



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 04:21 PM GMT

PDB ID : 4EIT
Title : Crystal structure of an enoyl-(acyl carrier protein) reductase from Bartonella henselae
Authors : Seattle Structural Genomics Center for Infectious Disease (SSGCID)
Deposited on : 2012-04-05
Resolution : 2.40 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
<http://wwpdb.org/validation/2016/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
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

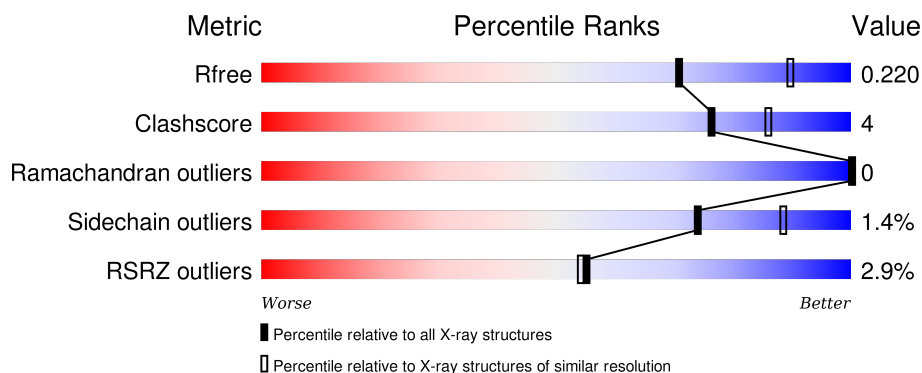
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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}	91344	2919 (2.40-2.40)
Clashscore	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)
RSRZ outliers	91569	2928 (2.40-2.40)

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	276	<div> <div>7%</div> <div> <div></div> <div>83%</div> <div>7%</div> <div>10%</div> </div> </div>
1	B	276	<div> <div>2%</div> <div> <div></div> <div>81%</div> <div>8%</div> <div>10%</div> </div> </div>
1	C	276	<div> <div>2%</div> <div> <div></div> <div>82%</div> <div>9%</div> <div>9%</div> </div> </div>
1	D	276	<div> <div>%</div> <div> <div></div> <div>83%</div> <div>7%</div> <div>10%</div> </div> </div>
1	E	276	<div> <div>%</div> <div> <div></div> <div>82%</div> <div>8%</div> <div>10%</div> </div> </div>

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
1	F	276	<div><div></div><div>3%</div><div>83%</div><div>7%</div><div>9%</div></div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 11186 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 Enoyl-[acyl-carrier-protein] reductase [NADH].

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	249	Total	C	N	O	S	Se	0	0	0
			1799	1148	297	345	2	7			
1	B	248	Total	C	N	O	S	Se	0	1	0
			1819	1159	299	351	2	8			
1	C	250	Total	C	N	O	S	Se	0	0	0
			1848	1175	310	354	2	7			
1	D	249	Total	C	N	O	S	Se	0	1	0
			1857	1184	311	352	2	8			
1	E	248	Total	C	N	O	S	Se	0	1	0
			1843	1175	307	351	2	8			
1	F	250	Total	C	N	O	S	Se	0	1	0
			1821	1156	304	351	2	8			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	GLY	-	EXPRESSION TAG	UNP Q6G4D5
A	-2	PRO	-	EXPRESSION TAG	UNP Q6G4D5
A	-1	GLY	-	EXPRESSION TAG	UNP Q6G4D5
A	0	SER	-	EXPRESSION TAG	UNP Q6G4D5
B	-3	GLY	-	EXPRESSION TAG	UNP Q6G4D5
B	-2	PRO	-	EXPRESSION TAG	UNP Q6G4D5
B	-1	GLY	-	EXPRESSION TAG	UNP Q6G4D5
B	0	SER	-	EXPRESSION TAG	UNP Q6G4D5
C	-3	GLY	-	EXPRESSION TAG	UNP Q6G4D5
C	-2	PRO	-	EXPRESSION TAG	UNP Q6G4D5
C	-1	GLY	-	EXPRESSION TAG	UNP Q6G4D5
C	0	SER	-	EXPRESSION TAG	UNP Q6G4D5
D	-3	GLY	-	EXPRESSION TAG	UNP Q6G4D5
D	-2	PRO	-	EXPRESSION TAG	UNP Q6G4D5
D	-1	GLY	-	EXPRESSION TAG	UNP Q6G4D5
D	0	SER	-	EXPRESSION TAG	UNP Q6G4D5
E	-3	GLY	-	EXPRESSION TAG	UNP Q6G4D5

