



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 05:30 PM UTC

PDB ID : 3QMV / pdb_00003qmv
Title : RedJ-Thioesterase from the Prodiginine biosynthetic pathway in *Streptomyces coelicolor*
Authors : Whicher, J.R.; Smith, J.L.
Deposited on : 2011-02-05
Resolution : 2.12 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

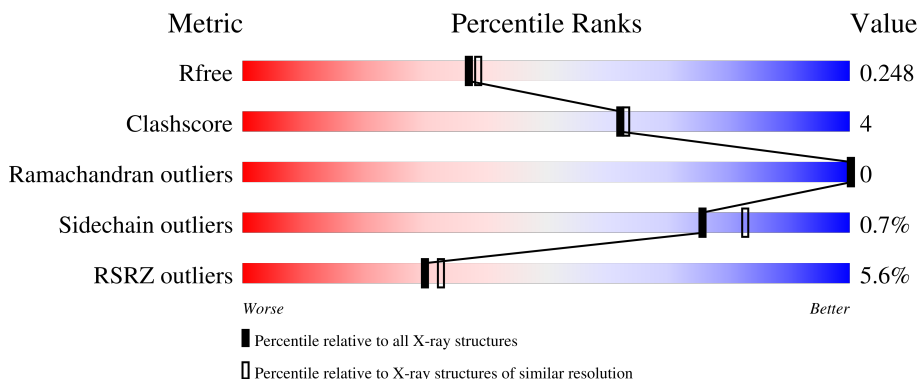
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

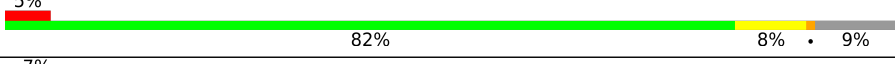
The reported resolution of this entry is 2.12 Å.

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	8290 (2.14-2.10)
Clashscore	190562	8817 (2.14-2.10)
Ramachandran outliers	187476	8738 (2.14-2.10)
Sidechain outliers	187428	8739 (2.14-2.10)
RSRZ outliers	180081	8294 (2.14-2.10)

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	280	
1	B	280	
1	C	280	
1	D	280	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8628 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 Thioesterase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	244	Total	C	N	O	S	Se	0	3	0
			1943	1225	369	342	4	3			
1	B	254	Total	C	N	O	S	Se	0	2	0
			2002	1261	378	356	4	3			
1	C	268	Total	C	N	O	S	Se	0	3	0
			2127	1340	399	381	4	3			
1	D	266	Total	C	N	O	S	Se	0	2	0
			2104	1326	393	378	4	3			

There are 96 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-18	MSE	-	expression tag	UNP O54157
A	-17	HIS	-	expression tag	UNP O54157
A	-16	HIS	-	expression tag	UNP O54157
A	-15	HIS	-	expression tag	UNP O54157
A	-14	HIS	-	expression tag	UNP O54157
A	-13	HIS	-	expression tag	UNP O54157
A	-12	HIS	-	expression tag	UNP O54157
A	-11	SER	-	expression tag	UNP O54157
A	-10	SER	-	expression tag	UNP O54157
A	-9	GLY	-	expression tag	UNP O54157
A	-8	VAL	-	expression tag	UNP O54157
A	-7	ASP	-	expression tag	UNP O54157
A	-6	LEU	-	expression tag	UNP O54157
A	-5	GLY	-	expression tag	UNP O54157
A	-4	THR	-	expression tag	UNP O54157
A	-3	GLU	-	expression tag	UNP O54157
A	-2	ASN	-	expression tag	UNP O54157
A	-1	LEU	-	expression tag	UNP O54157
A	0	TYR	-	expression tag	UNP O54157
A	1	PHE	-	expression tag	UNP O54157
A	2	GLN	-	expression tag	UNP O54157

