



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 12:41 PM UTC

PDB ID : 4QFW / pdb_00004qfw
Title : Crystal structure of acyl-CoA thioesterase tesB from Yersinia pestis
Authors : Swarbrick, C.M.D.; Forwood, J.K.
Deposited on : 2014-05-21
Resolution : 2.00 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
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

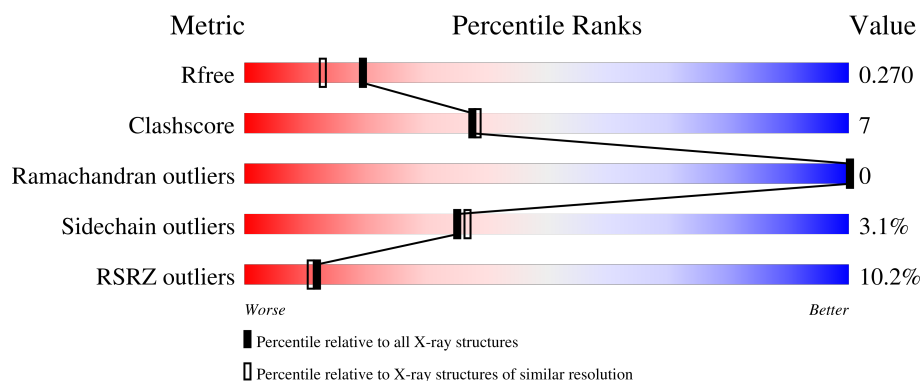
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	288	4% 83% 7% 9%
1	B	288	7% 77% 15% 8%
1	C	288	6% 78% 15% 7%
1	D	288	20% 72% 19% 9%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8667 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 Acyl-CoA thioesterase II.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	263	Total	C	N	O	S	0	0	0
			2097	1334	367	389	7			
1	B	265	Total	C	N	O	S	0	0	0
			2114	1346	369	392	7			
1	C	269	Total	C	N	O	S	0	0	0
			2152	1373	376	396	7			
1	D	263	Total	C	N	O	S	0	0	0
			2097	1334	367	389	7			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	287	SER	-	expression tag	UNP Q0WCE2
A	288	ASN	-	expression tag	UNP Q0WCE2
B	287	SER	-	expression tag	UNP Q0WCE2
B	288	ASN	-	expression tag	UNP Q0WCE2
C	287	SER	-	expression tag	UNP Q0WCE2
C	288	ASN	-	expression tag	UNP Q0WCE2
D	287	SER	-	expression tag	UNP Q0WCE2
D	288	ASN	-	expression tag	UNP Q0WCE2

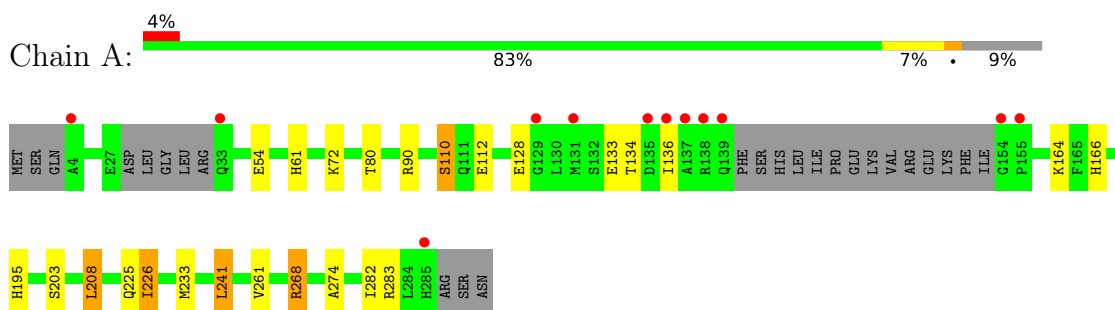
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	79	Total	O	0	0
			79	79		
2	B	42	Total	O	0	0
			42	42		
2	C	62	Total	O	0	0
			62	62		
2	D	24	Total	O	0	0
			24	24		

