



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 15, 2026 – 01:01 AM UTC

PDB ID : 3DMY / pdb_00003dmy
Title : Crystal Structure of a predicated acyl-CoA synthetase from E.coli
Authors : Sugadev, R.; Burley, S.K.; Swaminathan, S.; New York SGX Research Center
for Structural Genomics (NYSGXRC)
Deposited on : 2008-07-01
Resolution : 2.07 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
Mogul	:	2022.3.0, CSD as543be (2022)
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

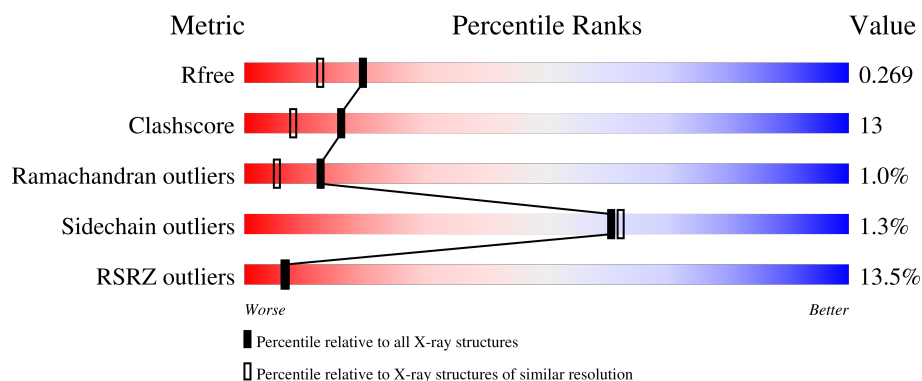
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.07 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	3774 (2.08-2.04)
Clashscore	190562	3883 (2.08-2.04)
Ramachandran outliers	187476	3860 (2.08-2.04)
Sidechain outliers	187428	3860 (2.08-2.04)
RSRZ outliers	180081	3775 (2.08-2.04)

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	480	15% 69% 22% 7%
1	B	480	10% 72% 19% 7%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6952 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 Protein fdxA.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	447	Total	C	N	O	S	Se	0	0	0
			3299	2073	581	628	7	10			
1	B	446	Total	C	N	O	S	Se	0	0	0
			3291	2069	580	625	7	10			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	556	GLY	-	expression tag	UNP Q47208
A	557	HIS	-	expression tag	UNP Q47208
A	558	PRO	-	expression tag	UNP Q47208
A	559	HIS	-	expression tag	UNP Q47208
A	560	HIS	-	expression tag	UNP Q47208
A	561	HIS	-	expression tag	UNP Q47208
B	556	GLY	-	expression tag	UNP Q47208
B	557	HIS	-	expression tag	UNP Q47208
B	558	PRO	-	expression tag	UNP Q47208
B	559	HIS	-	expression tag	UNP Q47208
B	560	HIS	-	expression tag	UNP Q47208
B	561	HIS	-	expression tag	UNP Q47208

