



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 08:27 AM UTC

PDB ID : 3P3B / pdb_00003p3b
Title : CRYSTAL STRUCTURE OF Galacturonate DEHYDRATASE FROM
GEOBACILLUS SP. COMPLEXED WITH D-TARTRATE
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GXRC)
Deposited on : 2010-10-04
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

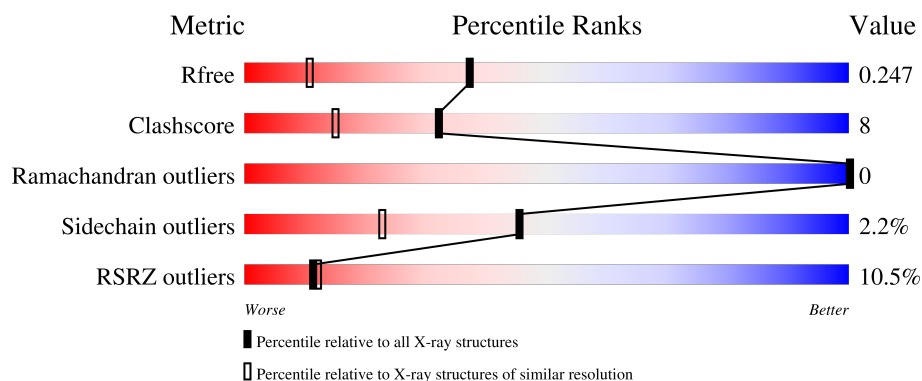
1 Overall quality at a glance

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	392	3% 83% 12% • •
1	B	392	17% 76% 16% • 7%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6208 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 Mandelate racemase/muconate lactonizing protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	377	Total	C	N	O	S	0	1	0
			2990	1893	522	565	10			
1	B	364	Total	C	N	O	S	0	0	0
			2878	1824	503	541	10			

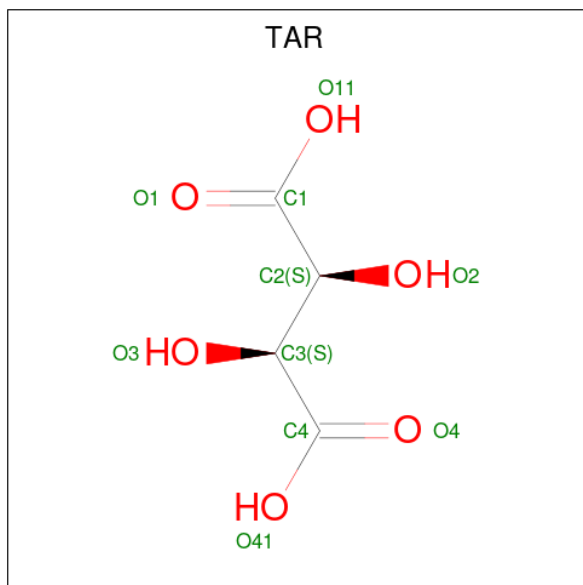
There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	2	SER	-	expression tag	UNP D3EID5
A	3	LEU	-	expression tag	UNP D3EID5
A	385	GLU	-	expression tag	UNP D3EID5
A	386	GLY	-	expression tag	UNP D3EID5
A	387	HIS	-	expression tag	UNP D3EID5
A	388	HIS	-	expression tag	UNP D3EID5
A	389	HIS	-	expression tag	UNP D3EID5
A	390	HIS	-	expression tag	UNP D3EID5
A	391	HIS	-	expression tag	UNP D3EID5
A	392	HIS	-	expression tag	UNP D3EID5
B	2	SER	-	expression tag	UNP D3EID5
B	3	LEU	-	expression tag	UNP D3EID5
B	385	GLU	-	expression tag	UNP D3EID5
B	386	GLY	-	expression tag	UNP D3EID5
B	387	HIS	-	expression tag	UNP D3EID5
B	388	HIS	-	expression tag	UNP D3EID5
B	389	HIS	-	expression tag	UNP D3EID5
B	390	HIS	-	expression tag	UNP D3EID5
B	391	HIS	-	expression tag	UNP D3EID5
B	392	HIS	-	expression tag	UNP D3EID5

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	2	Total Mg 2 2	0	0
2	B	2	Total Mg 2 2	0	0

- Molecule 3 is D(-)-TARTARIC ACID (CCD ID: TAR) (formula: C₄H₆O₆).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 10 4 6	0	0
3	B	1	Total C O 10 4 6	0	0