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
E	-2	PRO	-	EXPRESSION TAG	UNP Q6G4D5
E	-1	GLY	-	EXPRESSION TAG	UNP Q6G4D5
E	0	SER	-	EXPRESSION TAG	UNP Q6G4D5
F	-3	GLY	-	EXPRESSION TAG	UNP Q6G4D5
F	-2	PRO	-	EXPRESSION TAG	UNP Q6G4D5
F	-1	GLY	-	EXPRESSION TAG	UNP Q6G4D5
F	0	SER	-	EXPRESSION TAG	UNP Q6G4D5

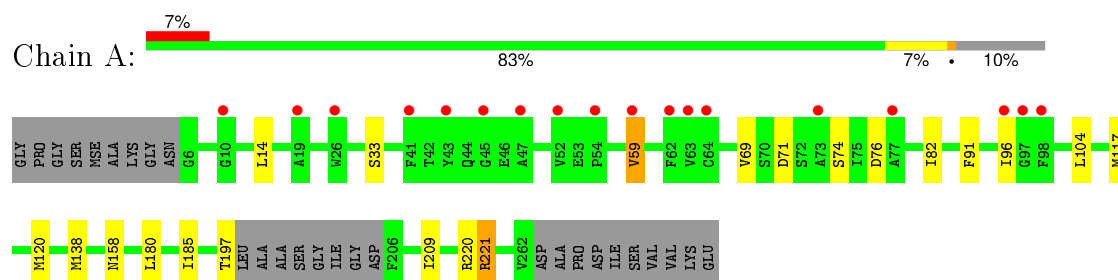
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	24	Total O 24 24	0	0
2	B	31	Total O 31 31	0	0
2	C	38	Total O 38 38	0	0
2	D	36	Total O 36 36	0	0
2	E	36	Total O 36 36	0	0
2	F	34	Total O 34 34	0	0

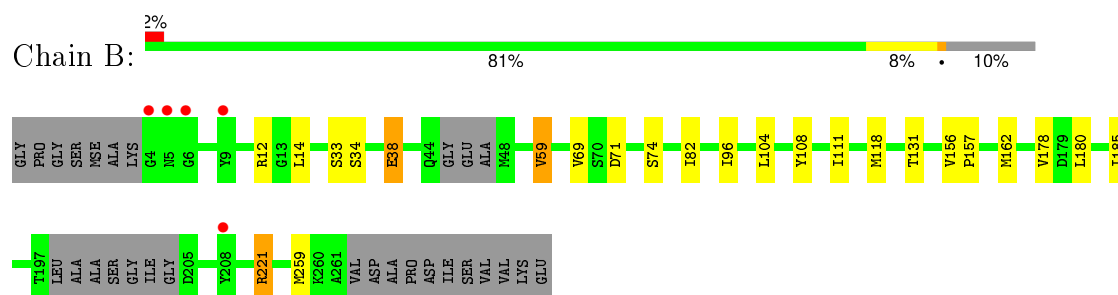
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

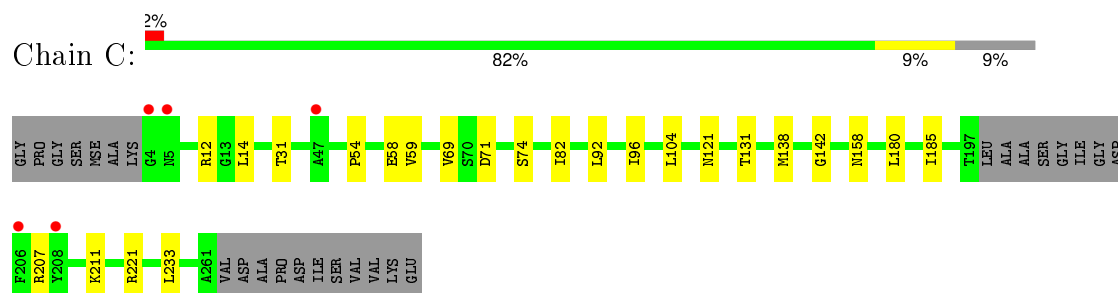
- Molecule 1: Enoyl-[acyl-carrier-protein] reductase [NADH]