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Chain	Residue	Modelled	Actual	Comment	Reference
A	3	SER	-	expression tag	UNP O54157
A	4	ASN	-	expression tag	UNP O54157
A	5	ALA	-	expression tag	UNP O54157
B	-18	MSE	-	expression tag	UNP O54157
B	-17	HIS	-	expression tag	UNP O54157
B	-16	HIS	-	expression tag	UNP O54157
B	-15	HIS	-	expression tag	UNP O54157
B	-14	HIS	-	expression tag	UNP O54157
B	-13	HIS	-	expression tag	UNP O54157
B	-12	HIS	-	expression tag	UNP O54157
B	-11	SER	-	expression tag	UNP O54157
B	-10	SER	-	expression tag	UNP O54157
B	-9	GLY	-	expression tag	UNP O54157
B	-8	VAL	-	expression tag	UNP O54157
B	-7	ASP	-	expression tag	UNP O54157
B	-6	LEU	-	expression tag	UNP O54157
B	-5	GLY	-	expression tag	UNP O54157
B	-4	THR	-	expression tag	UNP O54157
B	-3	GLU	-	expression tag	UNP O54157
B	-2	ASN	-	expression tag	UNP O54157
B	-1	LEU	-	expression tag	UNP O54157
B	0	TYR	-	expression tag	UNP O54157
B	1	PHE	-	expression tag	UNP O54157
B	2	GLN	-	expression tag	UNP O54157
B	3	SER	-	expression tag	UNP O54157
B	4	ASN	-	expression tag	UNP O54157
B	5	ALA	-	expression tag	UNP O54157
C	-18	MSE	-	expression tag	UNP O54157
C	-17	HIS	-	expression tag	UNP O54157
C	-16	HIS	-	expression tag	UNP O54157
C	-15	HIS	-	expression tag	UNP O54157
C	-14	HIS	-	expression tag	UNP O54157
C	-13	HIS	-	expression tag	UNP O54157
C	-12	HIS	-	expression tag	UNP O54157
C	-11	SER	-	expression tag	UNP O54157
C	-10	SER	-	expression tag	UNP O54157
C	-9	GLY	-	expression tag	UNP O54157
C	-8	VAL	-	expression tag	UNP O54157
C	-7	ASP	-	expression tag	UNP O54157
C	-6	LEU	-	expression tag	UNP O54157
C	-5	GLY	-	expression tag	UNP O54157
C	-4	THR	-	expression tag	UNP O54157

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-3	GLU	-	expression tag	UNP O54157
C	-2	ASN	-	expression tag	UNP O54157
C	-1	LEU	-	expression tag	UNP O54157
C	0	TYR	-	expression tag	UNP O54157
C	1	PHE	-	expression tag	UNP O54157
C	2	GLN	-	expression tag	UNP O54157
C	3	SER	-	expression tag	UNP O54157
C	4	ASN	-	expression tag	UNP O54157
C	5	ALA	-	expression tag	UNP O54157
D	-18	MSE	-	expression tag	UNP O54157
D	-17	HIS	-	expression tag	UNP O54157
D	-16	HIS	-	expression tag	UNP O54157
D	-15	HIS	-	expression tag	UNP O54157
D	-14	HIS	-	expression tag	UNP O54157
D	-13	HIS	-	expression tag	UNP O54157
D	-12	HIS	-	expression tag	UNP O54157
D	-11	SER	-	expression tag	UNP O54157
D	-10	SER	-	expression tag	UNP O54157
D	-9	GLY	-	expression tag	UNP O54157
D	-8	VAL	-	expression tag	UNP O54157
D	-7	ASP	-	expression tag	UNP O54157
D	-6	LEU	-	expression tag	UNP O54157
D	-5	GLY	-	expression tag	UNP O54157
D	-4	THR	-	expression tag	UNP O54157
D	-3	GLU	-	expression tag	UNP O54157
D	-2	ASN	-	expression tag	UNP O54157
D	-1	LEU	-	expression tag	UNP O54157
D	0	TYR	-	expression tag	UNP O54157
D	1	PHE	-	expression tag	UNP O54157
D	2	GLN	-	expression tag	UNP O54157
D	3	SER	-	expression tag	UNP O54157
D	4	ASN	-	expression tag	UNP O54157
D	5	ALA	-	expression tag	UNP O54157

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	134	Total O 134 134	0	0
2	B	106	Total O 106 106	0	0
2	C	118	Total O 118 118	0	0

