



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 07:06 PM UTC

PDB ID : 1U5H / pdb_00001u5h
Title : Structure of Citrate Lyase beta subunit from Mycobacterium tuberculosis
Authors : Goulding, C.W.; Legin, T.; Kim, C.Y.; Segelke, B.; Terwilliger, T.C.; Eisenberg, D.; TB Structural Genomics Consortium (TBSGC)
Deposited on : 2004-07-27
Resolution : 1.65 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

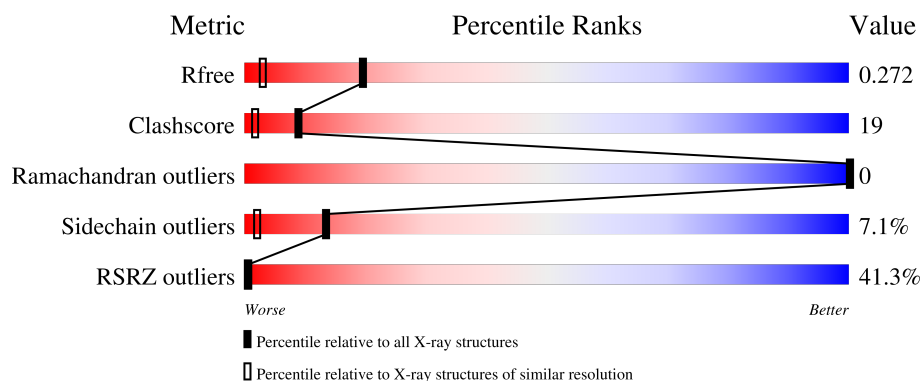
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.65 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	2563 (1.66-1.66)
Clashscore	190562	2662 (1.66-1.66)
Ramachandran outliers	187476	2621 (1.66-1.66)
Sidechain outliers	187428	2621 (1.66-1.66)
RSRZ outliers	180081	2564 (1.66-1.66)

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	273	34% 63% 15% • • 18%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 1776 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.

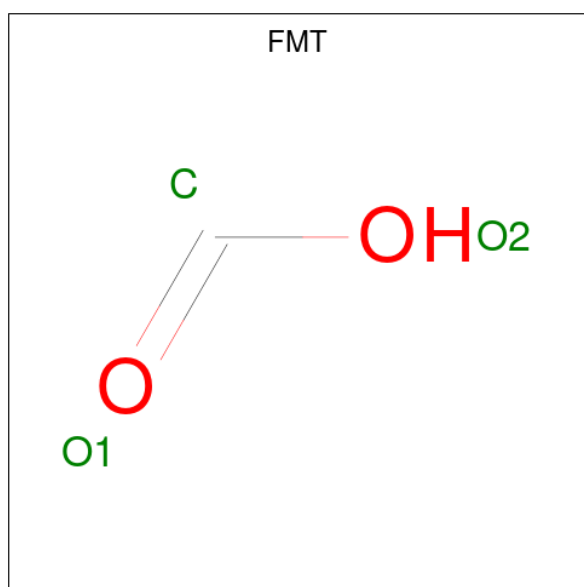
- Molecule 1 is a protein called citE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	223	1617	1006	290	314	7	0	0	0

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	17	ALA	ARG	engineered mutation	UNP O06162
A	20	ALA	ARG	engineered mutation	UNP O06162
A	23	ALA	LYS	engineered mutation	UNP O06162
A	222	ALA	ARG	engineered mutation	UNP O06162
A	223	ALA	PRO	engineered mutation	UNP O06162

- Molecule 2 is FORMIC ACID (CCD ID: FMT) (formula: CH₂O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
2	A	1	3	1	2	0	0