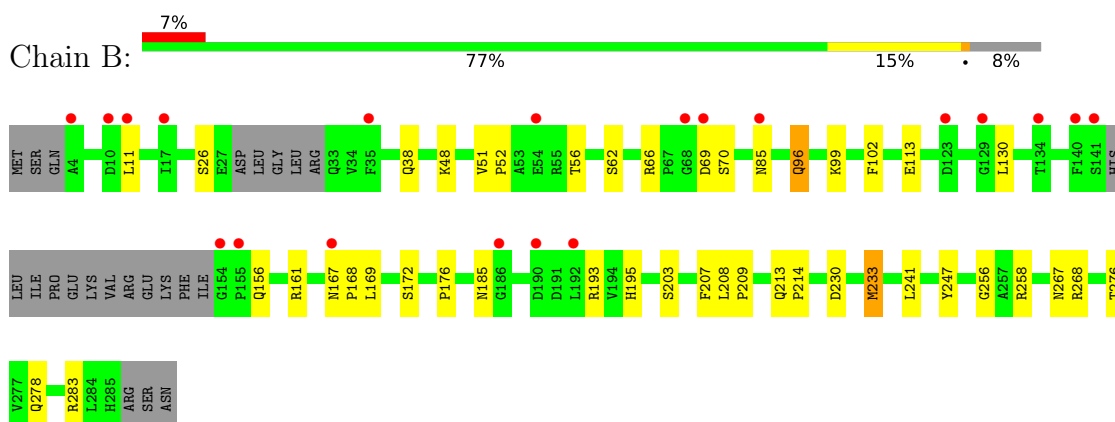
3 Residue-property plots [i](#)

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.