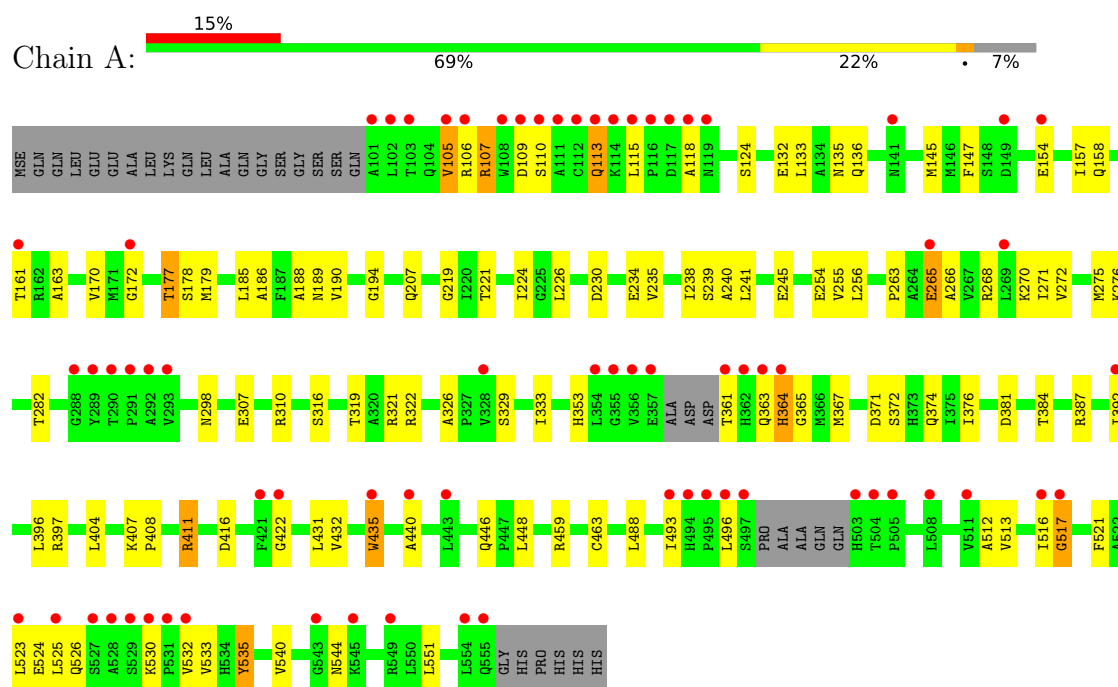
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	167	Total	O	0	0
			167	167		
2	B	195	Total	O	0	0
			195	195		

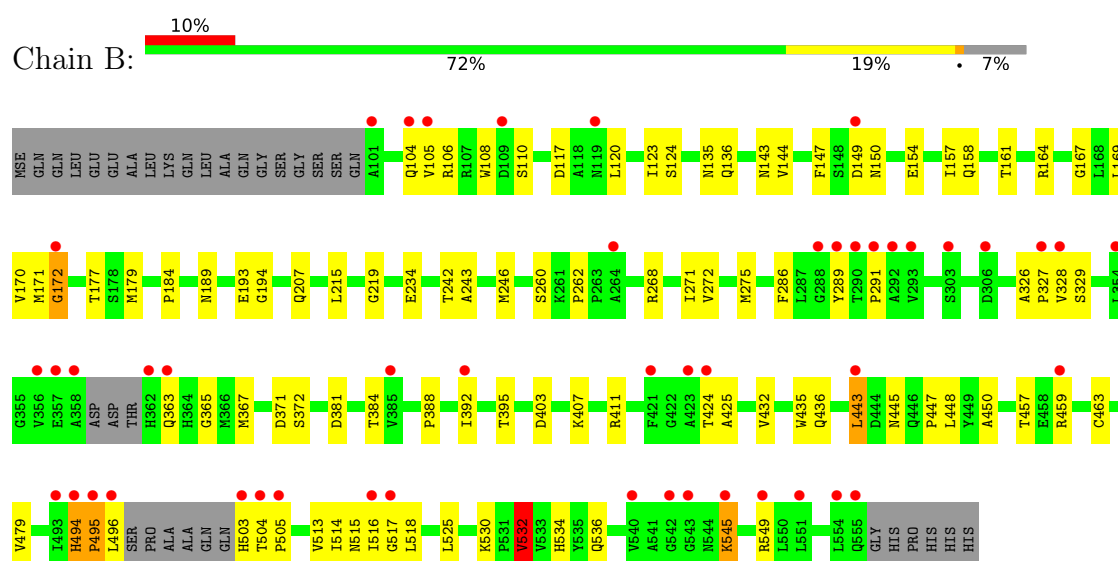
3 Residue-property plots [i](#)

