-	expression tag	UNP P76461
B	394	ASN	-	expression tag	UNP P76461
B	395	LEU	-	expression tag	UNP P76461
B	396	GLU	-	expression tag	UNP P76461
B	397	HIS	-	expression tag	UNP P76461
B	398	HIS	-	expression tag	UNP P76461
B	399	HIS	-	expression tag	UNP P76461

Continued on next page...

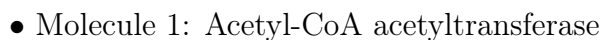
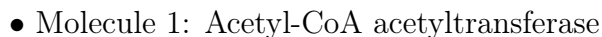
Continued from previous page...

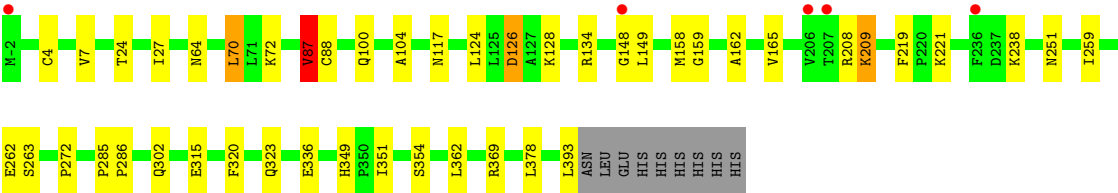
Chain	Residue	Modelled	Actual	Comment	Reference
B	400	HIS	-	expression tag	UNP P76461
B	401	HIS	-	expression tag	UNP P76461
B	402	HIS	-	expression tag	UNP P76461
C	-2	MET	-	initiating methionine	UNP P76461
C	-1	GLY	-	expression tag	UNP P76461
C	0	ALA	-	expression tag	UNP P76461
C	394	ASN	-	expression tag	UNP P76461
C	395	LEU	-	expression tag	UNP P76461
C	396	GLU	-	expression tag	UNP P76461
C	397	HIS	-	expression tag	UNP P76461
C	398	HIS	-	expression tag	UNP P76461
C	399	HIS	-	expression tag	UNP P76461
C	400	HIS	-	expression tag	UNP P76461
C	401	HIS	-	expression tag	UNP P76461
C	402	HIS	-	expression tag	UNP P76461
D	-2	MET	-	initiating methionine	UNP P76461
D	-1	GLY	-	expression tag	UNP P76461
D	0	ALA	-	expression tag	UNP P76461
D	394	ASN	-	expression tag	UNP P76461
D	395	LEU	-	expression tag	UNP P76461
D	396	GLU	-	expression tag	UNP P76461
D	397	HIS	-	expression tag	UNP P76461
D	398	HIS	-	expression tag	UNP P76461
D	399	HIS	-	expression tag	UNP P76461
D	400	HIS	-	expression tag	UNP P76461
D	401	HIS	-	expression tag	UNP P76461
D	402	HIS	-	expression tag	UNP P76461

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	214	Total O 214 214	0	0
2	B	200	Total O 200 200	0	0
2	C	144	Total O 144 144	0	0
2	D	134	Total O 134 134	0	0

- Molecule 1: Acetyl-CoA acetyltransferase





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	75.01Å 85.15Å 269.11Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.10 50.00 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.3 (50.00-2.10) 99.6 (50.00-2.10)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.49 (at 2.08Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.169 , 0.218 0.184 , 0.225	Depositor DCC
R_{free} test set	5100 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	28.4	Xtriage
Anisotropy	0.733	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 33.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	12024	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.89% 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	1.16	5/2865 (0.2%)	1.08	4/3874 (0.1%)
1	B	1.14	2/2865 (0.1%)	1.10	7/3874 (0.2%)
1	C	1.08	0/2865	1.05	5/3874 (0.1%)
1	D	1.09	2/2865 (0.1%)	1.06	4/3874 (0.1%)
All	All	1.12	9/11460 (0.1%)	1.07	20/15496 (0.1%)

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	C	0	1

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	284	VAL	C-N	6.48	1.41	1.33
1	A	201	VAL	N-CA	-6.01	1.40	1.46
1	A	12	THR	C-O	5.56	1.31	1.23
1	A	117	ASN	N-CA	-5.40	1.39	1.46
1	D	285	PRO	CA-C	5.35	1.54	1.51
1	D	165	VAL	C-O	-5.21	1.18	1.24
1	A	350	PRO	N-CD	5.15	1.54	1.47
1	B	285	PRO	N-CD	5.14	1.54	1.47
1	A	105	GLY	C-O	-5.03	1.17	1.24

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	351	ILE	N-CA-C	9.48	120.32	110.36
1	A	347	LEU	N-CA-C	7.01	119.96	111.82
1	B	294	VAL	CA-C-N	-6.72	112.07	119.19