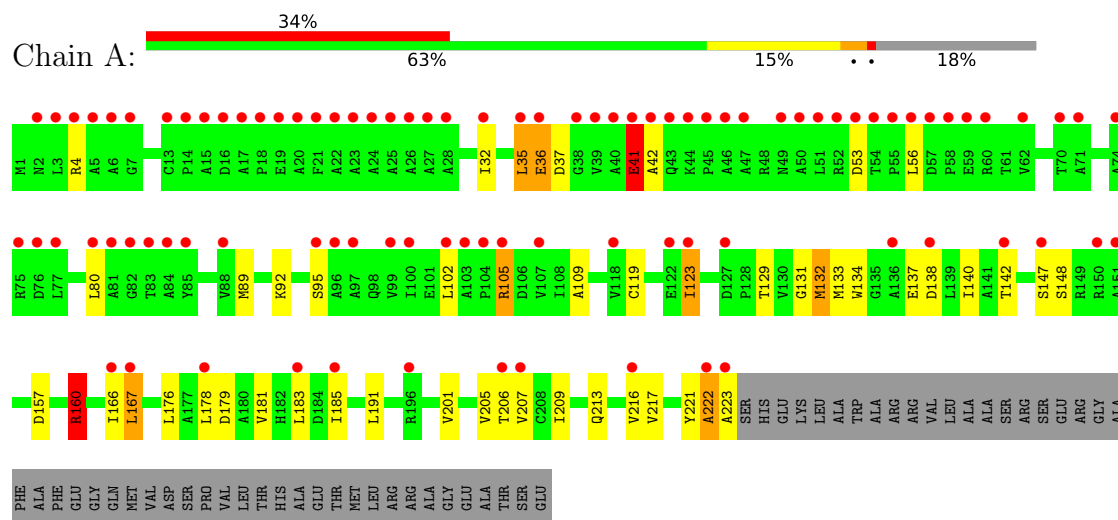
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	156	Total 156	O 156	0	0

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: citE



4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	91.21Å 91.21Å 221.23Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	19.67 – 1.65 19.67 – 1.65	Depositor EDS
% Data completeness (in resolution range)	96.8 (19.67-1.65) 96.7 (19.67-1.65)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.62 (at 1.50Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.268 , 0.272 0.266 , 0.272	Depositor DCC
R_{free} test set	4213 reflections (7.55%)	wwPDB-VP
Wilson B-factor (Å ²)	22.2	Xtriage
Anisotropy	0.266	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 37.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.014 for -h,1/3*h-1/3*k-1/3*l,-4/3*h-8/3*k +1/3*l 0.016 for -1/3*h+1/3*k+1/3*l,-k,8/3*h+4/ 3*k+1/3*l 0.020 for -2/3*h-1/3*k-1/3*l,-1/3*h-2/3*k+ 1/3*l,-4/3*h+4/3*k+1/3*l	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	1776	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.66% 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

Bond lengths and bond angles in the following residue types are not validated in this section: FMT

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	1.20	3/1641 (0.2%)	1.36	8/2241 (0.4%)

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	A	0	1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	41	GLU	CG-CD	24.87	2.14	1.52
1	A	132	MET	SD-CE	-13.84	1.45	1.79
1	A	160	ARG	CD-NE	-7.09	1.36	1.46

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	41	GLU	CB-CG-CD	33.95	170.31	112.60
1	A	41	GLU	CG-CD-OE1	-19.68	73.13	118.40
1	A	41	GLU	CG-CD-OE2	17.66	159.02	118.40
1	A	160	ARG	NE-CZ-NH2	-7.32	112.61	119.20
1	A	160	ARG	CD-NE-CZ	6.65	133.71	124.40
1	A	160	ARG	CG-CD-NE	-6.29	98.17	112.00
1	A	160	ARG	NE-CZ-NH1	6.26	127.76	121.50
1	A	222	ALA	N-CA-C	5.07	119.68	111.37

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	160	ARG	Sidechain

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	1617	0	1616	60	1
2	A	3	0	1	0	0
3	A	156	0	0	2	0
All	All	1776	0	1617	60	1

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 19.