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: Protein fdrA



• Molecule 1: Protein fdrA



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	63.57Å 107.78Å 153.60Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.97 – 2.07 48.97 – 2.07	Depositor EDS
% Data completeness (in resolution range)	86.7 (48.97-2.07) 86.8 (48.97-2.07)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.66 (at 1.98Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.237 , 0.270 0.237 , 0.269	Depositor DCC
R_{free} test set	2826 reflections (4.24%)	wwPDB-VP
Wilson B-factor (Å ²)	23.6	Xtriage
Anisotropy	0.518	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 37.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	6952	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.44% 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	0.41	0/3339	0.95	12/4524 (0.3%)
1	B	0.44	0/3331	0.99	13/4513 (0.3%)
All	All	0.42	0/6670	0.97	25/9037 (0.3%)

There are no bond length outliers.

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	513	VAL	N-CA-C	8.69	121.66	108.71
1	A	513	VAL	N-CA-C	8.60	120.21	108.17
1	B	516	ILE	N-CA-C	7.94	118.39	111.56
1	B	147	PHE	N-CA-C	-7.52	102.44	111.69
1	B	407	LYS	CA-C-N	7.35	126.88	119.24
1	B	407	LYS	C-N-CA	7.35	126.88	119.24
1	A	376	ILE	N-CA-C	7.18	118.22	108.17
1	A	307	GLU	N-CA-C	-6.90	103.69	111.14
1	B	177	THR	N-CA-C	5.88	119.05	109.06
1	A	221	THR	N-CA-C	-5.80	104.97	112.68
1	A	364	HIS	N-CA-C	5.65	119.36	111.90
1	A	224	ILE	N-CA-C	5.64	115.91	107.51
1	A	115	LEU	CA-C-N	5.47	124.93	119.24
1	A	115	LEU	C-N-CA	5.47	124.93	119.24
1	B	514	ILE	N-CA-C	-5.41	99.44	107.51
1	B	443	LEU	N-CA-C	-5.40	103.25	110.55
1	A	147	PHE	N-CA-C	-5.39	106.20	112.89
1	B	532	VAL	CB-CA-C	-5.39	103.81	111.40
1	A	177	THR	N-CA-C	5.32	118.11	109.06
1	B	123	ILE	N-CA-C	5.23	115.49	108.17
1	A	404	LEU	N-CA-C	-5.17	106.08	112.38
1	B	515	ASN	CA-C-N	-5.05	117.98	123.08
1	B	515	ASN	C-N-CA	-5.05	117.98	123.08
1	A	363	GLN	N-CA-C	5.02	117.58	110.50
1	B	457	THR	N-CA-C	5.00	117.49	110.14

There are no chirality outliers.

There are no planarity outliers.