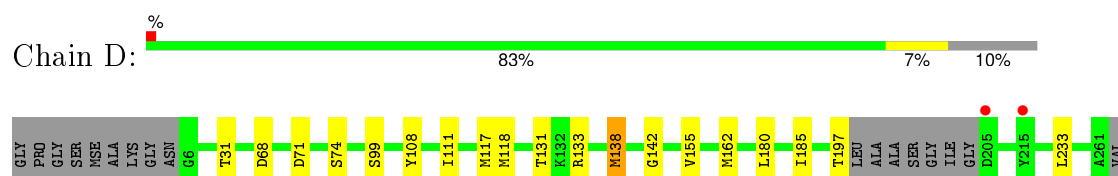
- Molecule 1: Enoyl-[acyl-carrier-protein] reductase [NADH]



- Molecule 1: Enoyl-[acyl-carrier-protein] reductase [NADH]

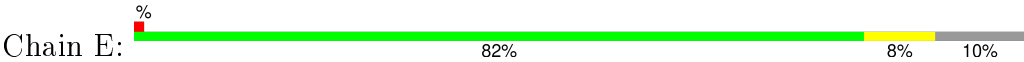


- Molecule 1: Enoyl-[acyl-carrier-protein] reductase [NADH]



ASP
ALA
PRO
ASP
ILE
SER
VAL
LYS
GLU

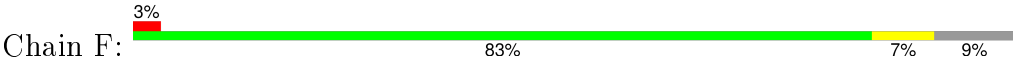
- Molecule 1: Enoyl-[acyl-carrier-protein] reductase [NADH]



GLY PRO GLY SER MSE ALA LYS GLY
MS L14 L31 SS4 C67 D68 V69 S70 D71 S74 V78 I82 L89 I96 Y108 I111 T131 M138 V156 P157 N158 L180 K183 H184 I185 T195
LYS THR LEU ALA ALA SER GLY ILE GLY

D206 F206 R207 Y208 L233 A261 VAL ASP ALA PRO ASP ILE SER VAL VAL LYS GLU

- Molecule 1: Enoyl-[acyl-carrier-protein] reductase [NADH]



GLY PRO GLY SER MSE ALA LYS GLY ASN G6 L14 G45 E46 A47 R51 V52 V59 C64 V69 S70 D71 S74 I82 T96 G97 F98 S99 F116 M117 M118 T119 M120 M138 S139 G142 L147 M162 G163 V164 L180 T185 T197
LEU

ALA ALA SER GLY ILE GLY D205 Y215 R221 V262 ASP ALA PRO ASP ILE SER VAL VAL LYS GLU

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	122.36 Å 76.86 Å 171.95 Å 90.00° 107.99° 90.00°	Depositor
Resolution (Å)	50.00 – 2.40 47.83 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.1 (50.00-2.40) 99.1 (47.83-2.40)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.20 (at 2.39 Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.196 , 0.226 0.189 , 0.220	Depositor DCC
R_{free} test set	2994 reflections (5.34%)	DCC
Wilson B-factor (Å ²)	38.0	Xtriage
Anisotropy	0.074	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 41.6	EDS
Estimated twinning fraction	0.017 for h,-k,-h-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 59082 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	11186	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.22% 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, 0.375 respectively for untwinned datasets, and 0.333, 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	0.66	0/1824	0.77	5/2474 (0.2%)
1	B	0.68	0/1845	0.74	2/2501 (0.1%)
1	C	0.66	0/1874	0.73	0/2535
1	D	0.72	0/1883	0.79	3/2543 (0.1%)
1	E	0.72	0/1872	0.78	2/2530 (0.1%)
1	F	0.70	0/1848	0.80	4/2504 (0.2%)
All	All	0.69	0/11146	0.77	16/15087 (0.1%)

There are no bond length outliers.

