



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 29, 2023 – 02:08 PM JST

PDB ID : 7YVY
Title : Crystal structure of thiolase PFC_04095 from *Pyrococcus furiosus*
Authors : Singh, R.; Das, A.K.
Deposited on : 2022-08-20
Resolution : 2.70 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 : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.35
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) : 2.35

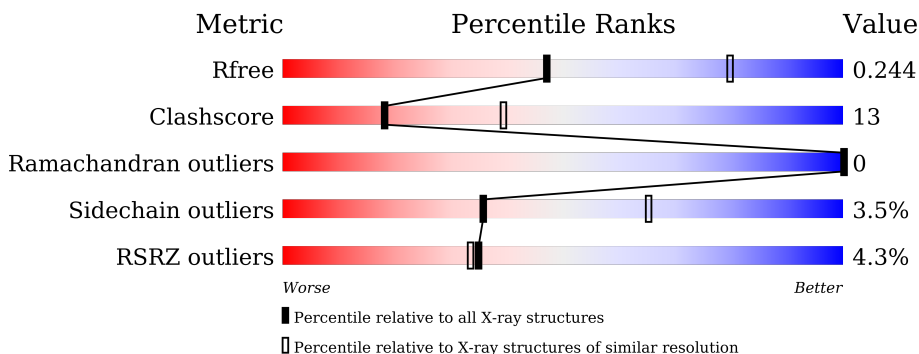
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	2808 (2.70-2.70)
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (2.70-2.70)

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	388	 6% 71% 22% • 5%
1	B	388	 2% 64% 29% • 6%
1	C	388	 4% 64% 28% • 7%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 7760 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.

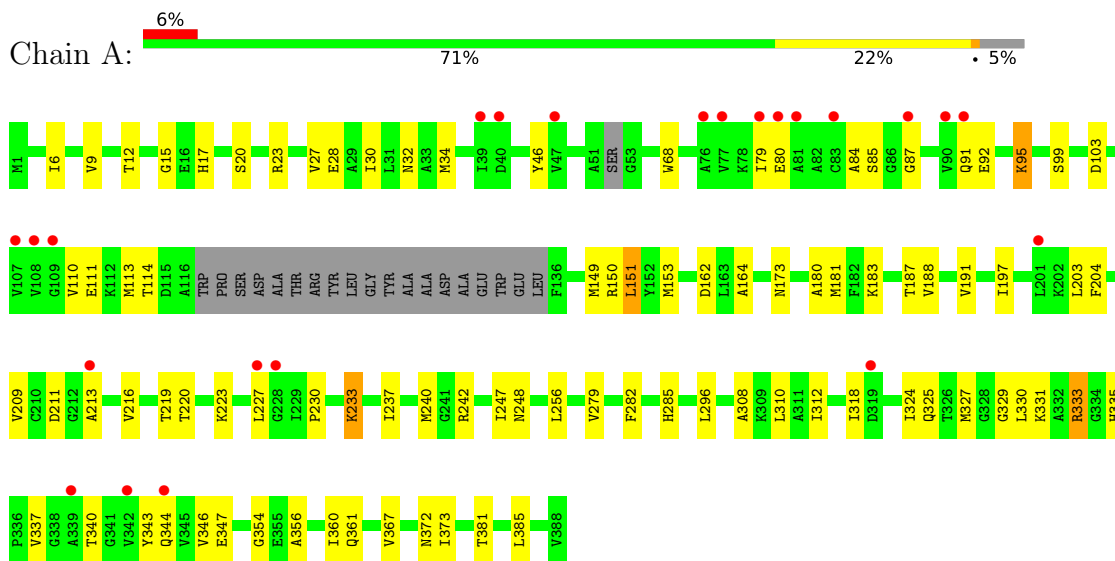
- Molecule 1 is a protein called Acetyl-CoA acetyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	368	Total 2637	C 1674	N 440	O 509	S 14	0	0	0
1	B	364	Total 2592	C 1643	N 435	O 501	S 13	0	0	0
1	C	359	Total 2531	C 1606	N 427	O 484	S 14	0	0	0

