



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2018 – 10:34 pm GMT

PDB ID : 4A10
Title : Apo-structure of 2-octenoyl-CoA carboxylase reductase CinF from streptomyces sp.
Authors : Quade, N.; Huo, L.; Rachid, S.; Heinz, D.W.; Muller, R.
Deposited on : 2011-09-13
Resolution : 2.25 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 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	:	trunk30967
Percentile statistics	:	20171227.v01 (using entries in the PDB archive December 27th 2017)
Refmac	:	5.8.0158
CCP4	:	7.0 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk30967

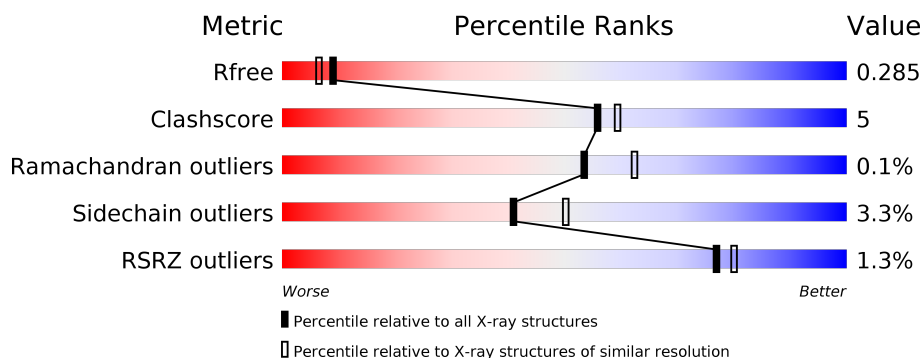
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.25 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}	111664	1178 (2.26-2.26)
Clashscore	122126	1286 (2.26-2.26)
Ramachandran outliers	120053	1253 (2.26-2.26)
Sidechain outliers	120020	1254 (2.26-2.26)
RSRZ outliers	108989	1158 (2.26-2.26)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	447	 % 81% 13% • 5%
1	B	447	 2% 83% 10% • •
1	C	447	 % 84% 9% • 7%
1	D	447	 % 83% 9% • 7%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 14058 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 OCTENOYL-COA REDUCTASE/CARBOXYLASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	424	Total	C	N	O	S	0	1	0
			3162	1988	572	586	16			
1	B	428	Total	C	N	O	S	0	0	0
			3197	2006	579	596	16			
1	C	417	Total	C	N	O	S	0	0	0
			3097	1943	562	576	16			
1	D	417	Total	C	N	O	S	0	0	0
			3087	1937	555	580	15			

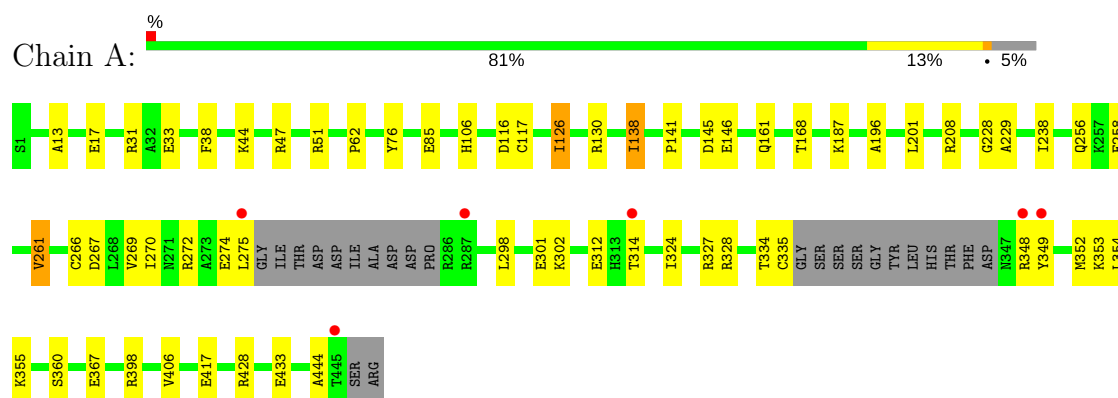
- Molecule 2 is water.