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	3299	0	3356	90	0
1	B	3291	0	3349	86	0
2	A	167	0	0	14	0
2	B	195	0	0	15	0
All	All	6952	0	6705	175	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 (175) 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:169:LEU:HD22	1:B:246:MSE:HE1	1.41	1.02
1:B:243:ALA:HA	1:B:246:MSE:HE3	1.42	1.00
1:B:242:THR:HG22	1:B:246:MSE:HE2	1.46	0.95
1:B:108:TRP:HE1	1:B:136:GLN:HE21	1.18	0.92
1:A:263:PRO:HG2	1:A:268:ARG:HB3	1.51	0.91
1:A:525:LEU:HB3	1:A:532:VAL:HG21	1.56	0.88
1:B:108:TRP:HE1	1:B:136:GLN:NE2	1.75	0.83
1:A:135:ASN:HD21	1:A:158:GLN:HE22	1.27	0.82
1:B:327:PRO:HB2	1:B:495:PRO:HB3	1.64	0.80
1:B:495:PRO:HG2	1:B:496:LEU:H	1.49	0.78
1:B:365:GLY:C	1:B:367:MSE:HE3	2.09	0.78
1:B:157:ILE:O	1:B:161:THR:HG23	1.85	0.76
1:A:365:GLY:C	1:A:367:MSE:HE2	2.11	0.76
1:B:327:PRO:HB2	1:B:495:PRO:CB	2.16	0.75
1:A:524:GLU:HB3	2:A:597:HOH:O	1.86	0.74
1:A:241:LEU:O	1:A:245:GLU:HG3	1.92	0.69
1:A:446:GLN:HG3	2:A:725:HOH:O	1.93	0.69
1:B:388:PRO:HB3	1:B:392:ILE:HD11	1.76	0.68
1:B:135:ASN:HD21	1:B:158:GLN:HE22	1.39	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:105:VAL:HG13	1:B:110:SER:HB2	1.78	0.65
1:A:272:VAL:HG12	1:A:276:LYS:HE3	1.79	0.64
1:A:365:GLY:HA2	1:A:367:MSE:CE	2.28	0.63
1:A:157:ILE:O	1:A:161:THR:HG23	2.00	0.62
1:A:408:PRO:HB2	1:A:446:GLN:HE21	1.64	0.62
1:B:271:ILE:CG2	1:B:275:MSE:HE3	2.29	0.62
1:A:525:LEU:CB	1:A:532:VAL:HG21	2.27	0.61
1:B:432:VAL:O	1:B:436:GLN:HG3	2.00	0.61
1:A:105:VAL:HG13	1:A:106:ARG:N	2.15	0.61
1:B:479:VAL:HA	2:B:753:HOH:O	2.00	0.61
1:A:321:ARG:HD2	1:A:488:LEU:HD21	1.82	0.60
1:A:266:ALA:HB3	2:A:698:HOH:O	2.01	0.60
1:A:154:GLU:H	1:A:154:GLU:CD	2.10	0.60
1:A:530:LYS:O	1:A:532:VAL:HG23	2.02	0.60
1:B:530:LYS:O	1:B:532:VAL:HG23	2.01	0.60
1:B:189:ASN:HD21	1:B:207:GLN:HE21	1.49	0.60
1:A:113:GLN:NE2	2:A:673:HOH:O	2.35	0.59
1:A:255:VAL:HG21	1:A:316:SER:HA	1.83	0.59
1:A:118:ALA:HA	1:A:512:ALA:HB3	1.85	0.58
1:B:105:VAL:CG1	1:B:110:SER:HB2	2.33	0.58
1:B:106:ARG:HG3	1:B:106:ARG:HH11	1.69	0.58
1:B:108:TRP:NE1	1:B:136:GLN:HE21	1.94	0.58
1:B:271:ILE:HG22	1:B:275:MSE:HE3	1.86	0.57
1:A:268:ARG:HD2	2:A:618:HOH:O	2.05	0.57
1:B:545:LYS:O	1:B:549:ARG:HG3	2.05	0.57
1:B:525:LEU:CB	1:B:532:VAL:HG21	2.35	0.56
1:A:256:LEU:HB2	1:A:282:THR:HG23	1.87	0.56
1:B:494:HIS:HB2	1:B:495:PRO:HA	1.86	0.56
1:A:353:HIS:HD2	2:A:637:HOH:O	1.88	0.56
1:B:365:GLY:CA	1:B:367:MSE:HE3	2.34	0.56
1:A:365:GLY:C	1:A:367:MSE:CE	2.77	0.56
1:A:435:TRP:HZ3	1:A:448:LEU:O	1.88	0.56
1:B:503:HIS:HB2	2:B:755:HOH:O	2.06	0.56
1:B:179:MSE:HE3	1:B:184:PRO:HB3	1.86	0.55
1:A:145:MSE:SE	1:A:178:SER:HB2	2.57	0.55
1:A:179:MSE:HG2	1:A:188:ALA:HB1	1.89	0.54
1:B:363:GLN:NE2	2:B:720:HOH:O	2.40	0.54