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	294	VAL	C-N-CA	-6.72	112.07	119.19
1	B	284	VAL	CA-C-N	-6.64	113.55	120.38
1	B	284	VAL	C-N-CA	-6.64	113.55	120.38
1	B	210	LYS	N-CA-C	6.28	116.73	107.88
1	B	235	ALA	N-CA-C	6.26	118.18	111.36
1	A	294	VAL	CA-C-N	-5.50	113.36	119.19
1	A	294	VAL	C-N-CA	-5.50	113.36	119.19
1	C	210	LYS	N-CA-C	5.41	116.49	108.60
1	C	209	LYS	N-CA-C	5.23	116.78	111.14
1	C	140	VAL	CB-CA-C	5.21	117.45	110.42
1	B	378	LEU	N-CA-C	5.10	116.60	108.79
1	C	125	LEU	CA-C-N	5.06	129.42	120.87
1	C	125	LEU	C-N-CA	5.06	129.42	120.87
1	D	219	PHE	CA-C-N	-5.03	114.60	119.78
1	D	219	PHE	C-N-CA	-5.03	114.60	119.78
1	A	338	VAL	N-CA-C	5.03	115.55	108.36
1	D	262	GLU	N-CA-C	5.01	117.12	111.11

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	350	PRO	Peptide

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	2833	0	2908	24	0
1	B	2833	0	2908	25	0
1	C	2833	0	2908	29	0
1	D	2833	0	2908	25	0
2	A	214	0	0	1	0
2	B	200	0	0	2	0
2	C	144	0	0	1	0
2	D	134	0	0	1	0
All	All	12024	0	11632	95	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 4.

All (95) 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:209:LYS:O	1:A:210:LYS:HG2	1.45	1.12
1:D:369:ARG:HG3	1:D:369:ARG:HH11	1.43	0.83
1:A:126:ASP:HB2	2:A:548:HOH:O	1.78	0.83
1:C:100:GLN:HE21	1:D:100:GLN:HE21	1.24	0.82
1:B:210:LYS:HG2	1:B:211:THR:H	1.45	0.81
1:B:172:THR:OG1	1:B:175:MET:HG3	1.81	0.80
1:A:209:LYS:C	1:A:210:LYS:HG2	2.09	0.76
1:D:349:HIS:HD2	1:D:354:SER:OG	1.71	0.74
1:C:209:LYS:C	1:C:210:LYS:HG2	2.16	0.71
1:C:349:HIS:HD2	1:C:354:SER:OG	1.74	0.69
1:C:126:ASP:HB2	2:C:527:HOH:O	1.93	0.68
1:A:349:HIS:HD2	1:A:354:SER:OG	1.78	0.67
1:B:126:ASP:HB2	2:B:512:HOH:O	1.94	0.66
1:C:209:LYS:O	1:C:210:LYS:HD3	1.96	0.66
1:D:320:PHE:H	1:D:323:GLN:HE21	1.43	0.65
1:A:251:ASN:HD22	1:A:349:HIS:H	1.45	0.64
1:A:206:VAL:HG23	1:A:211:THR:OG1	1.97	0.63
1:D:70:LEU:HD23	1:D:70:LEU:C	2.27	0.59
1:A:320:PHE:H	1:A:323:GLN:HE21	1.50	0.59
1:C:172:THR:OG1	1:C:175:MET:HG3	2.03	0.59
1:D:369:ARG:HG3	1:D:369:ARG:NH1	2.14	0.59
1:D:126:ASP:HB2	2:D:512:HOH:O	2.01	0.59
1:D:148:GLY:O	1:D:149:LEU:HD23	2.03	0.59
1:C:1:MET:HE1	1:D:104:ALA:HB1	1.85	0.58
1:A:251:ASN:ND2	1:A:349:HIS:H	2.02	0.58
1:B:210:LYS:HG2	1:B:211:THR:N	2.18	0.57
1:B:234:PRO:HB2	1:B:237:ASP:O	2.07	0.55
1:C:70:LEU:C	1:C:70:LEU:HD12	2.31	0.55
1:A:124:LEU:HD23	1:A:125:LEU:O	2.08	0.54
1:C:272:PRO:HB2	1:C:393:LEU:HD12	1.89	0.54
1:D:4:CYS:SG	1:D:259:ILE:CG2	2.97	0.53
1:C:251:ASN:ND2	1:C:349:HIS:H	2.07	0.53
1:A:219:PHE:N	1:A:220:PRO:CD	2.71	0.53
1:B:219:PHE:N	1:B:220:PRO:CD	2.73	0.52
1:D:369:ARG:NH1	1:D:369:ARG:CG	2.73	0.52
1:D:251:ASN:ND2	1:D:349:HIS:H	2.07	0.51
1:B:55:ASN:HD22	1:B:58:GLN:HG2	1.76	0.51