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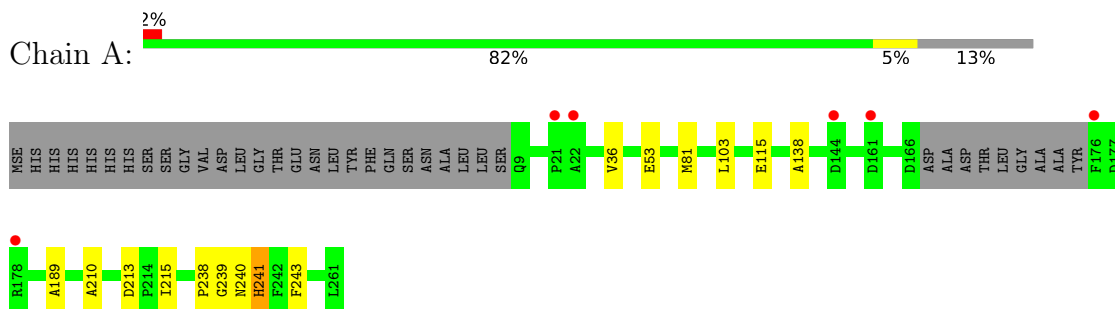
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	D	94	Total	O	0	0
			94	94		

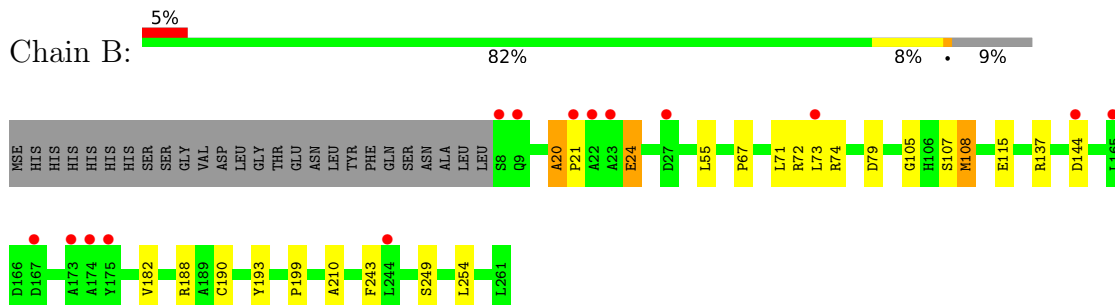
3 Residue-property plots [i](#)

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.