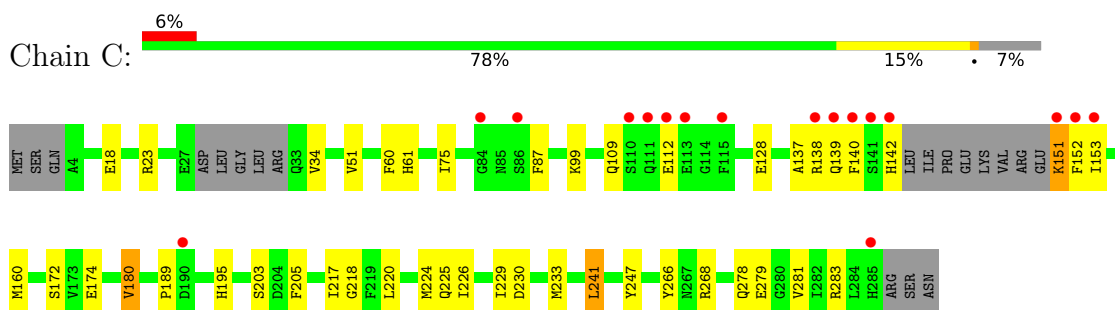
• Molecule 1: Acyl-CoA thioesterase II



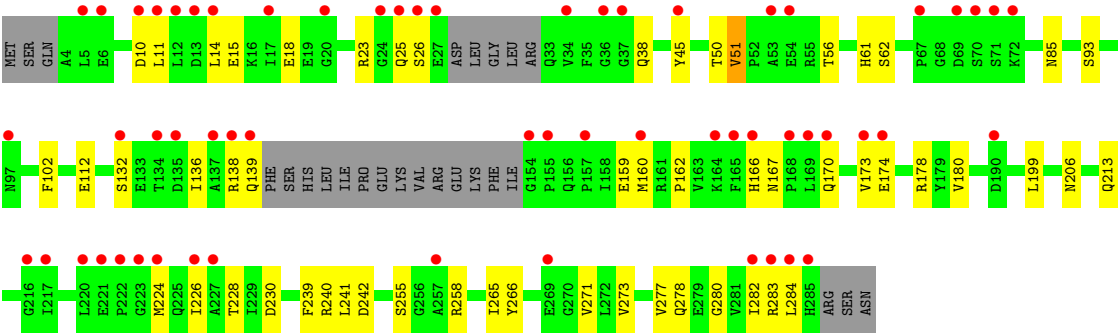
• Molecule 1: Acyl-CoA thioesterase II



• Molecule 1: Acyl-CoA thioesterase II



• Molecule 1: Acyl-CoA thioesterase II



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	51.23Å 171.62Å 73.70Å 90.00° 109.62° 90.00°	Depositor
Resolution (Å)	24.29 – 2.00 24.29 – 2.00	Depositor EDS
% Data completeness (in resolution range)	93.7 (24.29-2.00) 93.7 (24.29-2.00)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.36 (at 1.99Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.221 , 0.268 0.227 , 0.270	Depositor DCC
R_{free} test set	3827 reflections (4.75%)	wwPDB-VP
Wilson B-factor (Å ²)	22.9	Xtriage
Anisotropy	0.071	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 39.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.032 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	8667	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.36% 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	1.16	3/2152 (0.1%)	1.00	0/2915
1	B	1.14	2/2170 (0.1%)	1.06	2/2939 (0.1%)
1	C	1.21	1/2210 (0.0%)	1.10	4/2992 (0.1%)
1	D	1.15	5/2152 (0.2%)	1.04	2/2915 (0.1%)
All	All	1.16	11/8684 (0.1%)	1.05	8/11761 (0.1%)

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	50	THR	N-CA	6.23	1.54	1.46
1	B	233	MET	SD-CE	-6.10	1.64	1.79
1	A	274	ALA	N-CA	6.01	1.53	1.45
1	D	255	SER	CA-C	5.77	1.60	1.52
1	A	208	LEU	C-O	-5.54	1.19	1.24
1	C	266	TYR	CA-C	-5.47	1.46	1.52
1	A	261	VAL	C-O	5.21	1.29	1.24
1	D	162	PRO	N-CA	-5.12	1.41	1.46
1	B	213	GLN	C-O	-5.09	1.19	1.24
1	D	277	VAL	C-O	-5.05	1.19	1.24
1	D	160	MET	CA-C	5.01	1.58	1.52

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	153	ILE	N-CA-C	7.06	117.63	111.56
1	C	180	VAL	N-CA-C	6.31	117.38	108.17
1	C	218	GLY	N-CA-C	-5.91	104.09	112.37
1	B	258	ARG	CA-C-N	5.78	126.55	122.33
1	B	258	ARG	C-N-CA	5.78	126.55	122.33
1	C	229	ILE	N-CA-C	-5.26	106.68	111.67
1	D	51	VAL	CA-C-N	5.21	125.21	119.89
1	D	51	VAL	C-N-CA	5.21	125.21	119.89

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	2097	0	2024	19	0
1	B	2114	0	2038	38	0
1	C	2152	0	2078	35	0
1	D	2097	0	2024	35	0
2	A	79	0	0	0	0
2	B	42	0	0	0	0
2	C	62	0	0	1	0
2	D	24	0	0	0	0
All	All	8667	0	8164	117	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 7.

