



Full wwPDB X-ray Structure Validation Report ⓘ

May 16, 2020 – 06:47 am BST

PDB ID : 3GVC
Title : Crystal structure of probable short-chain dehydrogenase-reductase from Mycobacterium tuberculosis
Authors : Seattle Structural Genomics Center for Infectious Disease (SSGCID)
Deposited on : 2009-03-30
Resolution : 2.45 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	:	2.11
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)	:	2.11

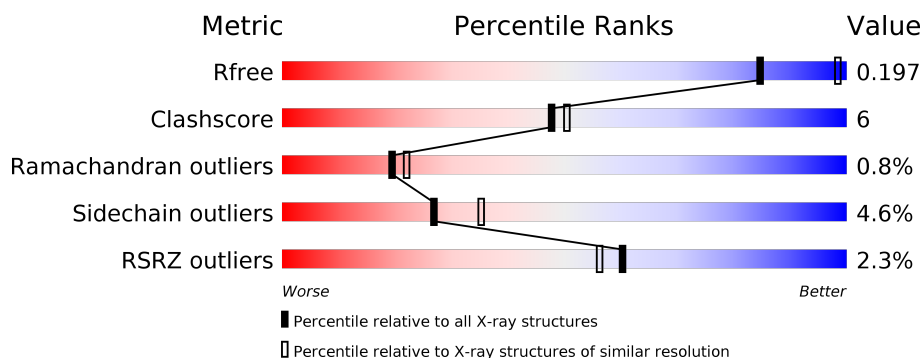
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.45 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	1544 (2.48-2.44)
Clashscore	141614	1613 (2.48-2.44)
Ramachandran outliers	138981	1598 (2.48-2.44)
Sidechain outliers	138945	1598 (2.48-2.44)
RSRZ outliers	127900	1523 (2.48-2.44)

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 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	277	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 11%, green 77%, grey 12%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 77% 11% • 12% </div> </div>
1	B	277	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 3%, orange 1%, yellow 11%, green 81%, grey 8%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 3% 81% 11% • 8% </div> </div>
1	C	277	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 3%, orange 1%, yellow 11%, green 78%, grey 10%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 3% 78% 11% • 10% </div> </div>
1	D	277	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 16%, green 73%, grey 10%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 73% 16% • 10% </div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7223 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 Probable short-chain type dehydrogenase/reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	245	Total	C	N	O	S	0	0	0
			1707	1056	305	330	16			
1	B	255	Total	C	N	O	S	0	0	0
			1794	1110	322	345	17			
1	C	250	Total	C	N	O	S	0	0	0
			1743	1080	310	337	16			
1	D	249	Total	C	N	O	S	0	0	0
			1739	1074	311	338	16			

There are 84 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	EXPRESSION TAG	UNP P95273
A	2	ALA	-	EXPRESSION TAG	UNP P95273
A	3	HIS	-	EXPRESSION TAG	UNP P95273
A	4	HIS	-	EXPRESSION TAG	UNP P95273
A	5	HIS	-	EXPRESSION TAG	UNP P95273
A	6	HIS	-	EXPRESSION TAG	UNP P95273
A	7	HIS	-	EXPRESSION TAG	UNP P95273
A	8	HIS	-	EXPRESSION TAG	UNP P95273
A	9	MET	-	EXPRESSION TAG	UNP P95273
A	10	GLY	-	EXPRESSION TAG	UNP P95273
A	11	THR	-	EXPRESSION TAG	UNP P95273
A	12	LEU	-	EXPRESSION TAG	UNP P95273
A	13	GLU	-	EXPRESSION TAG	UNP P95273
A	14	ALA	-	EXPRESSION TAG	UNP P95273
A	15	GLN	-	EXPRESSION TAG	UNP P95273
A	16	THR	-	EXPRESSION TAG	UNP P95273
A	17	GLN	-	EXPRESSION TAG	UNP P95273
A	18	GLY	-	EXPRESSION TAG	UNP P95273
A	19	PRO	-	EXPRESSION TAG	UNP P95273
A	20	GLY	-	EXPRESSION TAG	UNP P95273
A	21	SER	-	EXPRESSION TAG	UNP P95273