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:158:MET:HA	1:D:158:MET:HE2	1.93	0.50
1:B:349:HIS:HD2	1:B:354:SER:OG	1.95	0.50
1:C:251:ASN:HD22	1:C:349:HIS:H	1.59	0.50
1:D:64:ASN:HD22	1:D:64:ASN:C	2.20	0.50
1:D:272:PRO:HB2	1:D:393:LEU:HD12	1.94	0.50
1:D:124:LEU:HD23	1:D:124:LEU:C	2.38	0.49
1:B:70:LEU:HD12	1:B:70:LEU:C	2.37	0.49
1:D:251:ASN:HD22	1:D:349:HIS:H	1.61	0.49
1:C:373:LEU:HD23	1:C:391:GLU:HG3	1.93	0.49
1:D:7:VAL:HG12	1:D:7:VAL:O	2.12	0.49
1:D:126:ASP:HB3	1:D:128:LYS:H	1.78	0.49
1:D:70:LEU:C	1:D:70:LEU:CD2	2.86	0.48
1:D:315:GLU:HG3	1:D:362:LEU:HB2	1.95	0.47
1:A:52:ILE:O	1:A:112:ALA:HA	2.15	0.47
1:C:209:LYS:C	1:C:210:LYS:CG	2.85	0.47
1:B:251:ASN:ND2	1:B:349:HIS:H	2.13	0.47
1:B:393:LEU:HA	2:B:676:HOH:O	2.15	0.47
1:C:171:ILE:CD1	1:C:326:ALA:HB2	2.45	0.46
1:C:219:PHE:N	1:C:220:PRO:CD	2.78	0.46
1:C:320:PHE:H	1:C:323:GLN:HE21	1.62	0.46
1:A:1:MET:HE1	1:B:104:ALA:HB1	1.98	0.46
1:A:62:GLY:HA2	1:B:146:ARG:O	2.16	0.46
1:A:81:GLY:H	1:B:384:GLN:HE22	1.63	0.46
1:B:4:CYS:SG	1:B:259:ILE:CG2	3.05	0.45
1:B:17:PHE:O	1:B:18:ASN:HB2	2.17	0.44
1:B:357:ARG:C	1:B:357:ARG:HD2	2.41	0.44
1:C:41:ALA:O	1:C:42:LYS:HB2	2.18	0.44
1:C:95:VAL:HG12	1:C:388:MET:HE1	1.99	0.44
1:A:294:VAL:HB	1:A:295:PRO:CD	2.47	0.44
1:B:334:ASP:OD1	1:B:336:GLU:HG2	2.18	0.44
1:C:158:MET:HA	1:C:158:MET:HE2	2.00	0.44
1:A:371:LYS:HD3	1:A:371:LYS:HA	1.76	0.43
1:C:171:ILE:HD11	1:C:326:ALA:HA	2.00	0.43
1:C:98:ALA:CB	1:C:259:ILE:HD11	2.48	0.43
1:C:55:ASN:HD22	1:C:58:GLN:HG2	1.84	0.43
1:B:142:ASP:OD2	1:C:135:LEU:HD23	2.19	0.43
1:C:318:GLU:HB3	1:C:345:ILE:HG13	1.99	0.43
1:B:292:GLY:N	1:B:293:PRO:CD	2.82	0.43
1:C:179:LEU:HD13	1:C:179:LEU:C	2.43	0.43
1:A:159:GLY:O	1:A:162:ALA:HB3	2.19	0.42
1:A:209:LYS:C	1:A:210:LYS:CG	2.89	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:124:LEU:HD23	1:B:125:LEU:C	2.43	0.42
1:A:148:GLY:O	1:A:149:LEU:HD12	2.19	0.42
1:C:123:TYR:HA	1:C:141:TYR:O	2.20	0.42
1:D:209:LYS:HD2	1:D:209:LYS:HA	1.91	0.42
1:A:149:LEU:O	1:A:158:MET:HG2	2.19	0.42
1:B:237:ASP:HB3	1:B:240:GLY:HA3	2.00	0.41
1:D:24:THR:O	1:D:117:ASN:ND2	2.52	0.41
1:A:167:LYS:O	1:A:168:GLU:C	2.62	0.41
1:A:237:ASP:O	1:A:240:GLY:N	2.49	0.41
1:B:318:GLU:CD	1:B:343:GLY:HA3	2.46	0.41
1:C:294:VAL:HB	1:C:295:PRO:CD	2.51	0.41
1:B:148:GLY:O	1:B:149:LEU:HD23	2.20	0.41
1:C:373:LEU:CD2	1:C:391:GLU:HG3	2.51	0.41
1:A:55:ASN:HD21	1:A:61:LEU:HD12	1.87	0.40
1:C:64:ASN:HB2	1:D:87:VAL:CG2	2.50	0.40
1:D:159:GLY:O	1:D:162:ALA:HB3	2.22	0.40
1:A:64:ASN:HB2	1:B:87:VAL:CG2	2.52	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	394/405 (97%)	382 (97%)	11 (3%)	1 (0%)	36	36
1	B	394/405 (97%)	384 (98%)	9 (2%)	1 (0%)	36	36
1	C	394/405 (97%)	380 (96%)	13 (3%)	1 (0%)	36	36
1	D	394/405 (97%)	386 (98%)	7 (2%)	1 (0%)	36	36
All	All	1576/1620 (97%)	1532 (97%)	40 (2%)	4 (0%)	36	36