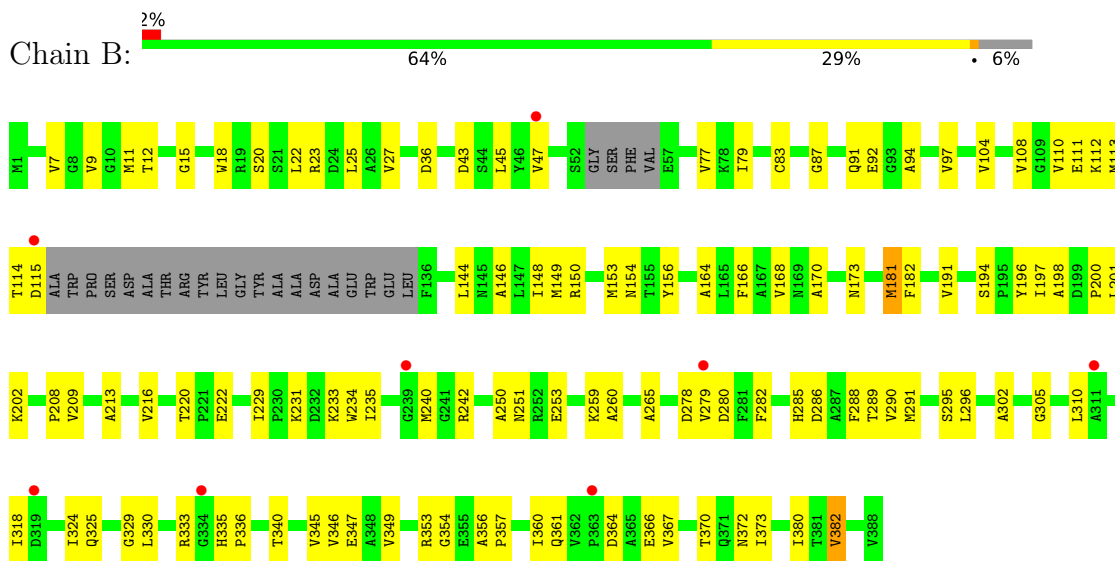
3 Residue-property plots i

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: Acetyl-CoA acetyltransferase