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Chain	Residue	Modelled	Actual	Comment	Reference
B	1	MET	-	EXPRESSION TAG	UNP P95273
B	2	ALA	-	EXPRESSION TAG	UNP P95273
B	3	HIS	-	EXPRESSION TAG	UNP P95273
B	4	HIS	-	EXPRESSION TAG	UNP P95273
B	5	HIS	-	EXPRESSION TAG	UNP P95273
B	6	HIS	-	EXPRESSION TAG	UNP P95273
B	7	HIS	-	EXPRESSION TAG	UNP P95273
B	8	HIS	-	EXPRESSION TAG	UNP P95273
B	9	MET	-	EXPRESSION TAG	UNP P95273
B	10	GLY	-	EXPRESSION TAG	UNP P95273
B	11	THR	-	EXPRESSION TAG	UNP P95273
B	12	LEU	-	EXPRESSION TAG	UNP P95273
B	13	GLU	-	EXPRESSION TAG	UNP P95273
B	14	ALA	-	EXPRESSION TAG	UNP P95273
B	15	GLN	-	EXPRESSION TAG	UNP P95273
B	16	THR	-	EXPRESSION TAG	UNP P95273
B	17	GLN	-	EXPRESSION TAG	UNP P95273
B	18	GLY	-	EXPRESSION TAG	UNP P95273
B	19	PRO	-	EXPRESSION TAG	UNP P95273
B	20	GLY	-	EXPRESSION TAG	UNP P95273
B	21	SER	-	EXPRESSION TAG	UNP P95273
C	1	MET	-	EXPRESSION TAG	UNP P95273
C	2	ALA	-	EXPRESSION TAG	UNP P95273
C	3	HIS	-	EXPRESSION TAG	UNP P95273
C	4	HIS	-	EXPRESSION TAG	UNP P95273
C	5	HIS	-	EXPRESSION TAG	UNP P95273
C	6	HIS	-	EXPRESSION TAG	UNP P95273
C	7	HIS	-	EXPRESSION TAG	UNP P95273
C	8	HIS	-	EXPRESSION TAG	UNP P95273
C	9	MET	-	EXPRESSION TAG	UNP P95273
C	10	GLY	-	EXPRESSION TAG	UNP P95273
C	11	THR	-	EXPRESSION TAG	UNP P95273
C	12	LEU	-	EXPRESSION TAG	UNP P95273
C	13	GLU	-	EXPRESSION TAG	UNP P95273
C	14	ALA	-	EXPRESSION TAG	UNP P95273
C	15	GLN	-	EXPRESSION TAG	UNP P95273
C	16	THR	-	EXPRESSION TAG	UNP P95273
C	17	GLN	-	EXPRESSION TAG	UNP P95273
C	18	GLY	-	EXPRESSION TAG	UNP P95273
C	19	PRO	-	EXPRESSION TAG	UNP P95273
C	20	GLY	-	EXPRESSION TAG	UNP P95273
C	21	SER	-	EXPRESSION TAG	UNP P95273

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Chain	Residue	Modelled	Actual	Comment	Reference
D	1	MET	-	EXPRESSION TAG	UNP P95273
D	2	ALA	-	EXPRESSION TAG	UNP P95273
D	3	HIS	-	EXPRESSION TAG	UNP P95273
D	4	HIS	-	EXPRESSION TAG	UNP P95273
D	5	HIS	-	EXPRESSION TAG	UNP P95273
D	6	HIS	-	EXPRESSION TAG	UNP P95273
D	7	HIS	-	EXPRESSION TAG	UNP P95273
D	8	HIS	-	EXPRESSION TAG	UNP P95273
D	9	MET	-	EXPRESSION TAG	UNP P95273
D	10	GLY	-	EXPRESSION TAG	UNP P95273
D	11	THR	-	EXPRESSION TAG	UNP P95273
D	12	LEU	-	EXPRESSION TAG	UNP P95273
D	13	GLU	-	EXPRESSION TAG	UNP P95273
D	14	ALA	-	EXPRESSION TAG	UNP P95273
D	15	GLN	-	EXPRESSION TAG	UNP P95273
D	16	THR	-	EXPRESSION TAG	UNP P95273
D	17	GLN	-	EXPRESSION TAG	UNP P95273
D	18	GLY	-	EXPRESSION TAG	UNP P95273
D	19	PRO	-	EXPRESSION TAG	UNP P95273
D	20	GLY	-	EXPRESSION TAG	UNP P95273
D	21	SER	-	EXPRESSION TAG	UNP P95273

