Full wwPDB X-ray Structure Validation Report (i)

Jan 30, 2021 – 02:12 PM EST

PDB ID	:	3DMY
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 (i) symbol.

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

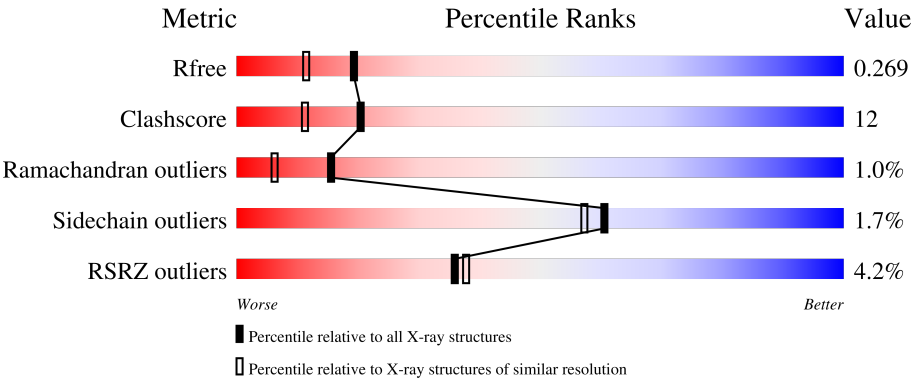
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.16
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.16

1 Overall quality at a glance (i)

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}	130704	2684 (2.08-2.04)
Clashscore	141614	2801 (2.08-2.04)
Ramachandran outliers	138981	2768 (2.08-2.04)
Sidechain outliers	138945	2768 (2.08-2.04)
RSRZ outliers	127900	2646 (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 <=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	4% 71% 20% · 7%
1	B	480	3% 76% 16% · 7%

2 Entry composition (i)

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 fdrA.

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.

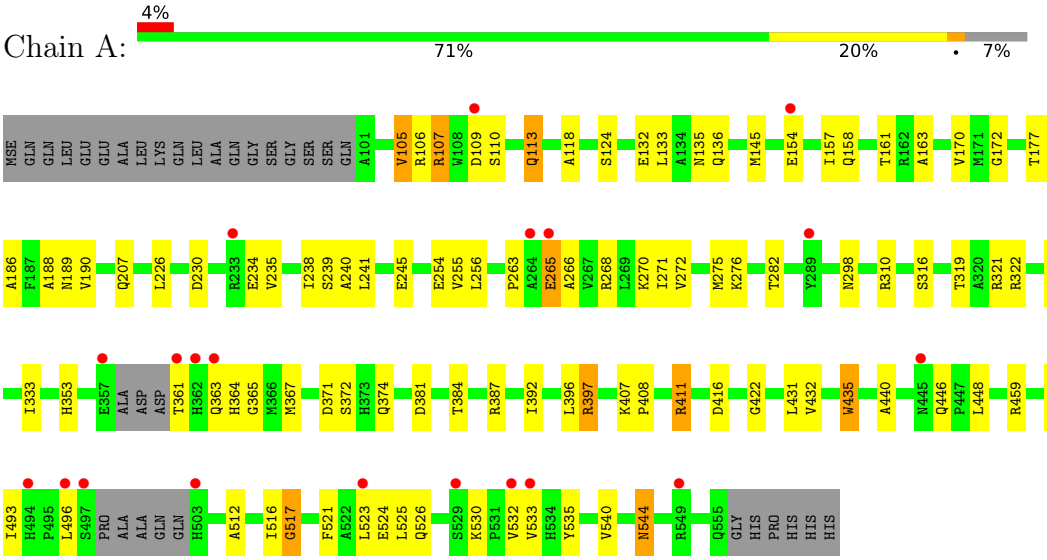
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	167	Total	O	0	0
			167	167		
2	B	195	Total	O	0	0
			195	195		

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.

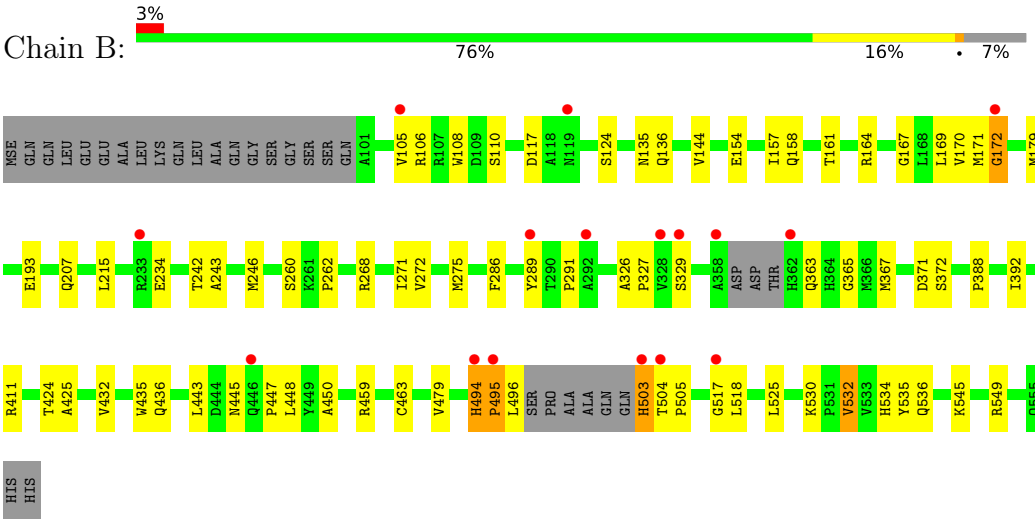
• Molecule 1: Protein fdrA

Chain A:



• Molecule 1: Protein fdrA

Chain B:



4 Data and refinement statistics (i)

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 – 1.99	Depositor EDS
% Data completeness (in resolution range)	86.7 (48.97-2.07) 81.6 (48.97-1.99)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [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 , 40.1	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	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.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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	0.32	0/3339	0.62	0/4524
1	B	0.35	0/3331	0.65	0/4513
All	All	0.34	0/6670	0.63	0/9037

There are no bond length outliers.

