



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 01:28 AM UTC

PDB ID : 7AGP / pdb_00007agp
Title : Structure of the AcylTransferase domain of Mycoberosic Acid Synthase from Mycobacterium tuberculosis
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Deposited on : 2020-09-23
Resolution : 2.40 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

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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

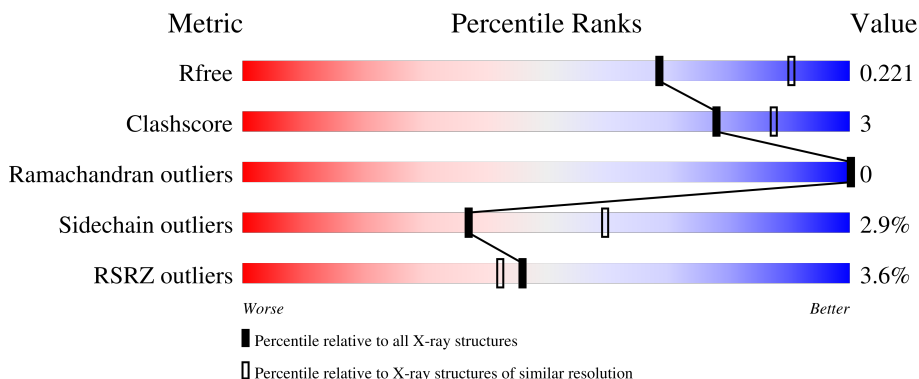
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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}	180053	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (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 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	439	4% 88% 10% .
1	B	439	3% 89% 9% ..

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6588 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 Mycocerosic acid synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	429	Total	C	N	O	S	0	3	0
			3151	1976	556	604	15			
1	B	432	Total	C	N	O	S	0	3	0
			3191	1998	567	610	16			

There are 28 discrepancies between the modelled and reference sequences:

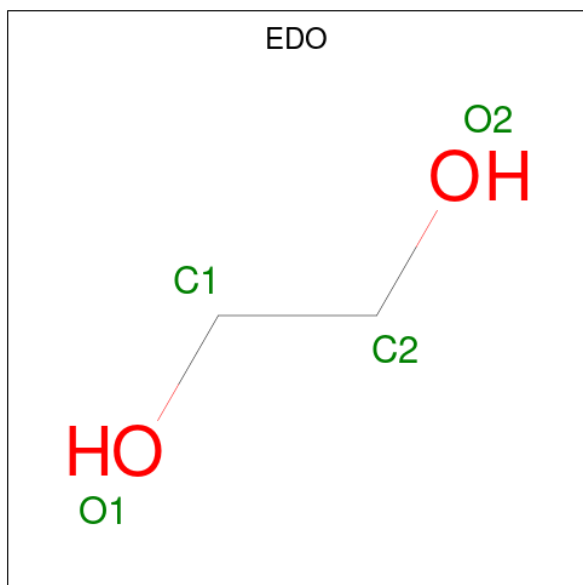
Chain	Residue	Modelled	Actual	Comment	Reference
A	435	MET	-	initiating methionine	UNP Q02251
A	436	GLY	-	expression tag	UNP Q02251
A	437	SER	-	expression tag	UNP Q02251
A	438	SER	-	expression tag	UNP Q02251
A	439	HIS	-	expression tag	UNP Q02251
A	440	HIS	-	expression tag	UNP Q02251
A	441	HIS	-	expression tag	UNP Q02251
A	442	HIS	-	expression tag	UNP Q02251
A	443	HIS	-	expression tag	UNP Q02251
A	444	HIS	-	expression tag	UNP Q02251
A	445	SER	-	expression tag	UNP Q02251
A	446	SER	-	expression tag	UNP Q02251
A	872	GLY	-	cloning artifact	UNP Q02251
A	873	SER	-	cloning artifact	UNP Q02251
B	435	MET	-	initiating methionine	UNP Q02251
B	436	GLY	-	expression tag	UNP Q02251
B	437	SER	-	expression tag	UNP Q02251
B	438	SER	-	expression tag	UNP Q02251
B	439	HIS	-	expression tag	UNP Q02251
B	440	HIS	-	expression tag	UNP Q02251
B	441	HIS	-	expression tag	UNP Q02251
B	442	HIS	-	expression tag	UNP Q02251
B	443	HIS	-	expression tag	UNP Q02251
B	444	HIS	-	expression tag	UNP Q02251
B	445	SER	-	expression tag	UNP Q02251