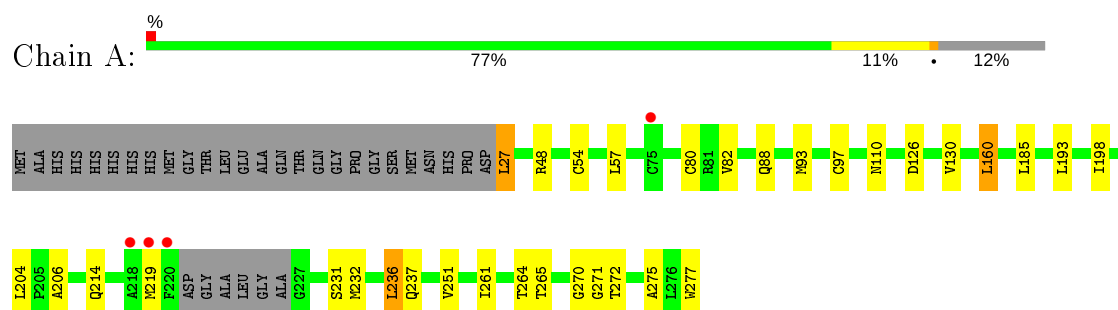
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	64	Total O 64 64	0	0
2	B	70	Total O 70 70	0	0
2	C	61	Total O 61 61	0	0
2	D	45	Total O 45 45	0	0

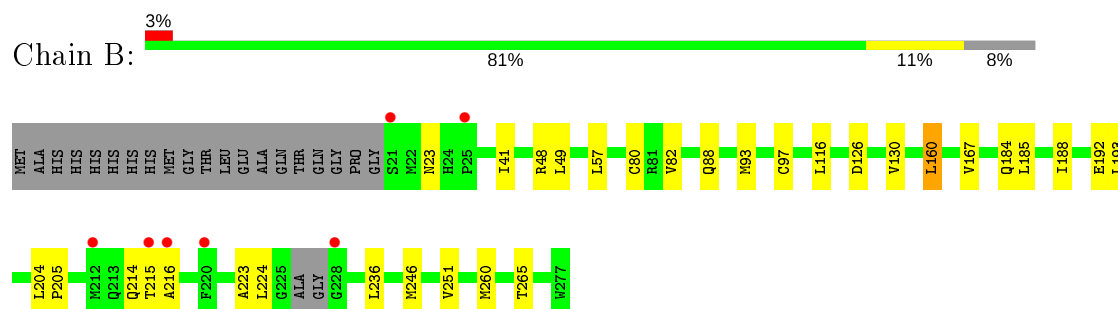
3 Residue-property plots

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.

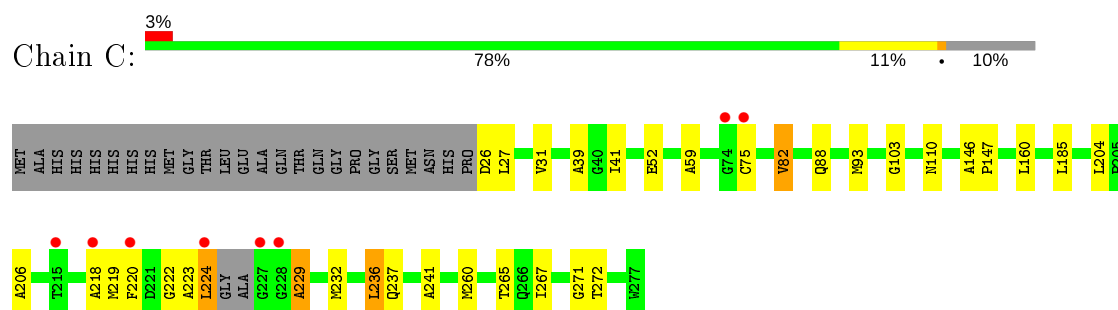
- Molecule 1: Probable short-chain type dehydrogenase/reductase



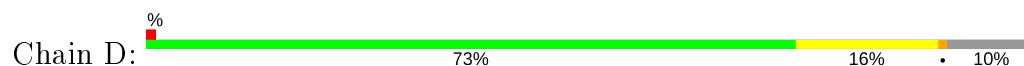
- Molecule 1: Probable short-chain type dehydrogenase/reductase