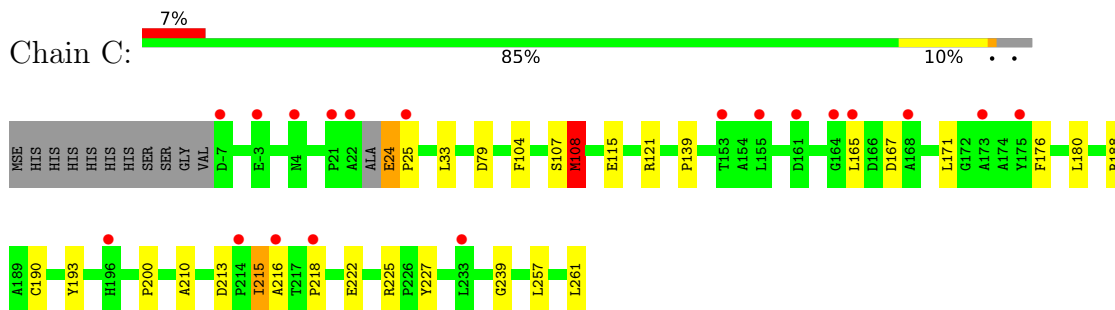
- Molecule 1: Thioesterase



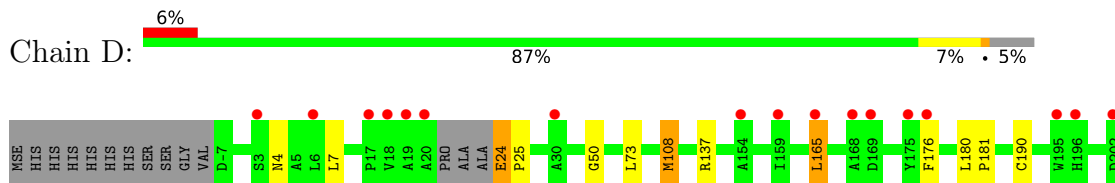
- Molecule 1: Thioesterase

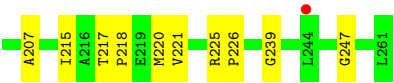


- Molecule 1: Thioesterase



- Molecule 1: Thioesterase





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	74.81Å 79.40Å 87.44Å 90.00° 90.12° 90.00°	Depositor
Resolution (Å)	50.00 – 2.12 50.00 – 2.12	Depositor EDS
% Data completeness (in resolution range)	(Not available) (50.00-2.12) 99.7 (50.00-2.12)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.01 (at 2.12Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.201 , 0.249 0.203 , 0.248	Depositor DCC
R_{free} test set	2945 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	28.4	Xtriage
Anisotropy	0.799	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 36.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.025 for h,-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	8628	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 42.98 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.8858e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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	0.82	1/1997 (0.1%)	0.82	5/2716 (0.2%)
1	B	0.75	5/2058 (0.2%)	0.80	0/2803
1	C	0.76	3/2185 (0.1%)	0.81	3/2975 (0.1%)
1	D	0.72	1/2160 (0.0%)	0.82	4/2938 (0.1%)
All	All	0.76	10/8400 (0.1%)	0.81	12/11432 (0.1%)

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	107	SER	C-O	-7.72	1.17	1.24
1	B	105	GLY	C-O	-6.37	1.17	1.24
1	B	107[A]	SER	C-O	-6.14	1.19	1.24
1	B	107[B]	SER	C-O	-6.14	1.19	1.24
1	B	20	ALA	CA-CB	-5.99	1.45	1.53
1	C	108	MSE	SE-CE	-5.64	1.78	1.95
1	A	238	PRO	C-O	-5.60	1.17	1.23
1	B	199	PRO	CA-C	5.56	1.54	1.51
1	C	215	ILE	CA-CB	-5.55	1.47	1.54
1	D	239	GLY	C-O	-5.49	1.18	1.23

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	215	ILE	N-CA-C	6.76	116.88	110.53
1	A	138	ALA	CA-C-N	5.52	124.98	119.24
1	A	138	ALA	C-N-CA	5.52	124.98	119.24
1	A	239	GLY	N-CA-C	5.39	117.54	112.04
1	C	24	GLU	CA-C-N	5.31	139.74	127.00
1	C	24	GLU	C-N-CA	5.31	139.74	127.00
1	C	24	GLU	C-N-CD	-5.26	109.03	120.60
1	A	241	HIS	N-CA-C	5.18	118.38	111.75
1	D	24	GLU	CA-C-N	5.17	139.41	127.00
1	D	24	GLU	C-N-CA	5.17	139.41	127.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	247	GLY	CA-C-N	5.05	125.08	119.32
1	D	247	GLY	C-N-CA	5.05	125.08	119.32