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Chain	Residue	Modelled	Actual	Comment	Reference
B	446	SER	-	expression tag	UNP Q02251
B	872	GLY	-	cloning artifact	UNP Q02251
B	873	SER	-	cloning artifact	UNP Q02251

- Molecule 2 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			4	2	2		
2	B	1	Total	C	O	0	0
			4	2	2		

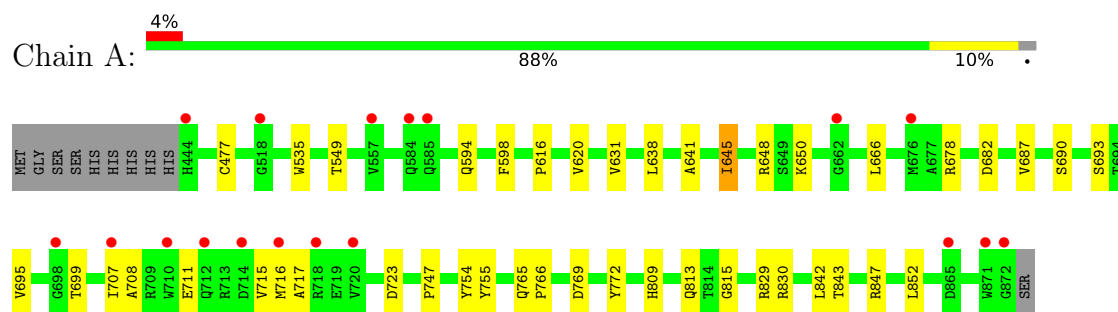
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	107	Total	O	0	0
			107	107		
3	B	131	Total	O	0	0
			131	131		

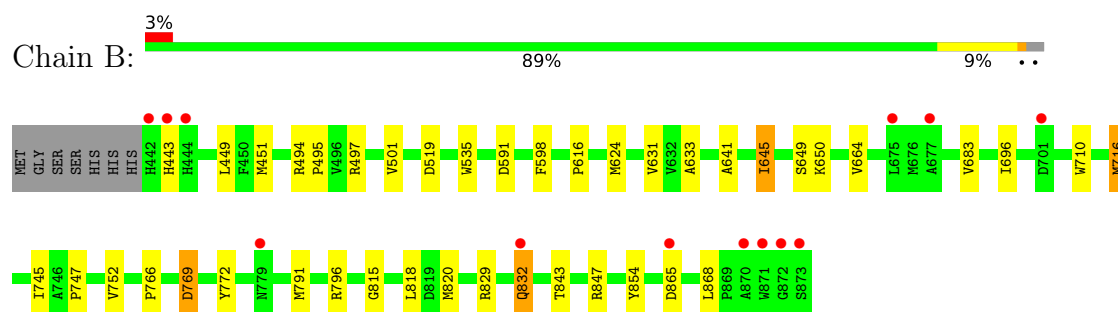
3 Residue-property plots

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: Mycocerosic acid synthase



- Molecule 1: Mycocerosic acid synthase



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	146.03Å 157.14Å 114.25Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.30 – 2.40 49.30 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.1 (49.30-2.40) 99.1 (49.30-2.40)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	0.04	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.10 (at 2.39Å)	Xtriage
Refinement program	BUSTER 2.10.3, REFMAC 5.8.0258	Depositor
R, R_{free}	0.197 , 0.219 0.193 , 0.221	Depositor DCC
R_{free} test set	2534 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å ²)	74.1	Xtriage
Anisotropy	0.346	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 68.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6588	wwPDB-VP
Average B, all atoms (Å ²)	85.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.75% 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: EDO