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	87	VAL
1	B	87	VAL
1	C	87	VAL
1	D	87	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	283/292 (97%)	270 (95%)	13 (5%)	24	25
1	B	283/292 (97%)	275 (97%)	8 (3%)	38	43
1	C	283/292 (97%)	269 (95%)	14 (5%)	22	22
1	D	283/292 (97%)	267 (94%)	16 (6%)	18	17
All	All	1132/1168 (97%)	1081 (96%)	51 (4%)	24	25

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	70	LEU
1	A	87	VAL
1	A	126	ASP
1	A	128	LYS
1	A	134	ARG
1	A	156	TYR
1	A	192	GLU
1	A	206	VAL
1	A	238	LYS
1	A	259	ILE
1	A	299	LYS
1	A	378	LEU
1	A	393	LEU
1	B	70	LEU
1	B	87	VAL
1	B	88	CYS
1	B	126	ASP
1	B	134	ARG

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	140	VAL
1	B	207	THR
1	B	307	GLN
1	C	70	LEU
1	C	87	VAL
1	C	88	CYS
1	C	124	LEU
1	C	126	ASP
1	C	131	SER
1	C	135	LEU
1	C	156	TYR
1	C	174	GLU
1	C	207	THR
1	C	208	ARG
1	C	259	ILE
1	C	378	LEU
1	C	393	LEU
1	D	27	ILE
1	D	70	LEU
1	D	72	LYS
1	D	87	VAL
1	D	88	CYS
1	D	126	ASP
1	D	134	ARG
1	D	208	ARG
1	D	209	LYS
1	D	221	LYS
1	D	238	LYS
1	D	263	SER
1	D	286	PRO
1	D	302	GLN
1	D	336	GLU
1	D	378	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (30) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	55	ASN
1	A	68	GLN
1	A	216	GLN
1	A	251	ASN
1	A	323	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	349	HIS
1	B	55	ASN
1	B	68	GLN
1	B	251	ASN
1	B	349	HIS
1	B	384	GLN
1	C	46	GLN
1	C	47	HIS
1	C	55	ASN
1	C	63	GLN
1	C	154	HIS
1	C	251	ASN
1	C	323	GLN
1	C	349	HIS
1	C	384	GLN
1	D	46	GLN
1	D	55	ASN
1	D	63	GLN
1	D	64	ASN
1	D	100	GLN
1	D	223	ASN
1	D	251	ASN
1	D	302	GLN
1	D	323	GLN
1	D	349	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	396/405 (97%)	-0.19	9 (2%) 61 64	21, 31, 51, 115	0
1	B	396/405 (97%)	-0.29	2 (0%) 87 88	21, 31, 51, 89	0
1	C	396/405 (97%)	-0.02	5 (1%) 75 77	22, 36, 58, 106	0
1	D	396/405 (97%)	-0.13	5 (1%) 75 77	24, 34, 56, 106	0
All	All	1584/1620 (97%)	-0.16	21 (1%) 75 77	21, 33, 55, 115	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	210	LYS	6.8
1	A	207	THR	6.4
1	A	211	THR	4.0
1	D	148	GLY	2.9
1	B	393	LEU	2.8
1	A	209	LYS	2.7
1	A	148	GLY	2.6
1	A	212	PHE	2.6
1	C	126	ASP	2.6
1	C	-2	MET	2.5
1	D	-2	MET	2.4
1	A	192	GLU	2.3
1	C	210	LYS	2.3
1	D	206	VAL	2.3
1	D	236	PHE	2.2
1	C	333	PHE	2.1
1	A	-2	MET	2.1
1	A	206	VAL	2.1
1	C	-1	GLY	2.1
1	D	207	THR	2.1
1	B	148	GLY	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.