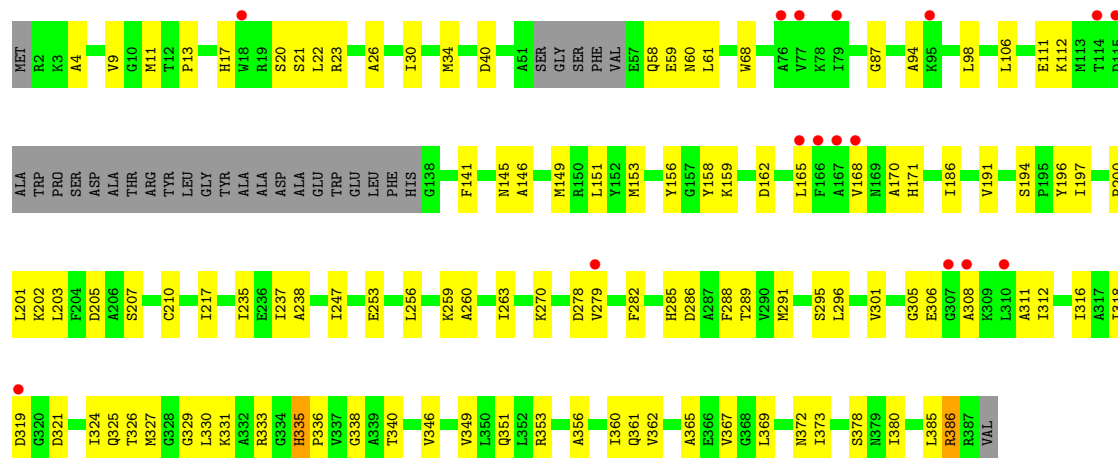


- Molecule 1: Acetyl-CoA acetyltransferase



- Molecule 1: Acetyl-CoA acetyltransferase

Chain C: 4% 64% 28% 7%



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	133.25Å 113.48Å 107.53Å 90.00° 123.03° 90.00°	Depositor
Resolution (Å)	48.03 – 2.70 48.03 – 2.35	Depositor EDS
% Data completeness (in resolution range)	99.3 (48.03-2.70) 91.0 (48.03-2.35)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.35 (at 2.34Å)	Xtrriage
Refinement program	PHENIX 1.19.2_4158	Depositor
R, R_{free}	0.209 , 0.252 0.208 , 0.244	Depositor DCC
R_{free} test set	2487 reflections (4.89%)	wwPDB-VP
Wilson B-factor (Å ²)	77.7	Xtrriage
Anisotropy	0.418	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 79.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7760	wwPDB-VP
Average B, all atoms (Å ²)	91.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.05% 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	0.34	0/2675	0.55	0/3640
1	B	0.35	0/2630	0.57	0/3585
1	C	0.36	0/2569	0.60	0/3506
All	All	0.35	0/7874	0.57	0/10731

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

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	336	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	2637	0	2614	58	0
1	B	2592	0	2538	76	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	2531	0	2463	72	0
All	All	7760	0	7615	206	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 13.