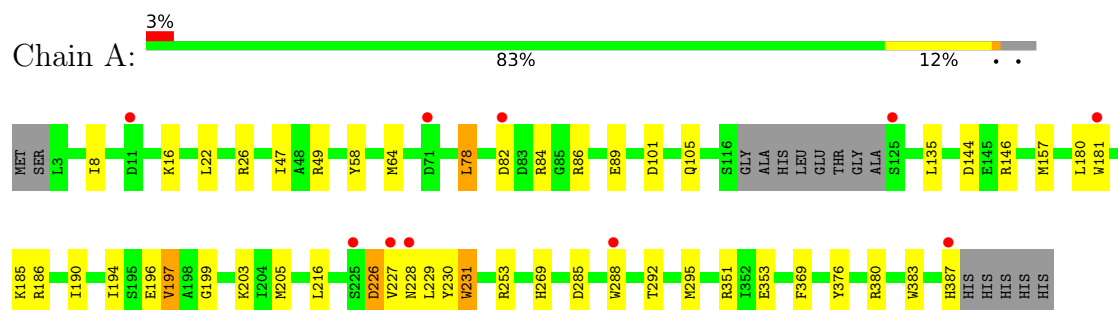
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	225	Total O 225 225	0	0
4	B	91	Total O 91 91	0	0

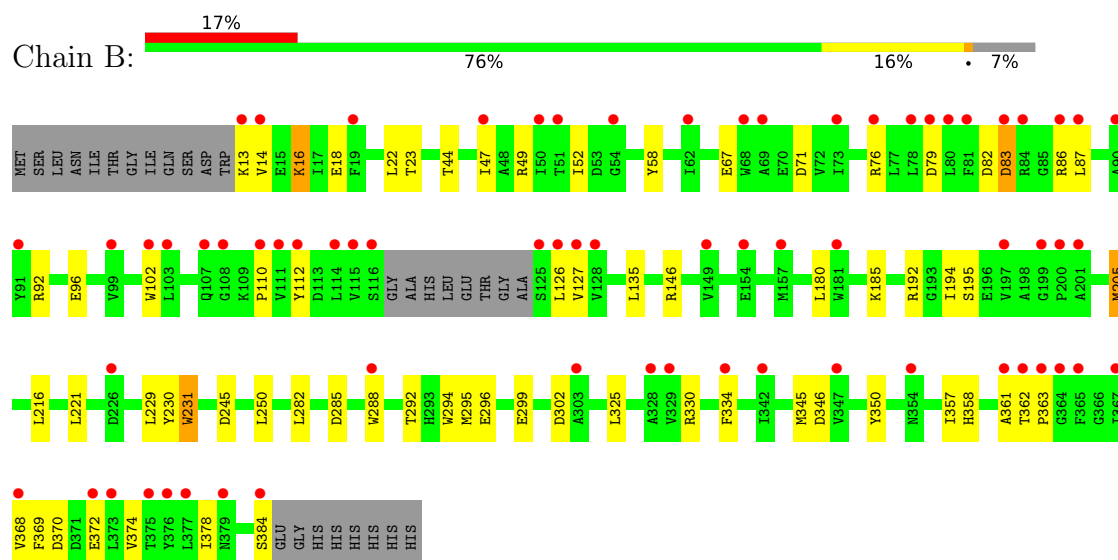
3 Residue-property plots

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: Mandelate racemase/muconate lactonizing protein



- Molecule 1: Mandelate racemase/muconate lactonizing protein



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	158.86Å 66.05Å 75.66Å 90.00° 95.48° 90.00°	Depositor
Resolution (Å)	41.20 – 1.65 41.20 – 1.65	Depositor EDS
% Data completeness (in resolution range)	93.3 (41.20-1.65) 93.3 (41.20-1.65)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.02 (at 1.65Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.5_2)	Depositor
R, R_{free}	0.216 , 0.252 0.214 , 0.247	Depositor DCC
R_{free} test set	4384 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	26.7	Xtriage
Anisotropy	0.128	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 30.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6208	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.13% 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](#)

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

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	0.52	0/3059	0.82	2/4145 (0.0%)
1	B	0.40	0/2944	0.74	0/3988
All	All	0.46	0/6003	0.78	2/8133 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	199	GLY	CA-C-N	6.27	126.18	119.28
1	A	199	GLY	C-N-CA	6.27	126.18	119.28

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	2990	0	2882	38	0
1	B	2878	0	2785	57	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	10	0	4	0	0
3	B	10	0	4	0	0
4	A	225	0	0	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	91	0	0	0	0
All	All	6208	0	5675	93	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 8.