All (60) 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:A:41:GLU:CD	1:A:41:GLU:CG	2.14	1.21
1:A:185:ILE:HD11	1:A:209:ILE:HG13	1.19	1.17
1:A:133:MET:HE3	1:A:207:VAL:HG11	1.33	1.08
1:A:185:ILE:HD11	1:A:209:ILE:CG1	1.92	1.00
1:A:41:GLU:CG	1:A:41:GLU:OE1	2.14	0.95
1:A:119:CYS:O	1:A:123:ILE:HD13	1.69	0.92
1:A:133:MET:CE	1:A:207:VAL:HG11	2.05	0.87
1:A:129:THR:HG21	1:A:132:MET:HE3	1.57	0.85
1:A:178:LEU:HD22	1:A:205:VAL:HB	1.58	0.84
1:A:134:TRP:HB3	1:A:166:ILE:HG12	1.59	0.82
1:A:129:THR:CG2	1:A:132:MET:HE3	2.13	0.78
1:A:191:LEU:HD23	1:A:216:VAL:HG11	1.64	0.77
1:A:167:LEU:HD11	1:A:201:VAL:HG12	1.70	0.72
1:A:185:ILE:CD1	1:A:209:ILE:HG13	2.10	0.72
1:A:102:LEU:O	1:A:105:ARG:HG2	1.89	0.72
1:A:206:THR:HG23	1:A:221:TYR:OH	1.90	0.71
1:A:138:ASP:O	1:A:142:THR:HG22	1.91	0.71
1:A:178:LEU:HD22	1:A:205:VAL:CB	2.25	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:41:GLU:OE1	1:A:41:GLU:HG3	1.95	0.66
1:A:102:LEU:O	1:A:105:ARG:HD2	1.96	0.65
1:A:222:ALA:O	1:A:223:ALA:C	2.42	0.63
1:A:131:GLY:O	1:A:132:MET:HE2	1.99	0.62
1:A:201:VAL:HG12	1:A:201:VAL:O	1.99	0.62
1:A:133:MET:HE3	1:A:207:VAL:CG1	2.21	0.61
1:A:95:SER:HA	1:A:123:ILE:HD11	1.82	0.60
1:A:37:ASP:OD1	3:A:408:HOH:O	2.18	0.57
1:A:109:ALA:HB3	1:A:132:MET:CE	2.34	0.57
1:A:206:THR:OG1	1:A:217:VAL:HG13	2.04	0.56
1:A:166:ILE:HD12	1:A:167:LEU:N	2.21	0.56
1:A:179:ASP:O	1:A:207:VAL:HG12	2.06	0.54
1:A:176:LEU:HD23	1:A:178:LEU:HD21	1.88	0.54
1:A:140:ILE:HG13	1:A:148:SER:HA	1.90	0.53
1:A:102:LEU:O	1:A:105:ARG:CG	2.57	0.52
1:A:137:GLU:HG3	1:A:183:LEU:HD21	1.91	0.52
1:A:191:LEU:CD2	1:A:216:VAL:HG11	2.37	0.52
1:A:109:ALA:HB3	1:A:132:MET:HE1	1.92	0.52
1:A:36:GLU:CD	1:A:37:ASP:H	2.19	0.51
1:A:137:GLU:CG	1:A:183:LEU:HD21	2.40	0.51
1:A:206:THR:OG1	1:A:217:VAL:CG1	2.60	0.50
1:A:181:VAL:CG2	1:A:207:VAL:HG13	2.42	0.50
1:A:123:ILE:CD1	1:A:123:ILE:N	2.76	0.49
1:A:183:LEU:HD13	3:A:279:HOH:O	2.13	0.49
1:A:35:LEU:HD12	1:A:35:LEU:N	2.28	0.49
1:A:41:GLU:HG3	1:A:42:ALA:N	2.14	0.48
1:A:129:THR:CG2	1:A:132:MET:CE	2.90	0.47
1:A:35:LEU:N	1:A:35:LEU:CD1	2.77	0.47
1:A:213:GLN:O	1:A:216:VAL:CG1	2.64	0.46
1:A:32:ILE:HD13	1:A:89:MET:HE1	1.98	0.46
1:A:80:LEU:HD23	1:A:105:ARG:HH11	1.82	0.45
1:A:102:LEU:O	1:A:105:ARG:CD	2.62	0.45
1:A:185:ILE:HD11	1:A:209:ILE:CD1	2.43	0.45
1:A:181:VAL:HG22	1:A:207:VAL:HG13	1.99	0.45
1:A:213:GLN:O	1:A:216:VAL:HG12	2.17	0.44
1:A:201:VAL:O	1:A:201:VAL:CG1	2.64	0.44
1:A:191:LEU:CD2	1:A:216:VAL:CG1	2.95	0.44
1:A:167:LEU:CD1	1:A:201:VAL:HG12	2.43	0.43
1:A:166:ILE:HD12	1:A:166:ILE:C	2.43	0.43
1:A:36:GLU:CG	1:A:37:ASP:H	2.33	0.41
1:A:185:ILE:HD13	1:A:213:GLN:NE2	2.36	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:92:LYS:NZ	1:A:138:ASP:OD2	2.47	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:157:ASP:OD2	1:A:160:ARG:CD[3_665]	2.13	0.07