1:A:326:ALA:HA	1:A:496:LEU:HD11	1.90	0.54
1:A:365:GLY:HA2	1:A:367:MSE:HE3	1.90	0.54
1:A:392:ILE:HD11	2:A:599:HOH:O	2.08	0.53
1:A:135:ASN:HD21	1:A:158:GLN:NE2	2.00	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:178:SER:C	1:A:179:MSE:HG3	2.34	0.53
1:A:265:GLU:HB2	2:A:593:HOH:O	2.09	0.53
1:A:107:ARG:HB3	1:A:107:ARG:HH11	1.73	0.53
1:A:540:VAL:HG12	1:A:544:ASN:HB3	1.91	0.52
1:A:133:LEU:HD22	1:A:516:ILE:HD13	1.91	0.52
1:B:193:GLU:HB3	2:B:591:HOH:O	2.09	0.52
1:A:329:SER:O	1:A:411:ARG:HD2	2.09	0.52
1:A:310:ARG:NH1	2:A:681:HOH:O	2.41	0.52
1:A:105:VAL:CG1	1:A:106:ARG:N	2.73	0.52
1:A:268:ARG:HH11	1:A:268:ARG:HG3	1.75	0.52
1:A:321:ARG:HD2	1:A:488:LEU:CD2	2.39	0.52
1:A:387:ARG:HD2	1:A:396:LEU:HD22	1.92	0.51
1:A:392:ILE:HG22	1:A:422:GLY:O	2.10	0.51
1:B:365:GLY:HA2	1:B:367:MSE:HE3	1.92	0.51
1:A:189:ASN:HD21	1:A:207:GLN:HE21	1.60	0.51
1:A:241:LEU:HD11	1:A:270:LYS:HD3	1.93	0.51
1:B:435:TRP:HH2	1:B:448:LEU:O	1.94	0.50
1:A:365:GLY:CA	1:A:367:MSE:CE	2.90	0.50
1:B:169:LEU:CD2	1:B:246:MSE:HE1	2.28	0.50
1:A:177:THR:CG2	1:A:179:MSE:HE2	2.42	0.50
1:B:234:GLU:N	1:B:234:GLU:OE1	2.45	0.49
1:B:525:LEU:HB3	1:B:532:VAL:HG21	1.94	0.49
1:A:521:PHE:O	1:A:524:GLU:HG2	2.12	0.49
1:A:268:ARG:O	1:A:272:VAL:HG23	2.13	0.49
1:B:154:GLU:H	1:B:154:GLU:CD	2.20	0.49
1:B:525:LEU:HB2	1:B:532:VAL:HG21	1.94	0.49
1:B:135:ASN:HD21	1:B:158:GLN:NE2	2.09	0.49
1:A:132:GLU:O	1:A:136:GLN:HG3	2.12	0.49
1:B:179:MSE:HG2	2:B:648:HOH:O	2.13	0.48
1:B:189:ASN:ND2	1:B:207:GLN:HE21	2.11	0.48
1:B:495:PRO:CG	1:B:496:LEU:H	2.17	0.48
1:A:118:ALA:HA	1:A:512:ALA:CB	2.43	0.48
1:B:443:LEU:HB3	1:B:445:ASN:OD1	2.14	0.48
1:A:105:VAL:HG13	1:A:107:ARG:H	1.79	0.47
1:A:124:SER:OG	1:A:517:GLY:HA3	2.15	0.47
1:B:403:ASP:HB3	2:B:592:HOH:O	2.13	0.47
1:B:144:VAL:HB	1:B:170:VAL:HG22	1.96	0.47
1:B:371:ASP:O	1:B:372:SER:HB2	2.15	0.47
1:B:167:GLY:HA2	2:B:707:HOH:O	2.15	0.47
1:A:397:ARG:NH2	1:A:416:ASP:OD1	2.48	0.46
1:B:367:MSE:HE2	1:B:367:MSE:HA	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:408:PRO:HB2	1:A:446:GLN:NE2	2.29	0.46
1:A:254:GLU:HB3	1:A:319:THR:HG21	1.98	0.46
1:A:440:ALA:HA	2:A:678:HOH:O	2.15	0.46
1:A:432:VAL:O	1:A:435:TRP:HD1	1.98	0.45
1:A:271:ILE:O	1:A:275:MSE:HG3	2.17	0.45
1:A:374:GLN:NE2	2:A:608:HOH:O	2.48	0.45
1:A:459:ARG:O	1:A:459:ARG:HG3	2.17	0.45
1:B:435:TRP:CH2	1:B:448:LEU:O	2.69	0.45
1:A:523:LEU:O	1:A:526:GLN:HB3	2.16	0.45
1:B:262:PRO:HG3	1:B:289:TYR:HB2	1.99	0.45
1:B:117:ASP:OD2	1:B:117:ASP:N	2.50	0.45
1:B:164:ARG:NH1	2:B:703:HOH:O	2.46	0.45
1:B:268:ARG:O	1:B:272:VAL:HG23	2.17	0.44
1:A:256:LEU:O	1:A:282:THR:HA	2.18	0.44
1:B:503:HIS:CD2	2:B:694:HOH:O	2.71	0.44
1:A:364:HIS:ND1	2:A:623:HOH:O	2.36	0.44