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	0.87	0/3222	1.12	9/4405 (0.2%)
1	B	0.90	1/3252 (0.0%)	1.13	6/4443 (0.1%)
All	All	0.88	1/6474 (0.0%)	1.13	15/8848 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	716	MET	SD-CE	-7.50	1.60	1.79

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	815	GLY	CA-C-N	6.38	128.83	120.28
1	B	815	GLY	C-N-CA	6.38	128.83	120.28
1	B	497	ARG	N-CA-C	5.68	118.94	110.14
1	A	815	GLY	CA-C-N	5.52	127.68	120.28
1	A	815	GLY	C-N-CA	5.52	127.68	120.28
1	B	443	HIS	CA-C-N	5.47	128.63	120.31
1	B	443	HIS	C-N-CA	5.47	128.63	120.31
1	A	638	LEU	CA-C-N	5.30	127.33	120.44
1	A	638	LEU	C-N-CA	5.30	127.33	120.44
1	A	842	LEU	CA-C-N	5.13	127.15	120.28
1	A	842	LEU	C-N-CA	5.13	127.15	120.28
1	B	769	ASP	CA-CB-CG	5.11	117.71	112.60
1	A	549	THR	N-CA-C	5.06	116.49	111.07
1	A	708	ALA	CA-C-N	5.00	126.98	120.28
1	A	708	ALA	C-N-CA	5.00	126.98	120.28

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	3151	0	3085	18	0
1	B	3191	0	3139	17	0
2	A	4	0	6	0	0
2	B	4	0	6	0	0
3	A	107	0	0	1	0
3	B	131	0	0	0	0
All	All	6588	0	6236	35	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 3.

All (35) 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:620:VAL:HG22	1:A:755:TYR:HB2	1.74	0.69
1:A:809:HIS:CE1	1:A:813:GLN:NE2	2.70	0.60
1:A:687:VAL:HB	1:A:695:VAL:HB	1.86	0.56
1:A:711:GLU:HG3	1:A:717:ALA:HB3	1.88	0.56
1:A:690:SER:OG	1:A:693:SER:HB3	2.10	0.52
1:B:449:LEU:HD23	1:B:501:VAL:HG22	1.92	0.52
1:A:682:ASP:HB2	1:A:699:THR:HG23	1.94	0.49
1:B:843:THR:O	1:B:847:ARG:HG3	2.14	0.47
1:B:451[B]:MET:HE1	1:B:854:TYR:HE1	1.78	0.47
1:B:631:VAL:CG2	1:B:641:ALA:CB	2.91	0.47
1:B:633:ALA:HA	1:B:752:VAL:HG22	1.97	0.47
1:A:535:TRP:CD1	1:A:616:PRO:HB3	2.50	0.47
1:A:631:VAL:CG2	1:A:641:ALA:CB	2.92	0.46
1:A:766:PRO:HB3	1:A:772:TYR:CE1	2.50	0.46
1:A:598:PHE:HB2	1:A:645:ILE:HD13	1.98	0.45
1:B:747:PRO:HG3	1:B:769:ASP:HB3	1.98	0.45
1:B:591:ASP:HA	1:B:649:SER:HB2	1.98	0.45
1:A:847:ARG:NH1	3:A:1003:HOH:O	2.49	0.45
1:B:683:VAL:HG11	1:B:696:ILE:HD12	1.99	0.45
1:B:598:PHE:HB2	1:B:645:ILE:HD13	1.99	0.44
1:A:594:GLN:HE22	1:A:648[A]:ARG:NH2	2.15	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:747:PRO:HG3	1:A:769:ASP:HB3	2.00	0.43
1:B:494:ARG:HB3	1:B:495:PRO:HD2	2.01	0.43
1:A:707:ILE:HG23	1:A:717:ALA:HB1	2.02	0.42
1:B:664:VAL:HG21	1:B:710:TRP:CD2	2.54	0.42
1:B:535:TRP:CD1	1:B:616:PRO:HB3	2.55	0.42
1:A:843:THR:O	1:A:847:ARG:HG3	2.20	0.42
1:B:791:MET:O	1:B:796:ARG:NH2	2.53	0.42
1:B:818:LEU:HB3	1:B:820:MET:HE3	2.02	0.41
1:A:809:HIS:CE1	1:A:813:GLN:HE22	2.39	0.41
1:B:829[A]:ARG:H	1:B:832[A]:GLN:NE2	2.19	0.41
1:A:765:GLN:HA	1:A:766:PRO:HD3	1.89	0.41
1:B:766:PRO:HB3	1:B:772:TYR:CE1	2.56	0.41
1:A:666:LEU:HG	1:A:715:VAL:HG11	2.03	0.40
1:B:624:MET:HE3	1:B:645:ILE:HD11	2.04	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	430/439 (98%)	423 (98%)	7 (2%)	0	100	100
1	B	433/439 (99%)	427 (99%)	6 (1%)	0	100	100
All	All	863/878 (98%)	850 (98%)	13 (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	316/335 (94%)	306 (97%)	10 (3%)	34	56
1	B	320/335 (96%)	311 (97%)	9 (3%)	38	60
All	All	636/670 (95%)	617 (97%)	19 (3%)	37	58