All (117) 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:C:138:ARG:O	1:C:139:GLN:HB2	1.68	0.90
1:B:26:SER:HA	1:B:38:GLN:HE22	1.35	0.89
1:B:69:ASP:H	1:B:96:GLN:HE22	1.21	0.85
1:A:203:SER:OG	1:A:233:MET:HE1	1.80	0.82
1:A:195:HIS:HD2	1:A:241:LEU:H	1.29	0.81
1:B:26:SER:HA	1:B:38:GLN:NE2	1.96	0.81
1:D:284:LEU:O	1:D:284:LEU:HD12	1.81	0.80
1:C:195:HIS:HD2	1:C:241:LEU:H	1.31	0.77
1:A:54:GLU:OE1	1:A:110:SER:HB3	1.86	0.76
1:C:99:LYS:HE2	2:C:320:HOH:O	1.85	0.75
1:C:112:GLU:OE1	1:D:283:ARG:NH1	2.22	0.73
1:B:11:LEU:HD21	1:B:38:GLN:NE2	2.03	0.72
1:A:195:HIS:CD2	1:A:241:LEU:H	2.08	0.72
1:B:203:SER:OG	1:B:233:MET:HE1	1.90	0.71
1:C:203:SER:OG	1:C:233:MET:HE1	1.91	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:136:ILE:O	1:D:139:GLN:HB3	1.92	0.69
1:D:10:ASP:OD1	1:D:11:LEU:N	2.26	0.68
1:A:203:SER:OG	1:A:233:MET:CE	2.41	0.67
1:D:166:HIS:ND1	1:D:174:GLU:OE2	2.29	0.64
1:C:195:HIS:CD2	1:C:241:LEU:H	2.15	0.64
1:D:15:GLU:OE1	1:D:23:ARG:NH2	2.31	0.63
1:D:178:ARG:HH11	1:D:206:ASN:HD22	1.46	0.63
1:C:151:LYS:HG3	1:C:220:LEU:HD11	1.80	0.62
1:A:112:GLU:OE1	1:B:283:ARG:NH2	2.32	0.62
1:C:160:MET:HE1	1:C:205:PHE:CB	2.30	0.62
1:A:128:GLU:OE2	1:A:268:ARG:NH2	2.33	0.62
1:C:225:GLN:NE2	1:C:283:ARG:HH21	1.97	0.61
1:C:138:ARG:O	1:C:139:GLN:CB	2.42	0.60
1:C:128:GLU:OE2	1:C:268:ARG:NH2	2.29	0.59
1:B:195:HIS:CD2	1:B:241:LEU:H	2.21	0.58
1:B:195:HIS:HD2	1:B:241:LEU:H	1.52	0.58
1:C:60:PHE:CZ	1:C:233:MET:HE3	2.39	0.58
1:D:166:HIS:CE1	1:D:174:GLU:OE2	2.57	0.57
1:B:268:ARG:HG2	1:B:268:ARG:HH11	1.70	0.57
1:D:226:ILE:HG22	1:D:282:ILE:HG22	1.86	0.57
1:C:217:ILE:HG21	1:C:224:MET:HE3	1.88	0.55
1:B:26:SER:CA	1:B:38:GLN:HE22	2.12	0.55
1:C:151:LYS:HD3	1:C:151:LYS:O	2.06	0.55
1:C:87:PHE:CE1	1:C:109:GLN:HG3	2.42	0.54
1:B:203:SER:CB	1:B:233:MET:HE1	2.37	0.54
1:C:172:SER:O	1:C:174:GLU:HG2	2.08	0.54
1:C:225:GLN:NE2	1:C:283:ARG:NH2	2.56	0.54
1:C:61:HIS:HE1	1:D:230:ASP:HB2	1.72	0.54
1:D:15:GLU:CD	1:D:23:ARG:HH21	2.17	0.53
1:A:226:ILE:HD12	1:A:282:ILE:HG22	1.90	0.53
1:B:203:SER:OG	1:B:233:MET:CE	2.56	0.52
1:D:11:LEU:HD11	1:D:25:GLN:O	2.09	0.52
1:B:130:LEU:O	1:B:161:ARG:HD3	2.09	0.52
1:D:167:ASN:ND2	1:D:170:GLN:OE1	2.43	0.52
1:B:167:ASN:ND2	1:B:169:LEU:H	2.08	0.51
1:A:133:GLU:HA	1:A:136:ILE:HG22	1.93	0.51
1:D:228:THR:HG23	1:D:278:GLN:CD	2.36	0.51
1:C:203:SER:OG	1:C:233:MET:CE	2.59	0.50
1:B:156:GLN:O	1:B:185:ASN:ND2	2.43	0.50
1:D:14:LEU:HB2	1:D:45:TYR:CE2	2.48	0.49
1:C:137:ALA:HA	1:C:140:PHE:CE1	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:199:LEU:HD22	1:D:239:PHE:CZ	2.48	0.49
1:A:128:GLU:CD	1:A:268:ARG:HH22	2.21	0.48
1:D:265:ILE:O	1:D:273:VAL:HG22	2.13	0.48
1:D:15:GLU:OE1	1:D:23:ARG:NE	2.46	0.48
1:C:225:GLN:HE22	1:C:283:ARG:NH2	2.12	0.48
1:D:167:ASN:HB3	1:D:170:GLN:HB3	1.96	0.48
1:A:112:GLU:CD	1:B:283:ARG:NH2	2.72	0.47
1:A:164:LYS:HD2	1:A:166:HIS:CE1	2.49	0.47