1:B:411:ARG:HD3	1:B:447:PRO:O	2.16	0.44
1:B:494:HIS:CB	1:B:495:PRO:HA	2.48	0.44
1:A:230:ASP:OD1	1:A:239:SER:HB2	2.18	0.44
1:B:243:ALA:CA	1:B:246:MSE:HE3	2.29	0.44
1:B:534:HIS:CD2	1:B:536:GLN:HG2	2.52	0.44
1:A:234:GLU:OE1	1:A:234:GLU:N	2.50	0.44
1:A:333:ILE:HG13	1:A:493:ILE:HG21	2.00	0.44
1:A:535:TYR:CD2	1:A:535:TYR:C	2.96	0.44
1:B:381:ASP:HA	1:B:384:THR:HG1	1.82	0.44
1:B:243:ALA:HA	1:B:246:MSE:CE	2.31	0.43
1:B:327:PRO:HB2	1:B:495:PRO:HB2	1.95	0.43
1:A:107:ARG:HH11	1:A:107:ARG:CB	2.31	0.43
1:A:326:ALA:CA	1:A:496:LEU:HD11	2.47	0.43
1:B:105:VAL:CG1	1:B:106:ARG:N	2.82	0.43
1:B:242:THR:CG2	1:B:246:MSE:HE2	2.33	0.43
1:A:268:ARG:HG3	1:A:268:ARG:NH1	2.34	0.43
1:A:322:ARG:NH1	1:A:496:LEU:HD23	2.34	0.43
1:B:104:GLN:CD	1:B:536:GLN:NE2	2.77	0.43
1:B:120:LEU:HD12	1:B:143:ASN:C	2.44	0.43
1:A:107:ARG:NH1	1:A:109:ASP:OD1	2.52	0.43
1:A:407:LYS:HE3	2:A:638:HOH:O	2.18	0.42
1:B:194:GLY:HA3	1:B:219:GLY:C	2.43	0.42
1:B:149:ASP:CG	1:B:150:ASN:H	2.27	0.42
1:A:163:ALA:HB2	1:A:170:VAL:HG23	2.01	0.42
1:A:371:ASP:O	1:A:372:SER:HB2	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:381:ASP:HA	1:A:384:THR:OG1	2.19	0.42
1:A:107:ARG:HB2	1:A:110:SER:OG	2.20	0.42
1:B:171:MSE:O	1:B:172:GLY:O	2.38	0.42
1:B:328:VAL:C	1:B:495:PRO:HG3	2.45	0.42
1:B:326:ALA:HA	2:B:640:HOH:O	2.20	0.42
1:B:381:ASP:CG	2:B:660:HOH:O	2.61	0.42
1:A:185:LEU:O	1:A:186:ALA:HB3	2.19	0.42
1:B:495:PRO:CG	1:B:496:LEU:N	2.80	0.42
1:A:361:THR:HG22	1:A:361:THR:O	2.19	0.42
1:B:189:ASN:HD21	1:B:207:GLN:NE2	2.17	0.42
1:B:260:SER:O	1:B:286:PHE:HA	2.20	0.42
1:A:113:GLN:HE21	1:A:113:GLN:HB2	1.59	0.41
1:A:235:VAL:O	1:A:238:ILE:HG13	2.20	0.41
1:B:124:SER:OG	1:B:518:LEU:N	2.42	0.41
1:B:164:ARG:HD3	2:B:741:HOH:O	2.20	0.41
1:B:494:HIS:CE1	2:B:679:HOH:O	2.73	0.41
1:A:431:LEU:HD23	1:A:431:LEU:HA	1.91	0.41
1:B:395:THR:HB	2:B:655:HOH:O	2.20	0.41
1:A:459:ARG:NH1	2:A:656:HOH:O	2.53	0.41
1:B:262:PRO:CG	1:B:289:TYR:HB2	2.51	0.41
1:B:424:THR:HG22	1:B:425:ALA:N	2.35	0.41
1:B:504:THR:HA	1:B:505:PRO:HD3	1.94	0.41
1:B:215:LEU:HD12	1:B:215:LEU:HA	1.94	0.41
1:A:226:LEU:HD13	1:A:240:ALA:HB2	2.02	0.41
1:A:107:ARG:HB3	1:A:107:ARG:NH1	2.36	0.41
1:A:190:VAL:HB	1:B:459:ARG:HB2	2.03	0.41
1:B:289:TYR:O	1:B:291:PRO:HD3	2.21	0.41
1:B:329:SER:HA	2:B:731:HOH:O	2.21	0.41
1:A:533:VAL:O	1:A:533:VAL:HG13	2.21	0.40
1:B:435:TRP:CZ3	1:B:450:ALA:HB2	2.57	0.40
1:A:276:LYS:HA	1:A:298:ASN:OD1	2.21	0.40
1:B:286:PHE:O	1:B:289:TYR:HB3	2.20	0.40
1:A:194:GLY:HA3	1:A:219:GLY:C	2.47	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	441/480 (92%)	419 (95%)	18 (4%)	4 (1%)	14	7
1	B	440/480 (92%)	418 (95%)	17 (4%)	5 (1%)	11	4
All	All	881/960 (92%)	837 (95%)	35 (4%)	9 (1%)	12	5