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	477	CYS
1	A	645	ILE
1	A	650	LYS
1	A	678	ARG
1	A	716	MET
1	A	723	ASP
1	A	754	TYR
1	A	829	ARG
1	A	830	ARG
1	A	852	LEU
1	B	519	ASP
1	B	645	ILE
1	B	650	LYS
1	B	716	MET
1	B	745	ILE
1	B	832[A]	GLN
1	B	832[B]	GLN
1	B	865	ASP
1	B	868	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	550	GLN
1	A	601	GLN
1	A	765	GLN
1	A	809	HIS
1	A	813	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](#)

2 ligands are 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	EDO	A	901	-	3,3,3	0.98	0	2,2,2	0.21	0
2	EDO	B	901	-	3,3,3	1.11	0	2,2,2	0.24	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	EDO	A	901	-	-	1/1/1/1	-
2	EDO	B	901	-	-	1/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	901	EDO	O1-C1-C2-O2
2	A	901	EDO	O1-C1-C2-O2

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	429/439 (97%)	0.38	18 (4%)	40 36	37, 84, 112, 134	2 (0%)
1	B	432/439 (98%)	0.27	13 (3%)	52 48	32, 82, 113, 131	3 (0%)
All	All	861/878 (98%)	0.33	31 (3%)	46 42	32, 83, 113, 134	5 (0%)

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	442	HIS	4.9
1	A	871	TRP	4.4
1	B	871	TRP	3.7
1	B	873	SER	3.6
1	A	444	HIS	3.2
1	B	675	LEU	3.1
1	A	584	GLN	3.0
1	B	444	HIS	3.0
1	A	710	TRP	2.9
1	B	832[A]	GLN	2.8
1	A	585	GLN	2.8
1	A	714	ASP	2.8
1	B	872	GLY	2.8
1	A	720	VAL	2.7
1	A	712	GLN	2.7
1	A	716	MET	2.6
1	B	677	ALA	2.6
1	A	718	ARG	2.6
1	B	701	ASP	2.6
1	A	865	ASP	2.4
1	B	779	ASN	2.2
1	B	443	HIS	2.2
1	B	870	ALA	2.2
1	A	707	ILE	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	676	MET	2.1
1	A	518	GLY	2.1
1	B	865	ASP	2.1
1	A	662	GLY	2.1
1	A	872	GLY	2.1
1	A	557	VAL	2.0
1	A	698	GLY	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.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	EDO	B	901	4/4	0.80	0.28	88,89,90,90	0
2	EDO	A	901	4/4	0.91	0.15	98,99,99,100	0

6.5 Other polymers [i](#)

There are no such residues in this entry.