All (206) 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:113:MET:HE1	1:B:209:VAL:HG12	1.44	0.97
1:A:91:GLN:HA	1:A:240:MET:HE1	1.46	0.97
1:C:9:VAL:HG21	1:C:349:VAL:HG21	1.55	0.88
1:A:15:GLY:HA2	1:A:181:MET:HG3	1.55	0.87
1:B:356:ALA:HB3	1:B:361:GLN:HG2	1.61	0.82
1:A:164:ALA:HB1	1:A:191:VAL:HG21	1.62	0.81
1:C:356:ALA:HB3	1:C:361:GLN:HG2	1.66	0.77
1:C:279:VAL:HA	1:C:367:VAL:HG13	1.68	0.75
1:C:327:MET:HG2	1:C:360:ILE:HD12	1.67	0.75
1:B:91:GLN:HE22	1:B:242:ARG:HB2	1.50	0.74
1:B:113:MET:CE	1:B:209:VAL:HG12	2.17	0.74
1:C:168:VAL:HG12	1:C:186:ILE:HG13	1.69	0.73
1:B:335:HIS:ND1	1:B:340:THR:HG21	2.04	0.72
1:A:237:ILE:HG21	1:A:240:MET:HE3	1.72	0.71
1:B:279:VAL:HG13	1:B:367:VAL:HG12	1.73	0.71
1:B:291:MET:O	1:B:295:SER:OG	2.11	0.69
1:B:357:PRO:HG2	1:B:360:ILE:HD11	1.73	0.69
1:B:345:VAL:O	1:B:349:VAL:HG12	1.93	0.68
1:A:151:LEU:HB3	1:A:256:LEU:HD21	1.73	0.68
1:B:47:VAL:HG23	1:B:108:VAL:HG13	1.76	0.67
1:C:259:LYS:HG2	1:C:263:ILE:HD11	1.74	0.67
1:A:87:GLY:H	1:A:372:ASN:HD22	1.44	0.66
1:B:265:ALA:HA	1:B:382:VAL:HG21	1.78	0.66
1:A:308:ALA:O	1:A:312:ILE:HD12	1.96	0.66
1:A:12:THR:HG22	1:A:28:GLU:HG2	1.78	0.65
1:B:15:GLY:HA2	1:B:181:MET:HG3	1.77	0.65
1:B:197:ILE:HD11	1:B:201:LEU:HD22	1.79	0.65
1:A:23:ARG:O	1:A:27:VAL:HG22	1.98	0.64
1:A:327:MET:HA	1:A:360:ILE:HD12	1.81	0.63
1:C:326:THR:HB	1:C:351:GLN:HE21	1.63	0.63
1:C:259:LYS:O	1:C:263:ILE:HD12	1.98	0.63
1:C:171:HIS:HE1	1:C:331:LYS:HE3	1.63	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:171:HIS:CE1	1:C:331:LYS:HE3	2.35	0.62
1:C:87:GLY:H	1:C:372:ASN:HD22	1.46	0.62
1:B:112:LYS:HE2	1:B:115:ASP:OD1	2.00	0.61
1:C:279:VAL:HG12	1:C:367:VAL:HG22	1.82	0.61
1:C:111:GLU:HG3	1:C:338:GLY:H	1.65	0.61
1:C:335:HIS:ND1	1:C:340:THR:HG21	2.16	0.61
1:A:242:ARG:HG3	1:A:381:THR:HG22	1.83	0.61
1:B:45:LEU:HD11	1:B:108:VAL:HG12	1.81	0.61
1:B:149:MET:HG3	1:B:290:VAL:CG2	2.31	0.60
1:B:150:ARG:HA	1:B:153:MET:HG2	1.85	0.59
1:C:197:ILE:HG13	1:C:201:LEU:HD23	1.84	0.59
1:C:201:LEU:HD11	1:C:288:PHE:CD1	2.37	0.59
1:C:9:VAL:HG13	1:C:346:VAL:HG13	1.83	0.59
1:A:333:ARG:HD3	1:A:347:GLU:OE2	2.02	0.59
1:C:165:LEU:HA	1:C:168:VAL:HG22	1.85	0.59
1:C:362:VAL:HG23	1:C:365:ALA:HB2	1.85	0.59
1:C:253:GLU:OE2	1:C:259:LYS:NZ	2.21	0.59
1:B:149:MET:HG3	1:B:290:VAL:HG22	1.83	0.58
1:A:149:MET:O	1:A:153:MET:HG3	2.03	0.58
1:C:260:ALA:HB3	1:C:380:ILE:HD12	1.86	0.58