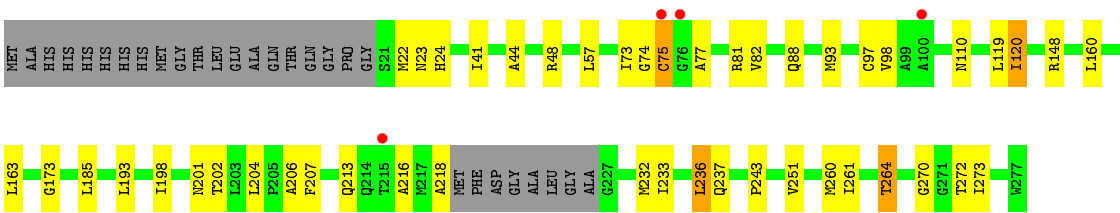


- Molecule 1: Probable short-chain type dehydrogenase/reductase



- Molecule 1: Probable short-chain type dehydrogenase/reductase





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	81.69 Å 51.38 Å 118.84 Å 90.00° 94.50° 90.00°	Depositor
Resolution (Å)	19.90 – 2.45 19.91 – 2.45	Depositor EDS
% Data completeness (in resolution range)	98.8 (19.90-2.45) 99.0 (19.91-2.45)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.42 (at 2.44 Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.192 , 0.247 0.198 , 0.197	Depositor DCC
R_{free} test set	1802 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	26.5	Xtriage
Anisotropy	0.124	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 34.5	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	0.93	EDS
Total number of atoms	7223	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.11% 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	0.54	1/1721 (0.1%)	0.62	0/2330
1	B	0.52	0/1811	0.63	0/2450
1	C	0.48	0/1758	0.61	0/2380
1	D	0.48	0/1755	0.61	0/2378
All	All	0.51	1/7045 (0.0%)	0.62	0/9538

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	54	CYS	CB-SG	-5.45	1.73	1.81

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	1707	0	1722	21	0
1	B	1794	0	1815	15	0
1	C	1743	0	1754	24	0
1	D	1739	0	1748	31	0
2	A	64	0	0	1	0
2	B	70	0	0	0	0
2	C	61	0	0	1	0
2	D	45	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	7223	0	7039	86	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 6.