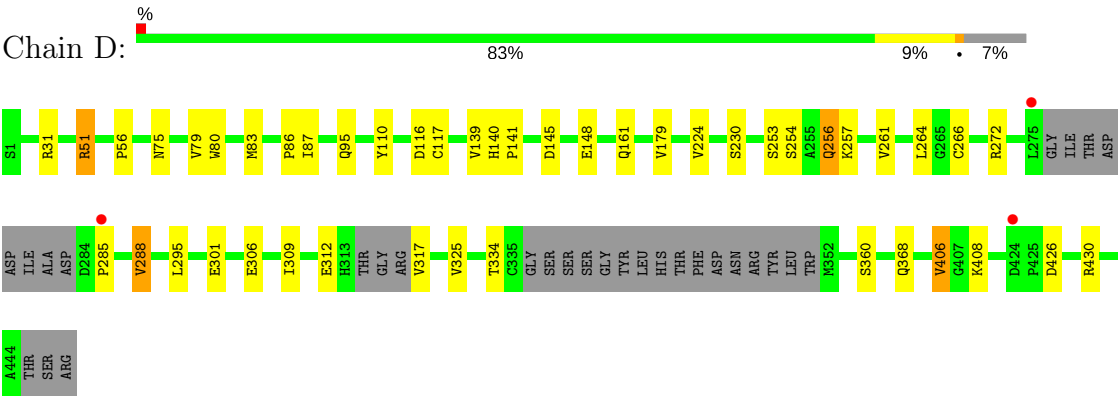
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	399	Total	O	0	0
			399	399		
2	B	404	Total	O	0	0
			404	404		
2	C	365	Total	O	0	0
			365	365		
2	D	347	Total	O	0	0
			347	347		

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA 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: OCTENOYL-COA REDUCTASE/CARBOXYLASE





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	101.64Å 85.22Å 113.97Å 90.00° 107.45° 90.00°	Depositor
Resolution (Å)	48.48 – 2.25 48.48 – 2.10	Depositor EDS
% Data completeness (in resolution range)	100.0 (48.48-2.25) 98.7 (48.48-2.10)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.36 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.209 , 0.269 0.226 , 0.285	Depositor DCC
R_{free} test set	5354 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	20.9	Xtriage
Anisotropy	0.288	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 27.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	14058	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.93% 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	0.51	0/3227	0.64	0/4387
1	B	0.53	0/3260	0.68	2/4428 (0.0%)
1	C	0.50	0/3155	0.62	0/4283
1	D	0.51	0/3146	0.63	0/4278
All	All	0.51	0/12788	0.64	2/17376 (0.0%)

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	B	0	1
1	C	0	1
All	All	0	2

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	336	GLY	N-CA-C	6.20	128.60	113.10
1	B	154	ASP	CB-CG-OD1	6.05	123.75	118.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	335	CYS	Peptide
1	C	274	GLU	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	3162	0	3135	44	0
1	B	3197	0	3170	37	0
1	C	3097	0	3084	30	0
1	D	3087	0	3039	34	0
2	A	399	0	0	8	0
2	B	404	0	0	6	0
2	C	365	0	0	8	0
2	D	347	0	0	8	0
All	All	14058	0	12428	137	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 5.