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	1943	0	1892	9	0
1	B	2002	0	1947	17	0
1	C	2127	0	2068	23	0
1	D	2104	0	2043	18	0
2	A	134	0	0	1	0
2	B	106	0	0	1	0
2	C	118	0	0	0	0
2	D	94	0	0	0	0
All	All	8628	0	7950	66	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 (66) 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:24:GLU:HB3	1:C:25:PRO:C	2.17	0.68
1:A:213:ASP:CG	1:A:241:HIS:HD1	2.01	0.68
1:D:24:GLU:HB3	1:D:25:PRO:HA	1.76	0.67
1:A:53[B]:GLU:H	1:A:53[B]:GLU:CD	2.01	0.67
1:C:24:GLU:HB3	1:C:25:PRO:CA	2.26	0.65
1:C:215:ILE:HG22	1:C:215:ILE:O	1.97	0.63
1:B:55:LEU:CD1	1:B:254:LEU:HD22	2.31	0.60
1:D:24:GLU:HG3	1:D:25:PRO:C	2.29	0.58
1:A:53[B]:GLU:CD	1:A:53[B]:GLU:N	2.62	0.57
1:D:108:MSE:HG3	1:D:190:CYS:SG	2.47	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:210:ALA:HB2	1:B:243:PHE:CD2	2.42	0.54
1:B:24:GLU:OE2	1:B:24:GLU:HA	2.08	0.54
1:D:137:ARG:HG2	1:D:220:MSE:HB3	1.90	0.54
1:B:55:LEU:HD11	1:B:254:LEU:HD22	1.91	0.53
1:C:222:GLU:OE2	1:C:225[B]:ARG:NH1	2.38	0.53
1:C:79:ASP:O	1:C:188:ARG:NH1	2.41	0.52
1:B:73:LEU:HD12	1:B:74:ARG:HD3	1.91	0.52
1:D:176:PHE:CE2	1:D:180:LEU:HD21	2.45	0.52
1:C:33:LEU:HD21	1:C:261:LEU:HD22	1.93	0.51
1:A:210:ALA:HB2	1:A:243:PHE:CD2	2.45	0.51
1:C:115:GLU:OE2	1:C:193:TYR:OH	2.27	0.51
1:B:21:PRO:O	1:B:21:PRO:HG2	2.12	0.49
1:A:81:MSE:HG3	1:A:115:GLU:HG3	1.95	0.48
1:C:213:ASP:HB3	1:C:216:ALA:HB3	1.96	0.47
1:C:165:LEU:HD11	1:C:176:PHE:HB2	1.96	0.47
1:D:4:ASN:HB3	1:D:7:LEU:HG	1.96	0.47
1:B:55:LEU:HD11	1:B:254:LEU:CD2	2.44	0.47
1:B:108:MSE:HG3	1:B:190:CYS:SG	2.54	0.47
1:D:165:LEU:HD11	1:D:176:PHE:CG	2.50	0.46
1:D:137:ARG:NH1	1:D:215:ILE:HG23	2.30	0.46
1:D:24:GLU:HB3	1:D:25:PRO:CA	2.44	0.46
1:D:165:LEU:CD1	1:D:176:PHE:CD1	2.98	0.46
1:A:81:MSE:HE2	1:A:189:ALA:HB1	1.97	0.46
1:C:167:ASP:HB3	1:C:171:LEU:HD12	1.98	0.46
1:C:24:GLU:HB3	1:C:25:PRO:HA	1.97	0.46
1:C:176:PHE:CE2	1:C:180:LEU:HD21	2.52	0.45
1:C:33:LEU:CD2	1:C:261:LEU:HD22	2.46	0.45
1:B:243:PHE:HB2	1:B:249:SER:HB2	1.98	0.45
1:B:67:PRO:HG3	1:B:72:ARG:NH2	2.32	0.45
1:A:240:ASN:HB2	2:A:358:HOH:O	2.17	0.45
1:A:210:ALA:HB3	1:A:213:ASP:HB2	1.98	0.45
1:C:108:MSE:O	1:C:108:MSE:HG2	2.14	0.44
1:D:137:ARG:HH12	1:D:215:ILE:HG23	1.81	0.44
1:B:20:ALA:HB1	1:B:21:PRO:HD2	2.00	0.44
1:B:115:GLU:OE2	1:B:193:TYR:OH	2.23	0.44
1:D:73:LEU:C	1:D:73:LEU:HD12	2.42	0.44
1:C:139:PRO:HD2	1:C:227:TYR:CZ	2.53	0.44
1:C:210:ALA:HB3	1:C:213:ASP:HB2	1.98	0.44
1:B:137:ARG:NH1	1:B:144:ASP:OD1	2.51	0.43
1:D:225:ARG:HB3	1:D:226:PRO:HD3	2.01	0.43
1:D:165:LEU:HD11	1:D:176:PHE:CD1	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:210:ALA:HB1	1:C:239:GLY:O	2.18	0.42
1:C:121:ARG:NH1	1:C:200:PRO:O	2.53	0.42
1:B:79:ASP:HB2	1:B:188:ARG:HD2	2.01	0.42
1:C:104:PHE:CD2	1:C:104:PHE:C	2.98	0.42
1:D:207:ALA:HB1	1:D:221:VAL:CG1	2.50	0.42
1:D:180:LEU:HB2	1:D:181:PRO:HD3	2.01	0.42
1:C:180:LEU:HD23	1:C:180:LEU:HA	1.94	0.41
1:B:55:LEU:HD13	1:B:254:LEU:HD22	2.01	0.41
1:D:217:THR:HB	1:D:218:PRO:HD2	2.02	0.41
1:B:20:ALA:HB2	1:D:50:GLY:HA2	2.01	0.41
1:B:182:VAL:HG23	2:B:300:HOH:O	2.19	0.41
1:C:218:PRO:O	1:C:222:GLU:HB2	2.21	0.41
1:C:108:MSE:HB2	1:C:190:CYS:SG	2.60	0.41
1:A:36:VAL:HB	1:A:103:LEU:HD23	2.03	0.41
1:C:257:LEU:O	1:C:261:LEU:HG	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	243/280 (87%)	240 (99%)	3 (1%)	0	100	100
1	B	254/280 (91%)	250 (98%)	4 (2%)	0	100	100
1	C	267/280 (95%)	263 (98%)	4 (2%)	0	100	100
1	D	264/280 (94%)	259 (98%)	5 (2%)	0	100	100
All	All	1028/1120 (92%)	1012 (98%)	16 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	197/219 (90%)	197 (100%)	0	100	100
1	B	202/219 (92%)	199 (98%)	3 (2%)	57	65
1	C	216/219 (99%)	215 (100%)	1 (0%)	81	87
1	D	214/219 (98%)	212 (99%)	2 (1%)	70	78
All	All	829/876 (95%)	823 (99%)	6 (1%)	76	83