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	221/273 (81%)	217 (98%)	4 (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	156/194 (80%)	145 (93%)	11 (7%)	13	2

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	ARG
1	A	35	LEU
1	A	36	GLU
1	A	41	GLU
1	A	53	ASP
1	A	56	LEU
1	A	105	ARG
1	A	123	ILE
1	A	147	SER
1	A	160	ARG
1	A	167	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	A	43	GLN
1	A	73	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

1 ligand is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	FMT	A	274	-	2,2,2	0.74	0	1,1,1	0.53	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	223/273 (81%)	1.95	92 (41%) 0 0	16, 25, 43, 53	0

All (92) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	223	ALA	6.0
1	A	40	ALA	6.0
1	A	23	ALA	5.4
1	A	42	ALA	5.4
1	A	53	ASP	5.3
1	A	6	ALA	5.2
1	A	17	ALA	5.2
1	A	222	ALA	5.1
1	A	3	LEU	5.1
1	A	4	ARG	4.9
1	A	22	ALA	4.5
1	A	105	ARG	4.2
1	A	41	GLU	4.1
1	A	166	ILE	4.1
1	A	207	VAL	4.0
1	A	151	ALA	4.0
1	A	185	ILE	3.9
1	A	51	LEU	3.9
1	A	81	ALA	3.9
1	A	13	CYS	3.9
1	A	70	THR	3.8
1	A	44	LYS	3.7
1	A	38	GLY	3.7
1	A	206	THR	3.7
1	A	196	ARG	3.6
1	A	18	PRO	3.5
1	A	5	ALA	3.5

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Mol	Chain	Res	Type	RSRZ
1	A	62	VAL	3.5
1	A	45	PRO	3.5
1	A	39	VAL	3.4
1	A	76	ASP	3.4
1	A	82	GLY	3.4
1	A	80	LEU	3.3
1	A	36	GLU	3.2
1	A	52	ARG	3.2
1	A	20	ALA	3.2
1	A	47	ALA	3.2
1	A	127	ASP	3.1
1	A	74	ALA	3.1
1	A	50	ALA	3.0
1	A	147	SER	3.0
1	A	83	THR	2.9
1	A	46	ALA	2.9
1	A	35	LEU	2.9
1	A	100	ILE	2.9
1	A	2	ASN	2.9
1	A	43	GLN	2.9
1	A	102	LEU	2.8
1	A	142	THR	2.8
1	A	216	VAL	2.8
1	A	57	ASP	2.8
1	A	58	PRO	2.7
1	A	56	LEU	2.7
1	A	54	THR	2.7
1	A	14	PRO	2.7
1	A	55	PRO	2.7
1	A	21	PHE	2.7
1	A	75	ARG	2.6
1	A	96	ALA	2.6
1	A	77	LEU	2.6
1	A	88	VAL	2.6
1	A	136	ALA	2.6
1	A	49	ASN	2.5
1	A	26	ALA	2.5
1	A	16	ASP	2.4
1	A	24	ALA	2.4
1	A	27	ALA	2.4
1	A	103	ALA	2.4
1	A	107	VAL	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	15	ALA	2.3
1	A	183	LEU	2.3
1	A	25	ALA	2.3
1	A	84	ALA	2.3
1	A	178	LEU	2.3
1	A	123	ILE	2.3
1	A	60	ARG	2.2
1	A	99	VAL	2.2
1	A	7	GLY	2.2
1	A	138	ASP	2.1
1	A	28	ALA	2.1
1	A	167	LEU	2.1
1	A	59	GLU	2.1
1	A	104	PRO	2.1
1	A	118	VAL	2.1
1	A	19	GLU	2.1
1	A	122	GLU	2.1
1	A	95	SER	2.1
1	A	97	ALA	2.0
1	A	32	ILE	2.0
1	A	85	TYR	2.0
1	A	150	ARG	2.0
1	A	71	ALA	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](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists the number of atoms with occupancy less than 0.9.

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	FMT	A	274	3/3	0.55	0.34	24,24,25,25	3

6.5 Other polymers [i](#)

There are no such residues in this entry.