All (93) 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:292:THR:HA	1:A:295:MET:HE2	1.46	0.98
1:A:157:MET:SD	1:A:197:VAL:HG13	2.14	0.87
1:B:292:THR:HA	1:B:295:MET:HE2	1.63	0.81
1:A:84:ARG:HD3	1:A:86:ARG:HH21	1.49	0.77
1:A:292:THR:HA	1:A:295:MET:CE	2.15	0.77
1:A:22:LEU:HD22	1:A:47:ILE:HD12	1.67	0.76
1:B:13:LYS:HG3	1:B:76:ARG:HD2	1.66	0.76
1:B:350:TYR:HB3	1:B:357:ILE:HD11	1.67	0.76
1:B:18:GLU:HB3	1:B:378:ILE:HD13	1.72	0.72
1:B:22:LEU:HD22	1:B:47:ILE:CD1	2.23	0.69
1:B:285:ASP:HB3	1:B:288:TRP:O	1.95	0.67
1:A:47:ILE:HD13	1:A:369:PHE:CE1	2.30	0.66
1:B:76:ARG:HB2	1:B:79:ASP:OD2	1.96	0.65
1:A:22:LEU:HD22	1:A:47:ILE:CD1	2.28	0.63
1:A:226:ASP:N	1:A:226:ASP:OD1	2.31	0.63
1:A:16:LYS:NZ	1:A:387:HIS:HA	2.13	0.62
1:B:16:LYS:HD3	1:B:18:GLU:HG3	1.81	0.61
1:A:285:ASP:HB3	1:A:288:TRP:O	2.03	0.58
1:B:185:LYS:NZ	1:B:185:LYS:HB3	2.18	0.57
1:B:346:ASP:HB3	1:B:368:VAL:HB	1.84	0.57
1:B:135:LEU:HD13	1:B:194:ILE:HD13	1.87	0.57
1:B:49:ARG:HD3	1:B:58:TYR:CZ	2.41	0.56
1:B:180:LEU:HD11	1:B:216:LEU:HD11	1.88	0.55
1:A:135:LEU:HD13	1:A:194:ILE:HD13	1.88	0.55
1:B:22:LEU:HD22	1:B:47:ILE:HD13	1.87	0.55
1:A:8:ILE:CG2	1:A:78:LEU:HD13	2.37	0.55
1:B:13:LYS:CG	1:B:76:ARG:HD2	2.35	0.52
1:A:47:ILE:HD13	1:A:369:PHE:HE1	1.72	0.52
1:B:302:ASP:HB3	1:B:330:ARG:NH2	2.25	0.52
1:A:16:LYS:HZ3	1:A:387:HIS:HA	1.74	0.51
1:B:23:THR:HG22	1:B:44:THR:HG22	1.92	0.51
1:B:345:MET:HG3	1:B:369:PHE:CE1	2.45	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:351:ARG:NH1	1:A:353:GLU:OE1	2.43	0.51
1:A:181:TRP:CZ2	1:A:185:LYS:HD2	2.47	0.50
1:B:86:ARG:CZ	1:B:296:GLU:HG2	2.42	0.50
1:B:282:LEU:HD22	1:B:282:LEU:N	2.26	0.50
1:A:84:ARG:HD3	1:A:86:ARG:NH2	2.23	0.49
1:A:203:LYS:HD3	1:A:230:TYR:CD1	2.47	0.49
1:A:180:LEU:CD2	1:B:216:LEU:HG	2.42	0.49
1:A:82:ASP:HB3	1:A:84:ARG:H	1.77	0.49
1:B:192:ARG:O	1:B:195:SER:HB3	2.12	0.49
1:B:245:ASP:OD1	1:B:245:ASP:C	2.55	0.48
1:B:362:THR:HB	1:B:363:PRO:CD	2.44	0.48
1:B:294:TRP:HB3	1:B:325:LEU:HD21	1.95	0.48
1:B:110:PRO:HG3	1:B:361:ALA:O	2.15	0.47
1:B:374:VAL:O	1:B:378:ILE:HG13	2.15	0.47
1:B:221:LEU:HD13	1:B:250:LEU:HD21	1.97	0.47
1:B:205:MET:HG2	1:B:231:TRP:CE2	2.49	0.47
1:B:230:TYR:CE2	1:B:334:PHE:HE2	2.32	0.46