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	24	GLU
1	B	71	LEU
1	B	108	MSE
1	C	108	MSE
1	D	108	MSE
1	D	165	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	95	HIS
1	B	196	HIS

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 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	241/280 (86%)	0.06	6 (2%) 58 61	8, 25, 47, 55	3 (1%)
1	B	251/280 (89%)	0.30	14 (5%) 30 33	8, 29, 50, 58	2 (0%)
1	C	265/280 (94%)	0.40	19 (7%) 21 23	11, 32, 53, 58	3 (1%)
1	D	263/280 (93%)	0.62	18 (6%) 23 25	10, 34, 57, 64	2 (0%)
All	All	1020/1120 (91%)	0.35	57 (5%) 30 33	8, 30, 52, 64	10 (0%)

All (57) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	22	ALA	4.2
1	A	21	PRO	3.9
1	C	175	TYR	3.8
1	D	20	ALA	3.8
1	D	176	PHE	3.6
1	B	22	ALA	3.6
1	A	22	ALA	3.5
1	D	165	LEU	3.4
1	D	159	ILE	3.1
1	D	168	ALA	3.0
1	B	73	LEU	3.0
1	C	165	LEU	3.0
1	C	21	PRO	2.8
1	B	9	GLN	2.8
1	B	144	ASP	2.8
1	B	167	ASP	2.8
1	D	19	ALA	2.8
1	B	8	SER	2.7
1	D	202	ASP	2.7
1	A	176	PHE	2.7
1	D	175	TYR	2.6

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Mol	Chain	Res	Type	RSRZ
1	C	196	HIS	2.6
1	C	216	ALA	2.6
1	A	144	ASP	2.6
1	C	-3	GLU	2.5
1	D	196	HIS	2.5
1	C	25	PRO	2.5
1	C	233	LEU	2.5
1	D	169	ASP	2.5
1	B	174	ALA	2.4
1	D	6	LEU	2.4
1	A	178	ARG	2.4
1	C	4	ASN	2.4
1	C	161	ASP	2.4
1	D	18	VAL	2.4
1	D	17	PRO	2.4
1	B	165	LEU	2.3
1	B	244	LEU	2.3
1	B	27	ASP	2.3
1	C	155	LEU	2.3
1	D	195	TRP	2.3
1	B	21	PRO	2.3
1	A	161	ASP	2.3
1	B	23	ALA	2.2
1	D	30	ALA	2.2
1	D	154	ALA	2.2
1	C	214	PRO	2.1
1	C	218	PRO	2.1
1	C	168	ALA	2.1
1	B	175	TYR	2.1
1	C	153	THR	2.1
1	B	173	ALA	2.1
1	C	173	ALA	2.1
1	D	3	SER	2.1
1	C	164	GLY	2.0
1	D	244	LEU	2.0
1	C	-7	ASP	2.0

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

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.