1:B:198:ALA:O	1:B:201:LEU:N	2.37	0.58
1:A:103:ASP:O	1:A:219:THR:HG22	2.04	0.57
1:B:302:ALA:HB2	1:B:310:LEU:HD11	1.86	0.57
1:C:30:ILE:HD13	1:C:106:LEU:HD21	1.86	0.57
1:A:279:VAL:HA	1:A:367:VAL:HB	1.87	0.57
1:B:280:ASP:OD2	1:B:366:GLU:N	2.31	0.56
1:B:20:SER:O	1:B:112:LYS:NZ	2.34	0.56
1:B:253:GLU:OE2	1:B:259:LYS:NZ	2.34	0.56
1:C:191:VAL:O	1:C:194:SER:OG	2.22	0.56
1:B:354:GLY:HA2	1:B:361:GLN:NE2	2.21	0.56
1:A:285:HIS:CD2	1:A:373:ILE:H	2.24	0.56
1:C:318:ILE:HD11	1:C:360:ILE:HG22	1.86	0.56
1:A:247:ILE:HD12	1:A:248:ASN:N	2.21	0.56
1:B:12:THR:HG22	1:B:213:ALA:H	1.71	0.56
1:B:285:HIS:HB3	1:B:373:ILE:HD12	1.88	0.56
1:C:9:VAL:HG21	1:C:349:VAL:CG2	2.32	0.56
1:B:250:ALA:O	1:B:251:ASN:ND2	2.39	0.55
1:C:235:ILE:HD11	1:C:353:ARG:HG2	1.89	0.55
1:C:196:TYR:CE1	1:C:202:LYS:HG2	2.42	0.55
1:B:79:ILE:HD12	1:B:92:GLU:HG3	1.87	0.55
1:B:191:VAL:O	1:B:194:SER:OG	2.20	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:36:ASP:OD2	1:B:353:ARG:NH2	2.40	0.54
1:C:34:MET:HG2	1:C:217:ILE:HD11	1.88	0.54
1:C:291:MET:O	1:C:295:SER:OG	2.23	0.54
1:C:351:GLN:OE1	1:C:361:GLN:HA	2.08	0.54
1:C:23:ARG:HG3	1:C:68:TRP:CD2	2.44	0.53
1:B:146:ALA:HB1	1:B:200:PRO:HG2	1.91	0.53
1:B:279:VAL:HB	1:B:282:PHE:CZ	2.43	0.53
1:C:158:TYR:CE1	1:C:308:ALA:HB3	2.42	0.53
1:B:354:GLY:HA2	1:B:361:GLN:HE21	1.73	0.53
1:C:326:THR:HB	1:C:351:GLN:NE2	2.23	0.53
1:A:113:MET:HE2	1:A:209:VAL:HB	1.91	0.53
1:C:151:LEU:HG	1:C:256:LEU:HD11	1.90	0.52
1:B:329:GLY:O	1:B:333:ARG:HB3	2.09	0.52
1:C:282:PHE:HD2	1:C:369:LEU:HD22	1.75	0.52
1:A:17:HIS:HB3	1:A:20:SER:HB2	1.90	0.52
1:A:164:ALA:CB	1:A:191:VAL:HG21	2.38	0.52
1:C:286:ASP:OD1	1:C:330:LEU:HB2	2.09	0.52
1:A:162:ASP:HB3	1:A:312:ILE:HD11	1.92	0.51
1:C:59:GLU:O	1:C:60:ASN:HB2	2.11	0.51
1:A:197:ILE:HD11	1:A:203:LEU:HD12	1.92	0.51
1:C:170:ALA:HB2	1:C:325:GLN:NE2	2.25	0.51
1:B:91:GLN:HE21	1:B:242:ARG:HH11	1.58	0.51
1:A:230:PRO:HG2	1:A:233:LYS:HG3	1.93	0.51
1:C:58:GLN:O	1:C:61:LEU:HB2	2.11	0.50
1:B:23:ARG:O	1:B:27:VAL:HG23	2.12	0.50
1:B:260:ALA:HB3	1:B:380:ILE:HD12	1.93	0.50
1:C:349:VAL:O	1:C:353:ARG:HG3	2.11	0.50
1:B:166:PHE:HB2	1:B:289:THR:HG22	1.95	0.49
1:A:80:GLU:HA	1:A:85:SER:HB3	1.94	0.49
1:B:144:LEU:O	1:B:148:ILE:HG12	2.12	0.49
1:B:9:VAL:HB	1:B:346:VAL:HG13	1.93	0.49
1:B:111:GLU:HG3	1:B:113:MET:HG3	1.95	0.49
1:B:9:VAL:HG11	1:B:349:VAL:HG11	1.95	0.49
1:C:373:ILE:HG23	1:C:380:ILE:HG12	1.94	0.49