1:A:157:MET:SD	1:A:197:VAL:CG1	2.96	0.46
1:B:83:ASP:N	1:B:83:ASP:OD1	2.48	0.46
1:B:82:ASP:HB3	1:B:86:ARG:H	1.79	0.46
1:A:49:ARG:HD3	1:A:58:TYR:CZ	2.50	0.46
1:B:13:LYS:HE3	1:B:76:ARG:HD3	1.98	0.46
1:B:146:ARG:N	1:B:146:ARG:HD2	2.30	0.46
1:A:186:ARG:O	1:A:190:ILE:HG12	2.16	0.46
1:B:92:ARG:HA	1:B:96:GLU:HB3	1.97	0.46
1:A:84:ARG:HD3	1:A:86:ARG:HE	1.81	0.45
1:A:180:LEU:HD22	1:B:216:LEU:HG	1.98	0.45
1:B:231:TRP:CD1	1:B:231:TRP:C	2.95	0.45
1:A:64:MET:HE3	1:A:383:TRP:CH2	2.51	0.45
1:B:205:MET:HG2	1:B:231:TRP:CD2	2.52	0.45
1:B:87:LEU:HD12	1:B:87:LEU:HA	1.84	0.45
1:B:146:ARG:N	1:B:146:ARG:CD	2.80	0.45
1:B:146:ARG:H	1:B:146:ARG:CD	2.29	0.45
1:A:376:TYR:CZ	1:A:380:ARG:HD3	2.53	0.44
1:B:205:MET:N	1:B:205:MET:HE3	2.32	0.44
1:B:14:VAL:HA	1:B:52:ILE:HG12	1.99	0.44
1:B:86:ARG:O	1:B:87:LEU:C	2.62	0.43
1:B:112:TYR:CG	1:B:126:LEU:HD22	2.54	0.43
1:A:16:LYS:HZ1	1:A:387:HIS:HA	1.82	0.43
1:A:101:ASP:O	1:A:105:GLN:HG2	2.19	0.42
1:A:144:ASP:OD1	1:A:146:ARG:HB2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:227:VAL:HG12	1:A:228:ASN:N	2.32	0.42
1:B:205:MET:HG2	1:B:231:TRP:CD1	2.54	0.42
1:B:86:ARG:NE	1:B:296:GLU:HG2	2.34	0.42
1:B:82:ASP:CG	1:B:83:ASP:N	2.77	0.42
1:A:185:LYS:HE2	1:A:185:LYS:HB3	1.64	0.42
1:B:127:VAL:HG22	1:B:358:HIS:ND1	2.35	0.42
1:B:370:ASP:OD1	1:B:370:ASP:C	2.63	0.42
1:B:82:ASP:CB	1:B:86:ARG:O	2.67	0.42
1:B:372:GLU:CD	1:B:372:GLU:N	2.78	0.41
1:A:205:MET:HG2	1:A:231:TRP:CG	2.54	0.41
1:B:67:GLU:CD	1:B:67:GLU:H	2.29	0.41
1:B:86:ARG:NH2	1:B:299:GLU:OE2	2.53	0.41
1:A:180:LEU:HD11	1:A:216:LEU:HD11	2.03	0.41
1:A:205:MET:HG2	1:A:231:TRP:CD1	2.56	0.41
1:A:269:HIS:HD2	4:A:477:HOH:O	2.04	0.41
1:B:52:ILE:HB	1:B:102:TRP:CE2	2.56	0.41
1:B:18:GLU:HG2	1:B:384:SER:HB3	2.03	0.40
1:A:26:ARG:HB2	4:A:555:HOH:O	2.21	0.40
1:B:205:MET:HG2	1:B:231:TRP:CG	2.56	0.40
1:A:253:ARG:HG3	4:A:597:HOH:O	2.21	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	374/392 (95%)	369 (99%)	5 (1%)	0	100	100
1	B	360/392 (92%)	349 (97%)	11 (3%)	0	100	100
All	All	734/784 (94%)	718 (98%)	16 (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	306/316 (97%)	299 (98%)	7 (2%)	44	21
1	B	294/316 (93%)	288 (98%)	6 (2%)	48	26
All	All	600/632 (95%)	587 (98%)	13 (2%)	45	23