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	172	GLY
1	B	172	GLY
1	B	495	PRO
1	A	265	GLU
1	A	517	GLY
1	B	494	HIS
1	B	517	GLY
1	B	463	CYS
1	A	463	CYS

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	345/359 (96%)	338 (98%)	7 (2%)	48	47
1	B	343/359 (96%)	341 (99%)	2 (1%)	78	82
All	All	688/718 (96%)	679 (99%)	9 (1%)	61	62

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	105	VAL
1	A	107	ARG
1	A	113	GLN
1	A	411	ARG
1	A	435	TRP
1	A	535	TYR
1	A	551	LEU
1	B	532	VAL
1	B	545	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	113	GLN
1	A	158	GLN
1	A	189	ASN
1	A	353	HIS
1	A	363	GLN
1	A	374	GLN
1	A	398	ASN
1	A	446	GLN
1	A	466	GLN
1	A	515	ASN
1	A	534	HIS
1	A	536	GLN
1	A	544	ASN
1	A	555	GLN
1	B	119	ASN
1	B	136	GLN
1	B	158	GLN
1	B	189	ASN
1	B	212	GLN
1	B	363	GLN
1	B	374	GLN
1	B	398	ASN
1	B	536	GLN

5.3.3 RNA ⓘ

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 oligosaccharides 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	437/480 (91%)	1.02	70 (16%) 5 5	18, 36, 59, 75	0
1	B	436/480 (90%)	0.75	48 (11%) 10 10	16, 31, 54, 74	0
All	All	873/960 (90%)	0.88	118 (13%) 7 6	16, 33, 58, 75	0