1:B:318:ILE:HD12	1:B:318:ILE:H	1.77	0.49
1:B:370:THR:O	1:B:382:VAL:HA	2.13	0.49
1:C:17:HIS:HB3	1:C:20:SER:OG	2.13	0.48
1:C:11:MET:SD	1:C:346:VAL:HG11	2.53	0.48
1:C:156:TYR:CD2	1:C:305:GLY:HA2	2.48	0.48
1:C:308:ALA:O	1:C:312:ILE:HG13	2.13	0.48
1:C:159:LYS:N	1:C:162:ASP:OD1	2.43	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:87:GLY:H	1:A:372:ASN:ND2	2.08	0.48
1:B:156:TYR:CD1	1:B:305:GLY:HA2	2.49	0.48
1:B:182:PHE:HE2	1:B:208:PRO:HG3	1.79	0.48
1:C:22:LEU:HD12	1:C:112:LYS:HB2	1.96	0.48
1:A:191:VAL:HG12	1:A:204:PHE:HB3	1.95	0.48
1:B:233:LYS:HA	1:B:353:ARG:HD3	1.97	0.47
1:B:110:VAL:HG23	1:B:213:ALA:HB2	1.96	0.47
1:C:111:GLU:CG	1:C:338:GLY:H	2.26	0.47
1:C:279:VAL:HG23	1:C:282:PHE:CZ	2.49	0.47
1:A:282:PHE:HB2	1:A:324:ILE:HG22	1.97	0.47
1:C:9:VAL:CG1	1:C:346:VAL:HG13	2.44	0.47
1:C:286:ASP:OD2	1:C:325:GLN:HG2	2.14	0.47
1:B:43:ASP:HB2	1:B:104:VAL:O	2.15	0.47
1:A:91:GLN:HA	1:A:240:MET:CE	2.31	0.47
1:B:336:PRO:HD2	1:B:340:THR:HG22	1.98	0.46
1:B:231:LYS:HA	1:B:234:TRP:CG	2.51	0.46
1:C:296:LEU:HD21	1:C:324:ILE:HG21	1.97	0.46
1:C:205:ASP:OD2	1:C:289:THR:OG1	2.26	0.46
1:A:32:ASN:N	1:A:32:ASN:HD22	2.14	0.46
1:C:149:MET:O	1:C:153:MET:HG3	2.16	0.45
1:B:170:ALA:HB2	1:B:325:GLN:NE2	2.30	0.45
1:B:173:ASN:OD1	1:B:318:ILE:HD11	2.16	0.45
1:B:9:VAL:HG11	1:B:349:VAL:CG1	2.46	0.45
1:C:316:ILE:HB	1:C:324:ILE:HG12	1.97	0.45
1:A:110:VAL:HG23	1:A:213:ALA:HB2	1.98	0.45
1:A:23:ARG:HG3	1:A:68:TRP:CD2	2.52	0.45
1:B:91:GLN:HG2	1:B:240:MET:HG2	1.99	0.45
1:A:92:GLU:O	1:A:95:LYS:HG2	2.17	0.45
1:B:286:ASP:OD2	1:B:330:LEU:N	2.47	0.45
1:C:329:GLY:O	1:C:333:ARG:HB3	2.17	0.45
1:C:165:LEU:HD23	1:C:311:ALA:O	2.17	0.45
1:C:26:ALA:O	1:C:30:ILE:HG13	2.18	0.44
1:C:146:ALA:HB1	1:C:200:PRO:HD2	1.98	0.44
1:C:238:ALA:HB2	1:C:386:ARG:HD2	1.99	0.44
1:A:324:ILE:HD12	1:A:325:GLN:HG3	1.99	0.44
1:C:141:PHE:O	1:C:145:ASN:ND2	2.49	0.44
1:A:223:LYS:HE3	1:A:227:LEU:HD21	2.00	0.44
1:A:318:ILE:HD12	1:A:318:ILE:H	1.83	0.44
1:B:333:ARG:NE	1:B:347:GLU:OE2	2.50	0.44
1:B:201:LEU:HD11	1:B:288:PHE:CD1	2.53	0.44
1:A:329:GLY:O	1:A:333:ARG:HB3	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:335:HIS:CE1	1:A:337:VAL:HG12	2.53	0.44
1:C:23:ARG:HG3	1:C:68:TRP:CG	2.52	0.43
1:C:270:LYS:HD3	1:C:270:LYS:HA	1.60	0.43
1:A:84:ALA:HA	1:A:372:ASN:ND2	2.33	0.43
1:B:22:LEU:HA	1:B:25:LEU:HD12	1.99	0.43
1:B:234:TRP:O	1:B:235:ILE:HD13	2.19	0.43