All (86) 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:57:LEU:HD22	1:D:97:CYS:HB2	1.73	0.71
1:C:26:ASP:HB2	1:C:52:GLU:OE1	1.93	0.69
1:C:204:LEU:HD13	1:C:265:THR:HG23	1.74	0.68
1:D:74:GLY:HA3	1:D:75:CYS:HB2	1.77	0.67
1:A:82:VAL:HG21	1:A:93:MET:SD	2.35	0.66
1:A:206:ALA:HB3	1:A:237:GLN:OE1	1.95	0.66
1:D:232:MET:HG2	1:D:236:LEU:HD22	1.78	0.66
1:A:110:ASN:ND2	1:A:160:LEU:HD22	2.11	0.65
1:D:41:ILE:HG21	1:D:160:LEU:HD21	1.78	0.65
1:A:204:LEU:HD13	1:A:265:THR:HG23	1.78	0.64
1:D:193:LEU:HD13	1:D:198:ILE:HD12	1.79	0.64
1:D:98:VAL:HG21	1:D:148:ARG:NH2	2.14	0.63
1:A:88:GLN:NE2	2:A:295:HOH:O	2.28	0.62
1:A:57:LEU:HD11	1:A:80:CYS:HB2	1.81	0.62
1:A:27:LEU:N	1:A:27:LEU:HD12	2.16	0.61
1:D:236:LEU:HG	1:D:272:THR:HG23	1.83	0.60
1:B:116:LEU:HD11	1:B:216:ALA:HA	1.84	0.59
1:D:218:ALA:HB1	2:D:298:HOH:O	2.03	0.59
1:A:48:ARG:HG3	1:A:251:VAL:HG21	1.83	0.58
1:A:204:LEU:CD1	1:A:265:THR:HG23	2.32	0.58
1:C:220:PHE:CE2	1:C:229:ALA:HB3	2.40	0.56
1:D:48:ARG:HG3	1:D:251:VAL:HG21	1.88	0.56
1:A:57:LEU:HD22	1:A:97:CYS:HB2	1.87	0.56
1:C:31:VAL:HG11	1:C:103:GLY:O	2.07	0.55
1:B:41:ILE:HG21	1:B:160:LEU:HD21	1.88	0.55
1:D:73:ILE:HG22	1:D:75:CYS:CB	2.37	0.55
1:D:110:ASN:OD1	1:D:160:LEU:HD22	2.07	0.55
1:C:110:ASN:OD1	1:C:160:LEU:HD22	2.08	0.54
1:C:241:ALA:HB2	1:D:260:MET:HE2	1.90	0.54
1:C:204:LEU:CD1	1:C:265:THR:HG23	2.37	0.53
1:B:224:LEU:N	1:B:224:LEU:HD22	2.23	0.53
1:A:110:ASN:HD22	1:A:160:LEU:HD22	1.71	0.53
1:C:222:GLY:O	1:C:223:ALA:HB3	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:204:LEU:HD13	1:B:265:THR:HG23	1.91	0.53
1:C:41:ILE:HG21	1:C:160:LEU:HD21	1.92	0.51
1:B:48:ARG:HG3	1:B:251:VAL:HG21	1.92	0.51
1:D:74:GLY:CA	1:D:75:CYS:HB2	2.41	0.51
1:A:204:LEU:HD13	1:A:265:THR:CG2	2.42	0.49
1:A:27:LEU:N	1:A:27:LEU:CD1	2.75	0.49
1:B:82:VAL:HG21	1:B:93:MET:SD	2.53	0.49
1:C:260:MET:HE1	1:D:270:GLY:HA3	1.95	0.48
1:C:146:ALA:HB3	1:C:147:PRO:HD3	1.95	0.48
1:D:206:ALA:HB3	1:D:237:GLN:OE1	2.13	0.48
1:C:59:ALA:HB1	1:C:82:VAL:CG2	2.44	0.48
1:A:236:LEU:HG	1:A:272:THR:HG23	1.95	0.48
1:D:163:LEU:HD11	1:D:273:ILE:CG2	2.43	0.47
1:D:120:ILE:HG23	2:D:296:HOH:O	2.14	0.47
1:A:275:ALA:HB3	1:A:277:TRP:CE2	2.50	0.47
1:A:270:GLY:N	1:B:260:MET:HE3	2.30	0.47
1:C:39:ALA:HB1	2:C:306:HOH:O	2.13	0.47
1:C:204:LEU:HD22	1:C:267:ILE:HD13	1.96	0.47
1:D:202:THR:HG22	1:D:204:LEU:HD12	1.96	0.47
1:B:126:ASP:O	1:B:130:VAL:HG23	2.15	0.46
1:D:75:CYS:O	1:D:75:CYS:SG	2.74	0.46
1:B:204:LEU:HD22	1:B:265:THR:CG2	2.46	0.46
1:B:192:GLU:O	1:B:193:LEU:HD23	2.15	0.46
1:D:73:ILE:HG22	1:D:75:CYS:HB2	1.96	0.46
1:C:204:LEU:HD22	1:C:267:ILE:CD1	2.46	0.45
1:C:59:ALA:HB1	1:C:82:VAL:HG23	1.97	0.45
1:B:57:LEU:HD22	1:B:97:CYS:HB2	1.98	0.45
1:C:26:ASP:OD1	1:C:52:GLU:OE1	2.35	0.45
1:D:73:ILE:HG22	1:D:75:CYS:HB3	1.99	0.45
1:B:49:LEU:HD23	1:B:251:VAL:HG22	1.99	0.44
1:A:271:GLY:O	1:A:272:THR:C	2.56	0.44
1:B:205:PRO:HB3	1:B:246:MET:SD	2.58	0.44
1:C:224:LEU:HD12	1:C:224:LEU:N	2.32	0.44
1:D:213:GLN:O	1:D:216:ALA:HB3	2.19	0.42
1:D:22:MET:CE	1:D:24:HIS:CE1	3.02	0.42
1:D:261:ILE:HG23	1:D:264:THR:HG21	2.00	0.42
1:B:184:GLN:O	1:B:188:ILE:HG22	2.18	0.42
1:C:271:GLY:O	1:C:272:THR:C	2.58	0.42
1:D:201:ASN:OD1	1:D:264:THR:HG23	2.20	0.42
1:D:44:ALA:HB2	1:D:243:PRO:HB2	2.02	0.42
1:C:82:VAL:HG21	1:C:93:MET:SD	2.60	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:202:THR:HG22	1:D:204:LEU:CD1	2.50	0.41
1:D:82:VAL:HG21	1:D:93:MET:SD	2.59	0.41
1:A:193:LEU:HB3	1:A:198:ILE:HB	2.02	0.41
1:D:207:PHE:CG	1:D:233:ILE:HG12	2.54	0.41
1:A:126:ASP:O	1:A:130:VAL:HG23	2.21	0.41
1:A:261:ILE:HG23	1:A:264:THR:HG21	2.02	0.41
1:C:206:ALA:HB3	1:C:237:GLN:OE1	2.20	0.41
1:A:236:LEU:HB3	1:A:272:THR:HG23	2.02	0.41
1:C:260:MET:CE	1:D:270:GLY:CA	2.99	0.41
1:C:260:MET:HE1	1:D:270:GLY:CA	2.51	0.40
1:B:41:ILE:HG21	1:B:160:LEU:CD2	2.50	0.40
1:C:232:MET:HG2	1:C:236:LEU:HD22	2.04	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	241/277 (87%)	230 (95%)	11 (5%)	0	100	100
1	B	251/277 (91%)	239 (95%)	11 (4%)	1 (0%)	34	41
1	C	246/277 (89%)	229 (93%)	13 (5%)	4 (2%)	9	8
1	D	245/277 (88%)	233 (95%)	9 (4%)	3 (1%)	13	12
All	All	983/1108 (89%)	931 (95%)	44 (4%)	8 (1%)	19	22