All (137) 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:D:264:LEU:HD12	2:D:737:HOH:O	1.65	0.95
1:B:138:ILE:HD11	1:B:196:ALA:HB1	1.56	0.86
1:A:138:ILE:HD11	1:A:196:ALA:HB1	1.62	0.81
1:C:352:MET:HB2	2:C:674:HOH:O	1.80	0.80
1:C:395:GLU:O	1:C:399:VAL:HG23	1.83	0.79
1:C:272:ARG:HD3	2:C:758:HOH:O	1.81	0.79
1:B:116:ASP:OD1	1:B:200:PRO:HB2	1.90	0.72
1:B:116:ASP:O	1:B:117:CYS:HB3	1.93	0.69
1:A:328:ARG:NH2	1:D:161:GLN:OE1	2.26	0.69
1:C:317:VAL:HG13	1:C:318:THR:HG23	1.74	0.69
1:A:335:CYS:CB	1:A:360:SER:O	2.41	0.68
1:B:295:LEU:HD23	1:B:325:VAL:HG11	1.74	0.68
1:A:444:ALA:O	2:A:501:HOH:O	2.12	0.68
1:B:347:ASN:ND2	2:B:503:HOH:O	2.27	0.67
1:B:46:VAL:HG22	1:B:82:ALA:HB3	1.76	0.67
1:C:130:ARG:NH2	1:D:145:ASP:OD2	2.28	0.66
1:A:130:ARG:NH2	1:B:145:ASP:OD2	2.26	0.66
1:A:145:ASP:OD2	1:B:130:ARG:NH2	2.26	0.65
1:C:139:VAL:HG11	1:C:179:VAL:HG11	1.78	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:13:ALA:O	1:A:17:GLU:HG2	1.95	0.65
1:B:38:PHE:HB3	1:B:44:LYS:HG2	1.79	0.64
1:D:230:SER:OG	1:D:257:LYS:HD3	1.99	0.63
1:A:335:CYS:HB3	1:A:360:SER:O	1.99	0.62
1:D:161:GLN:NE2	2:D:504:HOH:O	2.31	0.62
1:D:56:PRO:HG2	1:D:110:TYR:OH	1.99	0.62
1:D:139:VAL:HG11	1:D:179:VAL:HG11	1.81	0.62
1:D:301:GLU:HG3	2:D:535:HOH:O	1.99	0.62
1:D:80:TRP:CE3	1:D:83:MET:HE3	2.34	0.61
1:D:285:PRO:HA	1:D:288:VAL:HG13	1.82	0.61
1:D:31:ARG:HG3	1:D:51:ARG:HD2	1.81	0.61
1:D:261:VAL:CG1	1:D:266:CYS:HB3	2.31	0.61
1:A:335:CYS:HB2	1:A:360:SER:O	2.01	0.60
1:C:47:ARG:CZ	1:C:398:ARG:HD2	2.31	0.60
1:B:275:LEU:HG	1:B:291:THR:HG23	1.84	0.59
1:A:141:PRO:HG3	1:A:201:LEU:HD12	1.83	0.59
1:A:47:ARG:CZ	1:A:398:ARG:HD2	2.32	0.59
1:B:348:ARG:O	1:B:352:MET:HB2	2.02	0.59
1:A:126:ILE:HD11	1:B:126:ILE:HD11	1.85	0.58
1:A:348:ARG:HB3	1:A:352:MET:HE2	1.84	0.58
1:A:31:ARG:HH21	1:A:51:ARG:CZ	2.17	0.58
1:C:434:ASP:HB3	2:C:809:HOH:O	2.03	0.57
1:D:306:GLU:HG2	2:D:650:HOH:O	2.05	0.57
1:A:146:GLU:HB3	1:B:146:GLU:HB3	1.85	0.57
1:B:333:VAL:HG22	1:B:358:VAL:HB	1.87	0.57
1:D:295:LEU:HD23	1:D:325:VAL:HG11	1.89	0.55
1:B:61:ALA:HB1	1:B:62:PRO:HD2	1.88	0.55
1:B:46:VAL:HG22	1:B:82:ALA:CB	2.37	0.55
1:A:228:GLY:H	1:A:272:ARG:HH22	1.55	0.54
1:C:198:VAL:O	1:C:408:LYS:HE2	2.07	0.54
1:B:154:ASP:HB3	1:B:157:LEU:HD22	1.89	0.54
1:A:229:ALA:HB1	1:A:238:ILE:HD11	1.90	0.53
1:C:405:GLN:HE21	1:C:405:GLN:H	1.55	0.53
1:D:79:VAL:HG12	1:D:83:MET:HE2	1.90	0.53
1:D:116:ASP:OD1	2:D:501:HOH:O	2.19	0.53
1:A:353:LYS:HB2	1:A:355:LYS:HE3	1.90	0.53
1:A:398:ARG:HG3	1:A:398:ARG:O	2.08	0.52
1:C:26:LEU:HB3	1:C:175:GLU:HG3	1.91	0.52
1:C:306:GLU:HG3	2:C:599:HOH:O	2.11	0.51
1:D:261:VAL:HG13	1:D:266:CYS:HB3	1.93	0.51
1:B:216:ALA:HB2	1:B:333:VAL:HG21	1.91	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:85:GLU:HG2	2:A:826:HOH:O	2.11	0.50
1:A:324:ILE:HD12	1:A:349:TYR:CG	2.46	0.50
1:C:366:GLU:OE2	2:C:501:HOH:O	2.19	0.49
1:C:38:PHE:HB3	1:C:44:LYS:HG2	1.94	0.49
1:B:258:GLU:O	1:B:262:ARG:HG3	2.13	0.49
1:A:348:ARG:O	1:A:352:MET:HB2	2.13	0.48
1:A:312:GLU:O	1:A:334:THR:HA	2.13	0.48
1:D:140:HIS:CG	1:D:368:GLN:HG3	2.48	0.48
1:A:116:ASP:O	1:A:117:CYS:HB3	2.13	0.48
1:B:274:GLU:O	1:B:275:LEU:HB2	2.13	0.48
1:B:31:ARG:NH2	2:B:523:HOH:O	2.46	0.48