1:C:4:ALA:HB2	1:C:98:LEU:CD1	2.48	0.43
1:C:285:HIS:O	1:C:330:LEU:HD13	2.17	0.43
1:A:150:ARG:HD3	1:A:150:ARG:HA	1.81	0.43
1:A:340:THR:O	1:A:344:GLN:HG3	2.19	0.43
1:A:343:TYR:HA	1:A:346:VAL:HG22	2.00	0.43
1:A:330:LEU:HD12	1:A:333:ARG:NH2	2.33	0.43
1:B:164:ALA:O	1:B:168:VAL:HG13	2.19	0.43
1:C:207:SER:HB3	1:C:330:LEU:O	2.18	0.43
1:C:247:ILE:CD1	1:C:378:SER:HB3	2.49	0.43
1:B:149:MET:O	1:B:153:MET:HG2	2.19	0.43
1:A:113:MET:CE	1:A:209:VAL:HB	2.48	0.43
1:B:196:TYR:CE1	1:B:202:LYS:HG2	2.54	0.43
1:B:235:ILE:CG1	1:B:349:VAL:HG23	2.49	0.42
1:A:318:ILE:HD12	1:A:318:ILE:N	2.34	0.42
1:C:13:PRO:O	1:C:17:HIS:NE2	2.42	0.42
1:A:111:GLU:O	1:A:211:ASP:HA	2.20	0.42
1:C:40:ASP:OD1	1:C:40:ASP:N	2.53	0.42
1:B:296:LEU:HD21	1:B:324:ILE:HG21	2.00	0.42
1:B:182:PHE:CE2	1:B:208:PRO:HG3	2.54	0.42
1:B:113:MET:HB2	1:B:113:MET:HE2	1.52	0.42
1:A:114:THR:O	1:A:114:THR:OG1	2.38	0.42
1:A:296:LEU:HD23	1:A:310:LEU:HD12	2.01	0.42
1:A:356:ALA:HB3	1:A:361:GLN:OE1	2.19	0.42
1:A:9:VAL:HG12	1:A:216:VAL:HG22	2.01	0.41
1:A:237:ILE:HD13	1:A:385:LEU:HD21	2.02	0.41
1:B:18:TRP:CZ3	1:B:114:THR:HB	2.55	0.41
1:C:165:LEU:HA	1:C:168:VAL:CG2	2.49	0.41
1:A:173:ASN:CG	1:A:318:ILE:HD11	2.40	0.41
1:B:94:ALA:O	1:B:97:VAL:HG22	2.21	0.41
1:A:92:GLU:OE1	1:A:242:ARG:NH1	2.44	0.41
1:C:158:TYR:HH	1:C:306:GLU:HA	1.85	0.41
1:A:6:ILE:HD11	1:A:237:ILE:HD11	2.02	0.41
1:A:30:ILE:O	1:A:34:MET:HG3	2.20	0.41
1:B:7:VAL:HB	1:B:229:ILE:HG21	2.03	0.41
1:A:180:ALA:O	1:A:183:LYS:HE2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:47:VAL:HA	1:B:108:VAL:O	2.21	0.41
1:B:87:GLY:H	1:B:372:ASN:HD22	1.66	0.41
1:A:188:VAL:O	1:A:191:VAL:HG22	2.21	0.40
1:B:150:ARG:O	1:B:154:ASN:HB2	2.22	0.40
1:B:220:THR:OG1	1:B:222:GLU:OE2	2.38	0.40
1:B:222:GLU:CD	1:B:222:GLU:H	2.25	0.40
1:A:46:TYR:HD2	1:A:79:ILE:HD11	1.86	0.40
1:C:94:ALA:HB2	1:C:237:ILE:HD13	2.02	0.40
1:A:354:GLY:HA2	1:A:361:GLN:HE22	1.86	0.40
1:B:197:ILE:HD11	1:B:201:LEU:CD2	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	362/388 (93%)	350 (97%)	12 (3%)	0	100	100
1	B	358/388 (92%)	348 (97%)	10 (3%)	0	100	100
1	C	353/388 (91%)	340 (96%)	13 (4%)	0	100	100
All	All	1073/1164 (92%)	1038 (97%)	35 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	257/290 (89%)	249 (97%)	8 (3%)	40	69
1	B	249/290 (86%)	241 (97%)	8 (3%)	39	68
1	C	239/290 (82%)	229 (96%)	10 (4%)	30	58
All	All	745/870 (86%)	719 (96%)	26 (4%)	36	65