There are no bond angle outliers.

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	88	0
1	B	3291	0	3349	78	0
2	A	167	0	0	14	0
2	B	195	0	0	14	0
All	All	6952	0	6705	165	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 12.

All (165) 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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:157:ILE:O	1:B:161:THR:HG23	1.85	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:B:365:GLY:C	1:B:367:MSE:HE3	2.09	0.73
1:A:365:GLY:C	1:A:367:MSE:HE2	2.11	0.71
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
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: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:106:ARG:HG3	1:B:106:ARG:HH11	1.69	0.58
1:B:105:VAL:CG1	1:B:110:SER:HB2	2.33	0.58
1:B:271:ILE:HG22	1:B:275:MSE:HE3	1.86	0.57
1:B:365:GLY:CA	1:B:367:MSE:HE3	2.34	0.57
1:A:268:ARG:HD2	2:A:618:HOH:O	2.05	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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: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:A:154:GLU:CD	1:A:154:GLU:H	2.10	0.55
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
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:HG3	1:A:268:ARG:HH11	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:365:GLY:C	1:A:367:MSE:CE	2.77	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:OE1	1:B:234:GLU:N	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:525:LEU:HB2	1:B:532:VAL:HG21	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:178:SER:C	1:A:179:MSE:HG3	2.34	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:HA	1:B:367:MSE:HE2	1.97	0.46
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:HG3	1:A:459:ARG:O	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:N	1:B:117:ASP:OD2	2.50	0.45
1:B:495:PRO:CG	1:B:496:LEU:N	2.80	0.45
1:B:154:GLU:H	1:B:154:GLU:CD	2.20	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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: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: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
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:326:ALA:HA	2:B:640:HOH:O	2.20	0.42
1:A:185:LEU:O	1:A:186:ALA:HB3	2.19	0.42
1:A:361:THR:O	1:A:361:THR:HG22	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:HB2	1:A:113:GLN:HE21	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:HA	1:A:431:LEU:HD23	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:HA	1:B:215:LEU:HD12	1.94	0.41
1:A:226:LEU:HD13	1:A:240:ALA:HB2	2.02	0.41
1:A:190:VAL:HB	1:B:459:ARG:HB2	2.03	0.41
1:A:107:ARG:HB3	1:A:107:ARG:NH1	2.36	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

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%)	17 8
1	B	440/480 (92%)	418 (95%)	17 (4%)	5 (1%)	14 5
All	All	881/960 (92%)	837 (95%)	35 (4%)	9 (1%)	15 6

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%)	336 (97%)	9 (3%)	46 40
1	B	343/359 (96%)	340 (99%)	3 (1%)	78 78

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	688/718 (96%)	676 (98%)	12 (2%)	60 57

All (12) 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	363	GLN
1	A	397	ARG
1	A	411	ARG
1	A	435	TRP
1	A	535	TYR
1	A	544	ASN
1	B	503	HIS
1	B	532	VAL
1	B	535	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) 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	526	GLN
1	A	536	GLN
1	A	544	ASN
1	B	113	GLN
1	B	119	ASN
1	B	136	GLN
1	B	158	GLN
1	B	189	ASN
1	B	323	ASN
1	B	363	GLN
1	B	374	GLN

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Mol	Chain	Res	Type
1	B	398	ASN
1	B	536	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no 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, 95^{th} 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%)	0.45	21 (4%)	30	31	18, 36, 59, 75	0
1	B	436/480 (90%)	0.34	16 (3%)	41	43	16, 31, 54, 74	0
All	All	873/960 (90%)	0.40	37 (4%)	36	38	16, 33, 58, 75	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	292	ALA	6.2
1	A	361	THR	5.3
1	A	328	VAL	5.0
1	B	328	VAL	4.9
1	B	172	GLY	4.6
1	A	496	LEU	4.4
1	A	497	SER	4.0
1	B	495	PRO	4.0
1	B	503	HIS	3.8
1	A	494	HIS	3.3
1	A	363	GLN	3.2
1	B	494	HIS	3.2
1	A	503	HIS	3.2
1	A	233	ARG	3.0
1	B	358	ALA	2.9
1	B	504	THR	2.8
1	B	289	TYR	2.6
1	A	532	VAL	2.6
1	A	362	HIS	2.6
1	B	362	HIS	2.6
1	A	523	LEU	2.6
1	A	265	GLU	2.4
1	B	446	GLN	2.4
1	A	154	GLU	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	264	ALA	2.4
1	B	105	VAL	2.3
1	A	549	ARG	2.3
1	A	357	GLU	2.3
1	B	119	ASN	2.3
1	B	517	GLY	2.2
1	A	445	ASN	2.2
1	A	533	VAL	2.2
1	B	329	SER	2.1
1	A	109	ASP	2.1
1	B	233	ARG	2.1
1	A	289	TYR	2.1
1	A	529	SER	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.