1:B:85:ASN:O	1:B:85:ASN:OD1	2.33	0.46
1:C:225:GLN:HE21	1:C:283:ARG:HD2	1.81	0.46
1:D:26:SER:HA	1:D:38:GLN:NE2	2.30	0.45
1:B:233:MET:HE2	1:B:276:THR:CG2	2.46	0.45
1:B:233:MET:HE2	1:B:276:THR:HG22	1.99	0.45
1:D:226:ILE:O	1:D:226:ILE:HG13	2.16	0.45
1:B:66:ARG:HE	1:B:99:LYS:NZ	2.13	0.45
1:B:11:LEU:HD21	1:B:38:GLN:HE21	1.80	0.45
1:D:138:ARG:CZ	1:D:138:ARG:HB2	2.47	0.45
1:B:48:LYS:HE3	1:B:193:ARG:HD2	1.98	0.45
1:B:66:ARG:CZ	1:B:99:LYS:HE2	2.47	0.45
1:B:69:ASP:N	1:B:96:GLN:HE22	2.02	0.45
1:D:174:GLU:HG3	1:D:213:GLN:HB3	1.99	0.45
1:D:132:SER:OG	1:D:159:GLU:OE2	2.36	0.44
1:A:112:GLU:OE2	1:B:283:ARG:NH2	2.48	0.44
1:C:60:PHE:HZ	1:C:233:MET:HE3	1.79	0.44
1:C:283:ARG:NE	1:D:112:GLU:OE2	2.50	0.44
1:C:142:HIS:CD2	1:C:142:HIS:H	2.36	0.44
1:D:62:SER:HB2	1:D:102:PHE:CE1	2.53	0.44
1:A:208:LEU:HG	1:A:226:ILE:HD11	1.99	0.44
1:A:225:GLN:OE1	1:A:283:ARG:NH2	2.42	0.44
1:C:226:ILE:HA	1:C:281:VAL:O	2.18	0.44
1:D:258:ARG:HA	1:D:280:GLY:O	2.18	0.44
1:B:167:ASN:HD22	1:B:169:LEU:H	1.66	0.44
1:A:112:GLU:O	1:B:256:GLY:HA2	2.18	0.44
1:C:230:ASP:HB2	1:D:61:HIS:HE1	1.83	0.44
1:C:128:GLU:CD	1:C:268:ARG:HH22	2.23	0.43
1:C:189:PRO:O	1:C:195:HIS:HE1	2.01	0.43
1:C:230:ASP:OD1	1:C:279:GLU:HB3	2.18	0.43
1:D:14:LEU:HB2	1:D:45:TYR:CD2	2.54	0.43
1:A:80:THR:HA	1:A:90:ARG:HD3	2.00	0.43
1:B:203:SER:CB	1:B:233:MET:CE	2.95	0.43
1:C:230:ASP:O	1:C:278:GLN:HG3	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:230:ASP:HB2	1:D:61:HIS:CE1	2.53	0.43
1:A:61:HIS:HE1	1:B:230:ASP:HB2	1.84	0.42
1:B:167:ASN:HA	1:B:168:PRO:HD3	1.85	0.42
1:B:230:ASP:O	1:B:278:GLN:HG3	2.19	0.42
1:D:266:TYR:HA	1:D:271:VAL:O	2.19	0.42
1:B:62:SER:HB2	1:B:102:PHE:CE1	2.54	0.42
1:B:176:PRO:HB3	1:B:214:PRO:HG3	2.01	0.42
1:B:51:VAL:HG13	1:B:52:PRO:HD2	2.02	0.42
1:A:203:SER:CB	1:A:233:MET:CE	2.98	0.42
1:D:62:SER:HB2	1:D:102:PHE:CZ	2.55	0.42
1:B:208:LEU:N	1:B:209:PRO:CD	2.83	0.41
1:B:267:ASN:OD1	1:B:267:ASN:C	2.63	0.41
1:C:225:GLN:HE22	1:C:283:ARG:HH21	1.63	0.41
1:C:23:ARG:HB2	1:C:75:ILE:HD13	2.02	0.41
1:B:207:PHE:CE1	1:B:278:GLN:HB3	2.55	0.41
1:B:268:ARG:HH11	1:B:268:ARG:CG	2.32	0.41
1:C:151:LYS:CG	1:C:220:LEU:HD11	2.49	0.41
1:D:10:ASP:OD1	1:D:10:ASP:C	2.64	0.40
1:D:224:MET:HE2	1:D:224:MET:HA	2.03	0.40
1:D:240:ARG:HB3	1:D:242:ASP:OD1	2.21	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	257/288 (89%)	253 (98%)	4 (2%)	0	100	100
1	B	259/288 (90%)	254 (98%)	5 (2%)	0	100	100
1	C	263/288 (91%)	259 (98%)	4 (2%)	0	100	100
1	D	257/288 (89%)	250 (97%)	7 (3%)	0	100	100
All	All	1036/1152 (90%)	1016 (98%)	20 (2%)	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	224/248 (90%)	218 (97%)	6 (3%)	39	42
1	B	226/248 (91%)	220 (97%)	6 (3%)	39	42
1	C	230/248 (93%)	222 (96%)	8 (4%)	32	32
1	D	224/248 (90%)	216 (96%)	8 (4%)	31	31
All	All	904/992 (91%)	876 (97%)	28 (3%)	35	37