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	221	ARG	NE-CZ-NH1	-9.93	115.33	120.30
1	E	138	MSE	CA-CB-CG	-9.52	97.12	113.30
1	D	138	MSE	CA-CB-CG	-9.21	97.64	113.30
1	A	221	ARG	NE-CZ-NH2	8.00	124.30	120.30
1	D	117	MSE	CG-SE-CE	-7.33	82.78	98.90
1	F	147	LEU	CB-CG-CD1	-6.94	99.20	111.00
1	F	147	LEU	CB-CG-CD2	6.70	122.39	111.00
1	F	215	TYR	CA-CB-CG	5.87	124.54	113.40
1	A	220	ARG	NE-CZ-NH2	5.86	123.23	120.30
1	D	68	ASP	CB-CG-OD1	-5.75	113.13	118.30
1	F	138	MSE	CG-SE-CE	-5.39	87.05	98.90
1	B	221	ARG	NE-CZ-NH2	-5.24	117.68	120.30
1	A	220	ARG	NE-CZ-NH1	-5.24	117.68	120.30
1	B	221	ARG	NE-CZ-NH1	5.11	122.86	120.30
1	A	120	MSE	CG-SE-CE	-5.04	87.82	98.90
1	E	89	LEU	CB-CG-CD1	-5.01	102.47	111.00

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	1799	0	1710	12	0
1	B	1819	0	1722	18	0
1	C	1848	0	1779	20	0
1	D	1857	0	1813	11	0
1	E	1843	0	1793	9	0
1	F	1821	0	1738	16	0
2	A	24	0	0	0	0
2	B	31	0	0	0	0
2	C	38	0	0	0	0
2	D	36	0	0	0	0
2	E	36	0	0	0	0
2	F	34	0	0	0	0
All	All	11186	0	10555	82	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 (82) 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:156:VAL:HG13	1:B:157:PRO:HD2	1.42	1.00
1:C:138:MSE:HE3	1:C:142:GLY:HA3	1.44	0.98
1:F:138:MSE:HE3	1:F:142:GLY:HA3	1.52	0.91
1:F:116:PHE:CD2	1:F:117:MSE:HE2	2.17	0.80
1:F:138:MSE:CE	1:F:142:GLY:HA3	2.16	0.75
1:C:138:MSE:HE1	1:C:185:ILE:HG23	1.71	0.73
1:B:156:VAL:CG1	1:B:157:PRO:HD2	2.19	0.69
1:A:117:MSE:HE1	1:C:121:ASN:HA	1.76	0.67
1:A:69:VAL:HG11	1:A:96:ILE:HD13	1.77	0.66
1:A:117:MSE:HE3	1:C:121:ASN:OD1	1.96	0.65
1:C:138:MSE:HE2	1:C:185:ILE:HG12	1.81	0.62
1:C:138:MSE:CE	1:C:185:ILE:HG23	2.31	0.60
1:B:156:VAL:HG13	1:B:157:PRO:CD	2.23	0.60
1:B:33:SER:CB	1:B:59:VAL:HG22	2.32	0.59
1:C:69:VAL:HG11	1:C:96:ILE:HD13	1.86	0.58