All (118) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	172	GLY	7.5
1	B	496	LEU	6.6
1	A	361	THR	6.3
1	B	358	ALA	5.7
1	B	362	HIS	5.4
1	B	292	ALA	5.2
1	A	362	HIS	4.7
1	A	172	GLY	4.4
1	B	495	PRO	4.3
1	A	101	ALA	4.2
1	A	328	VAL	4.0
1	B	291	PRO	3.9
1	A	532	VAL	3.8
1	A	494	HIS	3.8
1	B	289	TYR	3.8
1	B	356	VAL	3.8
1	B	119	ASN	3.7
1	A	363	GLN	3.7
1	A	357	GLU	3.5
1	B	288	GLY	3.5
1	A	356	VAL	3.5
1	A	555	GLN	3.5
1	A	495	PRO	3.5
1	A	517	GLY	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	496	LEU	3.4
1	A	503	HIS	3.4
1	A	119	ASN	3.4
1	B	494	HIS	3.4
1	A	290	THR	3.4
1	A	497	SER	3.4
1	A	525	LEU	3.3
1	B	504	THR	3.3
1	A	289	TYR	3.2
1	A	293	VAL	3.1
1	B	290	THR	3.1
1	A	265	GLU	3.1
1	B	421	PHE	3.1
1	A	527	SER	3.1
1	A	549	ARG	3.0
1	B	363	GLN	2.9
1	B	493	ILE	2.9
1	A	292	ALA	2.9
1	B	293	VAL	2.9
1	A	493	ILE	2.8
1	A	116	PRO	2.8
1	A	516	ILE	2.8
1	B	503	HIS	2.8
1	B	328	VAL	2.8
1	A	113	GLN	2.8
1	B	549	ARG	2.7
1	B	517	GLY	2.7
1	B	443	LEU	2.6
1	B	357	GLU	2.6
1	A	110	SER	2.6
1	A	545	LYS	2.6
1	A	288	GLY	2.6
1	B	555	GLN	2.6
1	A	528	ALA	2.6
1	B	105	VAL	2.6
1	A	109	ASP	2.6
1	B	149	ASP	2.6
1	A	108	TRP	2.6
1	A	117	ASP	2.5
1	B	303	SER	2.5
1	A	102	LEU	2.5
1	A	435	TRP	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	392	ILE	2.5
1	A	105	VAL	2.5
1	B	545	LYS	2.4
1	A	504	THR	2.4
1	A	149	ASP	2.4
1	B	540	VAL	2.4
1	A	141	ASN	2.3
1	A	523	LEU	2.3
1	A	161	THR	2.3
1	B	306	ASP	2.3
1	A	529	SER	2.3
1	B	542	GLY	2.3
1	B	543	GLY	2.3
1	A	355	GLY	2.3
1	A	508	LEU	2.3
1	A	364	HIS	2.2
1	B	392	ILE	2.2
1	A	291	PRO	2.2
1	A	118	ALA	2.2
1	A	422	GLY	2.2
1	A	440	ALA	2.2
1	B	423	ALA	2.2
1	A	115	LEU	2.2
1	B	264	ALA	2.2
1	B	554	LEU	2.2
1	B	385	VAL	2.2
1	A	154	GLU	2.2
1	A	111	ALA	2.2
1	B	104	GLN	2.2
1	A	112	CYS	2.2
1	B	551	LEU	2.2
1	A	103	THR	2.2
1	A	421	PHE	2.1
1	B	101	ALA	2.1
1	B	505	PRO	2.1
1	B	516	ILE	2.1
1	A	511	VAL	2.1
1	B	459	ARG	2.1
1	B	354	LEU	2.1
1	B	109	ASP	2.1
1	A	531	PRO	2.1
1	A	106	ARG	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	354	LEU	2.1
1	B	424	THR	2.1
1	A	114	LYS	2.1
1	A	269	LEU	2.0
1	A	505	PRO	2.0
1	B	327	PRO	2.0
1	A	543	GLY	2.0
1	A	443	LEU	2.0
1	A	554	LEU	2.0
1	A	530	LYS	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.