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	72	LYS
1	A	110	SER
1	A	134	THR
1	A	226	ILE
1	A	241	LEU
1	A	268	ARG
1	B	56	THR
1	B	70	SER
1	B	96	GLN
1	B	113	GLU
1	B	172	SER
1	B	247	TYR
1	C	18	GLU
1	C	34	VAL
1	C	51	VAL
1	C	151	LYS
1	C	152	PHE
1	C	180	VAL
1	C	241	LEU
1	C	247	TYR
1	D	18	GLU
1	D	51	VAL

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Mol	Chain	Res	Type
1	D	56	THR
1	D	85	ASN
1	D	93	SER
1	D	173	VAL
1	D	180	VAL
1	D	241	LEU

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	111	GLN
1	A	170	GLN
1	A	195	HIS
1	A	196	GLN
1	A	215	HIS
1	A	285	HIS
1	B	38	GLN
1	B	49	GLN
1	B	96	GLN
1	B	195	HIS
1	B	215	HIS
1	C	38	GLN
1	C	42	GLN
1	C	49	GLN
1	C	139	GLN
1	C	142	HIS
1	C	156	GLN
1	C	195	HIS
1	C	215	HIS
1	C	225	GLN
1	C	285	HIS
1	D	111	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	263/288 (91%)	0.21	12 (4%)	37 36	11, 24, 48, 97	0
1	B	265/288 (92%)	0.56	20 (7%)	20 19	14, 30, 49, 67	0
1	C	269/288 (93%)	0.17	17 (6%)	26 24	11, 22, 44, 65	0
1	D	263/288 (91%)	1.27	59 (22%)	2 2	19, 34, 60, 86	0
All	All	1060/1152 (92%)	0.55	108 (10%)	12 11	11, 28, 53, 97	0