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	78	LEU
1	A	89	GLU
1	A	196	GLU
1	A	197	VAL
1	A	226	ASP
1	A	229	LEU
1	A	231	TRP
1	B	16	LYS
1	B	71	ASP
1	B	83	ASP
1	B	205	MET
1	B	229	LEU
1	B	231	TRP

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	107	GLN
1	B	107	GLN

5.3.3 RNA ⓘ

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](#)

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 for Mogul analysis.

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	$\# Z > 2$
3	TAR	A	395	2	9,9,9	1.19	0	12,12,12	1.30	1 (8%)
3	TAR	B	395	2	9,9,9	1.04	0	12,12,12	1.22	1 (8%)

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
3	TAR	A	395	2	-	4/12/12/12	-
3	TAR	B	395	2	-	4/12/12/12	-

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	395	TAR	C3-C2-C1	2.35	115.04	109.82
3	A	395	TAR	O11-C1-C2	2.35	119.83	113.31

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	395	TAR	O11-C1-C2-C3
3	A	395	TAR	O1-C1-C2-C3
3	B	395	TAR	O1-C1-C2-C3
3	B	395	TAR	O11-C1-C2-C3
3	A	395	TAR	O1-C1-C2-O2
3	A	395	TAR	O11-C1-C2-O2
3	B	395	TAR	O1-C1-C2-O2
3	B	395	TAR	O11-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	377/392 (96%)	0.17	10 (2%) 56 61	16, 26, 41, 60	1 (0%)
1	B	364/392 (92%)	1.17	68 (18%) 3 4	23, 41, 60, 86	0
All	All	741/784 (94%)	0.66	78 (10%) 11 12	16, 32, 57, 86	1 (0%)

All (78) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	197	VAL	4.1
1	B	181	TRP	4.0
1	B	288	TRP	3.9
1	B	157	MET	3.7
1	A	181	TRP	3.6
1	A	82	ASP	3.2
1	B	126	LEU	3.2
1	B	111	VAL	3.2
1	B	125	SER	3.1
1	B	373	LEU	3.1
1	A	228	ASN	3.1
1	B	78	LEU	3.0
1	B	372	GLU	3.0
1	A	387	HIS	3.0
1	B	47	ILE	3.0
1	B	329	VAL	2.9
1	B	73	ILE	2.9
1	B	226	ASP	2.9
1	B	76	ARG	2.8
1	B	377	LEU	2.8
1	A	11	ASP	2.8
1	A	288	TRP	2.7
1	B	127	VAL	2.7
1	B	69	ALA	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	87	LEU	2.7
1	A	125	SER	2.7
1	B	51	THR	2.7
1	B	80	LEU	2.6
1	B	86	ARG	2.6
1	B	347	VAL	2.6
1	B	375	THR	2.6
1	B	13	LYS	2.6
1	B	116	SER	2.6
1	B	199	GLY	2.5
1	B	81	PHE	2.5
1	B	365	PHE	2.5
1	A	227	VAL	2.5
1	B	342	ILE	2.5
1	B	68	TRP	2.5
1	B	14	VAL	2.5
1	B	361	ALA	2.5
1	B	102	TRP	2.4
1	B	334	PHE	2.4
1	B	50	ILE	2.4
1	B	368	VAL	2.4
1	B	303	ALA	2.4
1	B	107	GLN	2.4
1	B	200	PRO	2.4
1	A	71	ASP	2.4
1	B	115	VAL	2.4
1	B	154	GLU	2.3
1	B	103	LEU	2.3
1	B	91	TYR	2.3
1	B	354	ASN	2.3
1	B	367	ILE	2.3
1	B	364	GLY	2.3
1	B	384	SER	2.3
1	B	376	TYR	2.3
1	B	110	PRO	2.2
1	B	84	ARG	2.2
1	B	114	LEU	2.2
1	B	79	ASP	2.2
1	B	363	PRO	2.2
1	B	112	TYR	2.1
1	B	54	GLY	2.1
1	B	379	ASN	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	225	SER	2.1
1	B	83	ASP	2.1
1	B	128	VAL	2.1
1	B	90	ALA	2.0
1	B	201	ALA	2.0
1	B	328	ALA	2.0
1	B	149	VAL	2.0
1	B	362	THR	2.0
1	B	108	GLY	2.0
1	B	19	PHE	2.0
1	B	62	ILE	2.0
1	B	99	VAL	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(Å ²)	Q<0.9
3	TAR	B	395	10/10	0.93	0.08	26,30,34,36	0
3	TAR	A	395	10/10	0.96	0.06	19,23,25,29	0
2	MG	B	393	1/1	0.97	0.06	28,28,28,28	0
2	MG	B	394	1/1	0.97	0.07	27,27,27,27	0
2	MG	A	393	1/1	0.98	0.07	21,21,21,21	0
2	MG	A	394	1/1	1.00	0.07	22,22,22,22	0

6.5 Other polymers [i](#)

There are no such residues in this entry.