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	95	LYS
1	A	99	SER
1	A	151	LEU
1	A	187	THR
1	A	220	THR
1	A	233	LYS
1	A	331	LYS
1	A	333	ARG
1	B	11	MET
1	B	77	VAL
1	B	83	CYS
1	B	181	MET
1	B	216	VAL
1	B	278	ASP
1	B	364	ASP
1	B	382	VAL
1	C	21	SER
1	C	203	LEU
1	C	210	CYS
1	C	278	ASP
1	C	301	VAL
1	C	319	ASP
1	C	321	ASP
1	C	335	HIS
1	C	385	LEU
1	C	386	ARG

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

Mol	Chain	Res	Type
1	A	32	ASN
1	A	169	ASN
1	A	285	HIS

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Mol	Chain	Res	Type
1	A	325	GLN
1	A	372	ASN
1	B	91	GLN
1	B	251	ASN
1	C	58	GLN
1	C	372	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 monosaccharides 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	368/388 (94%)	0.40	23 (6%) 20 19	57, 86, 119, 160	0
1	B	364/388 (93%)	0.16	8 (2%) 62 63	65, 87, 111, 145	0
1	C	359/388 (92%)	0.31	16 (4%) 33 31	72, 93, 118, 136	0
All	All	1091/1164 (93%)	0.29	47 (4%) 35 33	57, 89, 117, 160	0

All (47) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	40	ASP	4.4
1	C	167	ALA	3.6
1	A	319	ASP	3.2
1	A	227	LEU	3.2
1	B	363	PRO	3.1
1	B	239	GLY	3.1
1	C	310	LEU	3.0
1	C	114	THR	2.9
1	A	107	VAL	2.9
1	B	115	ASP	2.9
1	A	47	VAL	2.8
1	C	76	ALA	2.8
1	A	80	GLU	2.8
1	C	166	PHE	2.8
1	B	334	GLY	2.8
1	A	228	GLY	2.7
1	C	319	ASP	2.5
1	B	319	ASP	2.5
1	C	307	GLY	2.4
1	B	279	VAL	2.4
1	A	91	GLN	2.4
1	A	39	ILE	2.4
1	A	81	ALA	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	79	ILE	2.4
1	A	342	VAL	2.3
1	A	76	ALA	2.3
1	A	344	GLN	2.3
1	A	87	GLY	2.2
1	C	77	VAL	2.2
1	A	77	VAL	2.2
1	C	115	ASP	2.2
1	C	165	LEU	2.2
1	A	213	ALA	2.2
1	A	90	VAL	2.1
1	A	339	ALA	2.1
1	B	47	VAL	2.1
1	C	308	ALA	2.1
1	C	168	VAL	2.1
1	C	279	VAL	2.1
1	B	311	ALA	2.1
1	A	83	CYS	2.1
1	C	95	LYS	2.1
1	A	201	LEU	2.0
1	C	18	TRP	2.0
1	A	109	GLY	2.0
1	A	108	VAL	2.0
1	C	79	ILE	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 monosaccharides 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.