1:A:274:GLU:O	1:A:275:LEU:HB2	2.13	0.48
1:B:237:ALA:O	1:B:241:VAL:HG23	2.13	0.48
1:A:208:ARG:HG3	1:A:367:GLU:OE2	2.14	0.47
1:D:285:PRO:HA	1:D:288:VAL:CG1	2.44	0.47
1:C:289:VAL:HG23	2:C:502:HOH:O	2.13	0.47
1:C:417:GLU:HG2	1:C:420:LEU:HD11	1.95	0.47
1:D:317:VAL:O	2:D:502:HOH:O	2.20	0.47
1:D:140:HIS:HB2	1:D:141:PRO:HD2	1.97	0.47
1:A:261:VAL:HG13	1:A:266:CYS:HB3	1.97	0.46
2:B:522:HOH:O	1:C:352:MET:HA	2.13	0.46
1:C:45:ASP:HB3	1:C:48:LYS:HG3	1.98	0.46
1:D:254:SER:OG	1:D:257:LYS:HG3	2.15	0.46
1:A:428:ARG:NH2	1:A:433:GLU:HG3	2.31	0.46
1:A:324:ILE:HD12	1:A:349:TYR:CD2	2.51	0.46
1:C:228:GLY:H	1:C:272:ARG:HH22	1.64	0.45
1:B:198:VAL:O	1:B:408:LYS:HE2	2.17	0.45
1:D:312:GLU:O	1:D:334:THR:HA	2.16	0.45
1:A:270:ILE:HD13	1:A:302:LYS:HE2	1.97	0.45
1:A:138:ILE:HD13	1:A:187:LYS:HA	1.99	0.45
1:B:98:ARG:NH2	2:B:528:HOH:O	2.49	0.44
1:B:229:ALA:HB1	1:B:238:ILE:HD11	2.00	0.44
1:C:187:LYS:HE3	2:C:538:HOH:O	2.16	0.44
1:C:228:GLY:H	1:C:272:ARG:NH2	2.15	0.44
1:C:242:LYS:HA	1:C:242:LYS:HD2	1.72	0.44
1:D:261:VAL:HG12	1:D:266:CYS:HB3	1.99	0.43
1:B:46:VAL:HG11	1:B:397:CYS:SG	2.57	0.43
1:D:75:ASN:OD1	1:D:408:LYS:NZ	2.51	0.43
1:B:132:LYS:HB3	1:B:132:LYS:HE3	1.73	0.43
1:B:258:GLU:OE2	1:B:262:ARG:NH1	2.50	0.43
1:D:95:GLN:HB3	2:D:664:HOH:O	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:31:ARG:HH21	1:A:51:ARG:NE	2.17	0.43
1:B:348:ARG:HB3	1:B:352:MET:CE	2.49	0.43
1:C:195:GLU:O	1:C:198:VAL:HG22	2.18	0.43
1:A:106:HIS:CE1	1:A:168:THR:HB	2.54	0.43
1:A:62:PRO:HD2	2:A:505:HOH:O	2.18	0.43
1:B:404:ARG:HD3	2:B:634:HOH:O	2.18	0.43
1:C:139:VAL:HG22	1:C:184:LEU:HD23	2.00	0.43
1:A:126:ILE:HD12	2:A:885:HOH:O	2.19	0.42
1:C:139:VAL:CG1	1:C:179:VAL:HG11	2.46	0.42
1:B:198:VAL:CG1	1:B:410:ALA:HB2	2.50	0.42
1:A:270:ILE:HD11	1:A:298:LEU:HB3	2.01	0.42
1:A:353:LYS:HD2	2:A:823:HOH:O	2.19	0.42
1:D:224:VAL:HG22	1:D:309:ILE:HB	2.00	0.42
1:C:318:THR:HA	2:C:506:HOH:O	2.19	0.42
1:B:116:ASP:O	1:B:117:CYS:CB	2.65	0.42
1:B:346:ASP:N	2:B:537:HOH:O	2.52	0.42
1:B:347:ASN:OD1	1:C:336:GLY:HA3	2.19	0.42
1:A:258:GLU:HG3	1:A:269:VAL:HG11	2.01	0.42
1:A:38:PHE:HB3	1:A:44:LYS:HG2	1.99	0.42
1:D:256:GLN:H	1:D:256:GLN:HG2	1.56	0.42
1:D:264:LEU:CD1	2:D:737:HOH:O	2.44	0.42
1:A:76:TYR:HE1	2:A:586:HOH:O	2.03	0.42
1:D:230:SER:HB2	1:D:406:VAL:HG22	2.01	0.42
1:A:354:LEU:HD23	1:D:360:SER:HA	2.01	0.41
1:B:258:GLU:HG3	1:B:269:VAL:HG11	2.02	0.41
1:C:75:ASN:ND2	1:C:116:ASP:OD1	2.53	0.41
1:A:314:THR:HG23	2:A:654:HOH:O	2.21	0.41
1:D:86:PRO:O	1:D:87:ILE:HD13	2.20	0.41
1:B:254:SER:OG	1:B:256:GLN:HG2	2.21	0.41
1:D:79:VAL:HG12	1:D:83:MET:CE	2.49	0.41
1:D:253:SER:HB3	1:D:272:ARG:HD3	2.03	0.41
1:A:47:ARG:NE	1:A:398:ARG:HD2	2.36	0.41
1:C:31:ARG:NH1	1:C:51:ARG:HH12	2.19	0.41
1:B:274:GLU:O	1:B:275:LEU:CB	2.68	0.40
1:A:301:GLU:HG3	2:A:682:HOH:O	2.21	0.40
1:C:116:ASP:O	1:C:117:CYS:HB3	2.22	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	419/447 (94%)	404 (96%)	15 (4%)	0	100	100
1	B	422/447 (94%)	405 (96%)	16 (4%)	1 (0%)	49	57
1	C	409/447 (92%)	396 (97%)	13 (3%)	0	100	100
1	D	409/447 (92%)	394 (96%)	15 (4%)	0	100	100
All	All	1659/1788 (93%)	1599 (96%)	59 (4%)	1 (0%)	53	62