All (108) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	137	ALA	5.5
1	D	11	LEU	5.0
1	C	152	PHE	4.8
1	D	284	LEU	4.7
1	D	285	HIS	4.6
1	A	135	ASP	4.6
1	B	154	GLY	4.6
1	C	153	ILE	4.6
1	D	155	PRO	4.5
1	A	4	ALA	4.3
1	B	85	ASN	4.2
1	D	36	GLY	4.2
1	D	169	LEU	4.1
1	A	139	GLN	4.1
1	D	221	GLU	4.0
1	D	17	ILE	3.9
1	D	71	SER	3.8
1	D	12	LEU	3.7
1	A	155	PRO	3.5
1	A	154	GLY	3.5
1	D	10	ASP	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	136	ILE	3.4
1	D	137	ALA	3.3
1	D	13	ASP	3.3
1	C	141	SER	3.3
1	D	135	ASP	3.2
1	D	223	GLY	3.2
1	D	72	LYS	3.1
1	C	84	GLY	3.1
1	D	170	GLN	3.1
1	D	222	PRO	3.1
1	A	138	ARG	3.1
1	D	160	MET	3.1
1	D	154	GLY	3.0
1	D	69	ASP	3.0
1	C	285	HIS	3.0
1	C	110	SER	3.0
1	D	132	SER	3.0
1	D	134	THR	3.0
1	C	111	GLN	2.9
1	D	174	GLU	2.9
1	B	129	GLY	2.9
1	D	139	GLN	2.8
1	D	220	LEU	2.8
1	B	134	THR	2.8
1	B	190	ASP	2.8
1	D	24	GLY	2.8
1	B	155	PRO	2.8
1	C	140	PHE	2.7
1	D	53	ALA	2.7
1	D	224	MET	2.7
1	D	70	SER	2.6
1	D	217	ILE	2.6
1	C	139	GLN	2.6
1	D	67	PRO	2.6
1	B	123	ASP	2.6
1	D	138	ARG	2.6
1	C	112	GLU	2.6
1	D	166	HIS	2.6
1	B	11	LEU	2.6
1	D	269	GLU	2.5
1	D	227	ALA	2.5
1	D	26	SER	2.5

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Mol	Chain	Res	Type	RSRZ
1	C	113	GLU	2.5
1	B	140	PHE	2.5
1	D	45	TYR	2.5
1	D	164	LYS	2.5
1	D	34	VAL	2.4
1	D	54	GLU	2.4
1	B	192	LEU	2.4
1	D	14	LEU	2.4
1	C	138	ARG	2.4
1	B	167	ASN	2.4
1	B	17	ILE	2.4
1	D	157	PRO	2.4
1	B	68	GLY	2.4
1	D	216	GLY	2.3
1	D	37	GLY	2.3
1	C	142	HIS	2.3
1	D	173	VAL	2.3
1	D	97	ASN	2.3
1	B	186	GLY	2.3
1	A	131	MET	2.3
1	D	226	ILE	2.2
1	D	168	PRO	2.2
1	D	5	LEU	2.2
1	B	54	GLU	2.2
1	D	27	GLU	2.2
1	C	190	ASP	2.2
1	D	257	ALA	2.2
1	A	285	HIS	2.2
1	C	151	LYS	2.2
1	C	115	PHE	2.2
1	C	86	SER	2.2
1	D	283	ARG	2.2
1	D	190	ASP	2.2
1	D	25	GLN	2.1
1	B	4	ALA	2.1
1	B	10	ASP	2.1
1	D	282	ILE	2.1
1	D	20	GLY	2.1
1	B	141	SER	2.1
1	A	33	GLN	2.1
1	A	129	GLY	2.1
1	D	165	PHE	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	69	ASP	2.0
1	B	35	PHE	2.0
1	D	6	GLU	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.