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	229	ALA
1	D	75	CYS
1	C	219	MET

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Mol	Chain	Res	Type
1	C	75	CYS
1	C	218	ALA
1	B	223	ALA
1	D	77	ALA
1	D	173	GLY

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	165/192 (86%)	157 (95%)	8 (5%)	25	33
1	B	176/192 (92%)	167 (95%)	9 (5%)	24	31
1	C	168/192 (88%)	162 (96%)	6 (4%)	35	46
1	D	170/192 (88%)	162 (95%)	8 (5%)	26	34
All	All	679/768 (88%)	648 (95%)	31 (5%)	27	35

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	27	LEU
1	A	160	LEU
1	A	185	LEU
1	A	214	GLN
1	A	219	MET
1	A	231	SER
1	A	232	MET
1	A	236	LEU
1	B	23	ASN
1	B	80	CYS
1	B	88	GLN
1	B	160	LEU
1	B	167	VAL
1	B	185	LEU
1	B	214	GLN
1	B	215	THR

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Mol	Chain	Res	Type
1	B	236	LEU
1	C	27	LEU
1	C	82	VAL
1	C	88	GLN
1	C	185	LEU
1	C	224	LEU
1	C	236	LEU
1	D	23	ASN
1	D	81	ARG
1	D	88	GLN
1	D	119	LEU
1	D	120	ILE
1	D	185	LEU
1	D	236	LEU
1	D	264	THR

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

Mol	Chain	Res	Type
1	A	110	ASN
1	C	214	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

There are no such residues in this entry.

5.8 Polymer linkage issues

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	245/277 (88%)	-0.42	4 (1%) 72 69	13, 20, 37, 62	0
1	B	255/277 (92%)	-0.33	7 (2%) 54 50	12, 20, 54, 65	0
1	C	250/277 (90%)	-0.35	8 (3%) 47 44	13, 22, 55, 73	0
1	D	249/277 (89%)	-0.18	4 (1%) 72 69	17, 27, 51, 59	0
All	All	999/1108 (90%)	-0.32	23 (2%) 60 56	12, 22, 52, 73	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	218	ALA	5.4
1	C	218	ALA	4.9
1	D	75	CYS	4.6
1	C	220	PHE	3.8
1	C	224	LEU	3.5
1	C	227	GLY	3.4
1	B	220	PHE	3.4
1	D	215	THR	3.3
1	B	25	PRO	3.3
1	C	75	CYS	3.2
1	C	74	GLY	3.1
1	C	228	GLY	3.0
1	D	76	GLY	2.9
1	B	215	THR	2.9
1	A	220	PHE	2.8
1	B	228	GLY	2.6
1	C	215	THR	2.5
1	A	219	MET	2.5
1	A	75	CYS	2.4
1	B	216	ALA	2.4
1	B	21	SER	2.2

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Mol	Chain	Res	Type	RSRZ
1	D	100	ALA	2.2
1	B	212	MET	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.