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	275	LEU

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	322/350 (92%)	312 (97%)	10 (3%)	43	52
1	B	327/350 (93%)	311 (95%)	16 (5%)	27	30
1	C	317/350 (91%)	309 (98%)	8 (2%)	50	60
1	D	314/350 (90%)	306 (98%)	8 (2%)	50	60
All	All	1280/1400 (91%)	1238 (97%)	42 (3%)	41	50

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	33	GLU
1	A	126	ILE
1	A	138	ILE
1	A	161	GLN
1	A	256	GLN
1	A	261	VAL
1	A	267	ASP
1	A	327	ARG
1	A	406	VAL
1	A	417	GLU
1	B	33	GLU
1	B	85	GLU
1	B	116	ASP
1	B	117	CYS
1	B	126	ILE
1	B	132	LYS
1	B	138	ILE
1	B	256	GLN
1	B	274	GLU
1	B	275	LEU
1	B	312	GLU
1	B	327	ARG
1	B	328	ARG
1	B	347	ASN
1	B	417	GLU
1	B	433	GLU
1	C	116	ASP
1	C	132	LYS
1	C	287	ARG
1	C	398	ARG
1	C	403	SER
1	C	405	GLN
1	C	434	ASP
1	C	435	ARG
1	D	51	ARG
1	D	117	CYS
1	D	148	GLU
1	D	256	GLN
1	D	288	VAL
1	D	406	VAL
1	D	426	ASP
1	D	430	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no

such sidechains identified.

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 carbohydrates 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	424/447 (94%)	0.20	6 (1%) 75 78	9, 17, 34, 53	1 (0%)
1	B	428/447 (95%)	0.17	9 (2%) 63 67	9, 17, 37, 57	0
1	C	417/447 (93%)	0.06	4 (0%) 82 84	8, 20, 40, 56	1 (0%)
1	D	417/447 (93%)	0.01	3 (0%) 87 89	11, 17, 34, 52	1 (0%)
All	All	1686/1788 (94%)	0.11	22 (1%) 77 80	8, 18, 36, 57	3 (0%)

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	275	LEU	6.1
1	D	285	PRO	3.9
1	C	275	LEU	3.7
1	B	349	TYR	3.4
1	A	349	TYR	3.3
1	A	275	LEU	3.2
1	A	445	THR	3.2
1	C	285	PRO	3.1
1	B	314	THR	2.8
1	B	315	GLY	2.8
1	B	337	SER	2.7
1	A	287	ARG	2.6
1	B	346	ASP	2.5
1	C	433	GLU	2.4
1	B	348	ARG	2.3
1	B	289	VAL	2.3
1	B	285	PRO	2.2
1	D	424	ASP	2.2
1	A	348	ARG	2.2
1	C	1	SER	2.1
1	A	314	THR	2.1
1	B	287	ARG	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 carbohydrates 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.