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:209:ILE:HG23	1:B:259:MSE:HE1	1.85	0.57
1:B:14:LEU:HD13	1:B:82:ILE:HD11	1.87	0.57
1:F:116:PHE:CZ	1:F:120:MSE:HE2	2.41	0.56
1:D:131:THR:CG2	1:D:180:LEU:HD11	2.36	0.55
1:C:31:THR:HG22	1:C:233:LEU:HD22	1.89	0.55
1:A:59:VAL:HG13	1:A:59:VAL:O	2.07	0.55
1:B:156:VAL:CG1	1:B:157:PRO:CD	2.84	0.53
1:B:59:VAL:HG13	1:B:59:VAL:O	2.08	0.53
1:A:33:SER:CB	1:A:59:VAL:HG22	2.39	0.53
1:F:59:VAL:HG13	1:F:59:VAL:O	2.07	0.53
1:F:14:LEU:HD13	1:F:82:ILE:HD11	1.90	0.53
1:D:99:SER:HB2	1:D:162:MSE:HE2	1.91	0.53
1:F:118[B]:MSE:HE2	1:F:118[B]:MSE:HA	1.91	0.52
1:C:207:ARG:O	1:C:211:LYS:HG2	2.09	0.52
1:E:69:VAL:HG11	1:E:96:ILE:HD13	1.92	0.52
1:A:91:PHE:HA	1:A:138:MSE:CE	2.40	0.51
1:C:138:MSE:HE3	1:C:142:GLY:CA	2.29	0.51
1:D:31:THR:HG22	1:D:233:LEU:HD22	1.92	0.50
1:F:99:SER:HB2	1:F:162:MSE:HE2	1.93	0.50
1:F:69:VAL:HG11	1:F:96:ILE:HD13	1.93	0.49
1:B:108:TYR:O	1:B:111:ILE:HG23	2.13	0.49
1:D:108:TYR:O	1:D:111:ILE:HG23	2.13	0.48
1:E:108:TYR:O	1:E:111:ILE:HG23	2.13	0.48
1:F:120:MSE:HE1	1:F:164:VAL:HG12	1.96	0.48
1:F:59:VAL:CG1	1:F:59:VAL:O	2.63	0.47
1:B:178:VAL:HG21	1:D:155:VAL:HG11	1.96	0.47
1:C:59:VAL:HG13	1:C:59:VAL:O	2.14	0.46
1:C:138:MSE:CE	1:C:142:GLY:HA3	2.32	0.46
1:A:180:LEU:HB3	1:A:185:ILE:HB	1.97	0.45
1:F:116:PHE:HD2	1:F:117:MSE:HE2	1.73	0.45
1:B:131:THR:CG2	1:B:180:LEU:HD11	2.46	0.45
1:F:138:MSE:HE2	1:F:185:ILE:HG12	1.99	0.45
1:B:69:VAL:HG11	1:B:96:ILE:HD13	1.99	0.45
1:E:156:VAL:HG13	1:E:157:PRO:HD2	1.98	0.45
1:F:180:LEU:HB3	1:F:185:ILE:HB	1.98	0.44
1:C:14:LEU:HD22	1:C:82:ILE:CD1	2.48	0.44
1:D:180:LEU:HB3	1:D:185:ILE:HB	2.00	0.44
1:B:180:LEU:HB3	1:B:185:ILE:HB	1.99	0.44
1:A:59:VAL:CG1	1:A:59:VAL:O	2.66	0.44
1:E:131:THR:CG2	1:E:180:LEU:HD11	2.48	0.44
1:C:180:LEU:HB3	1:C:185:ILE:HB	1.99	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:71:ASP:O	1:A:74:SER:HB3	2.19	0.43
1:C:59:VAL:O	1:C:59:VAL:CG1	2.66	0.43
1:B:71:ASP:O	1:B:74:SER:HB3	2.18	0.43
1:D:99:SER:CB	1:D:162:MSE:HE2	2.48	0.43
1:E:180:LEU:HB3	1:E:185:ILE:HB	1.99	0.43
1:A:14:LEU:HD22	1:A:82:ILE:CD1	2.49	0.42
1:E:71:ASP:O	1:E:74:SER:HB3	2.19	0.42
1:C:92:LEU:C	1:C:92:LEU:HD23	2.38	0.42
1:F:99:SER:CB	1:F:162:MSE:HE2	2.50	0.42
1:B:12:ARG:HG3	1:B:38:GLU:HG3	2.02	0.42
1:F:71:ASP:O	1:F:74:SER:HB3	2.19	0.42
1:E:31:THR:HG22	1:E:233:LEU:HD22	2.01	0.42
1:D:118[B]:MSE:HA	1:D:118[B]:MSE:CE	2.49	0.42
1:E:14:LEU:HD13	1:E:82:ILE:HD11	2.01	0.42
1:B:59:VAL:CG1	1:B:59:VAL:O	2.68	0.41
1:D:138:MSE:CE	1:D:142:GLY:HA3	2.50	0.41
1:D:71:ASP:O	1:D:74:SER:HB3	2.20	0.41
1:B:104:LEU:HD23	1:B:104:LEU:HA	1.86	0.41
1:C:104:LEU:O	1:C:158:ASN:HA	2.20	0.41
1:C:71:ASP:O	1:C:74:SER:HB3	2.20	0.41
1:A:104:LEU:O	1:A:158:ASN:HA	2.21	0.41
1:C:131:THR:CG2	1:C:180:LEU:HD11	2.50	0.41
1:E:67:CYS:HB2	1:E:78:VAL:HG21	2.03	0.41
1:B:118[B]:MSE:O	1:B:118[B]:MSE:HG3	2.20	0.40
1:D:133:ARG:HA	1:D:133:ARG:HD2	1.84	0.40
1:C:54:PRO:O	1:C:58:GLU:HG2	2.21	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	245/276 (89%)	238 (97%)	7 (3%)	0	100	100
1	B	243/276 (88%)	237 (98%)	6 (2%)	0	100	100
1	C	246/276 (89%)	240 (98%)	6 (2%)	0	100	100
1	D	246/276 (89%)	240 (98%)	6 (2%)	0	100	100
1	E	245/276 (89%)	239 (98%)	6 (2%)	0	100	100
1	F	247/276 (90%)	240 (97%)	7 (3%)	0	100	100
All	All	1472/1656 (89%)	1434 (97%)	38 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	176/212 (83%)	172 (98%)	4 (2%)	58	78
1	B	180/212 (85%)	175 (97%)	5 (3%)	51	72
1	C	185/212 (87%)	183 (99%)	2 (1%)	80	92
1	D	188/212 (89%)	187 (100%)	1 (0%)	92	97
1	E	187/212 (88%)	185 (99%)	2 (1%)	80	92
1	F	181/212 (85%)	180 (99%)	1 (1%)	90	96
All	All	1097/1272 (86%)	1082 (99%)	15 (1%)	74	88

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	59	VAL
1	A	76	ASP
1	A	197	THR
1	A	221	ARG
1	B	34	SER
1	B	38	GLU
1	B	59	VAL

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	162	MSE
1	B	221	ARG
1	C	12	ARG
1	C	221	ARG
1	D	197	THR
1	E	34	SER
1	E	183	LYS
1	F	221	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	44	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 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(Å²)	Q<0.9
1	A	242/276 (87%)	0.28	18 (7%)	17	17	22, 41, 66, 87	0
1	B	241/276 (87%)	-0.12	5 (2%)	67	66	19, 34, 57, 72	0
1	C	243/276 (88%)	-0.11	5 (2%)	67	66	20, 35, 60, 77	0
1	D	242/276 (87%)	-0.18	2 (0%)	87	87	20, 31, 57, 78	0
1	E	241/276 (87%)	-0.11	4 (1%)	73	72	17, 30, 53, 63	0
1	F	243/276 (88%)	0.10	8 (3%)	50	50	18, 34, 59, 81	0
All	All	1452/1656 (87%)	-0.02	42 (2%)	55	54	17, 34, 61, 87	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	47	ALA	5.7
1	B	9	TYR	4.5
1	A	45	GLY	4.3
1	A	62	PHE	3.9
1	E	205	ASP	3.9
1	C	47	ALA	3.6
1	A	64	CYS	3.6
1	F	64	CYS	3.2
1	B	5	ASN	3.2
1	F	45	GLY	3.1
1	A	10	GLY	3.1
1	A	77	ALA	2.9
1	F	52	VAL	2.9
1	A	59	VAL	2.8
1	C	4	GLY	2.6
1	F	139	SER	2.6
1	A	73	ALA	2.6
1	D	215	TYR	2.6
1	A	26	TRP	2.5

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	208	TYR	2.5
1	A	41	PHE	2.4
1	C	208	TYR	2.4
1	B	6	GLY	2.4
1	A	97	GLY	2.4
1	A	63	VAL	2.4
1	C	5	ASN	2.3
1	A	19	ALA	2.3
1	A	96	ILE	2.3
1	F	51	ARG	2.3
1	A	43	TYR	2.3
1	D	205	ASP	2.3
1	A	52	VAL	2.3
1	F	47	ALA	2.2
1	A	54	PRO	2.2
1	E	208	TYR	2.2
1	C	206	PHE	2.2
1	B	4	GLY	2.1
1	E	206	PHE	2.1
1	A	98	PHE	2.1
1	F	98	PHE	2.0
1	E	158	ASN	2.0
1	F	215	TYR	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.