



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 13, 2026 – 11:26 PM UTC

PDB ID : 4ADM / pdb_00004adm
Title : Crystal structure of Rv1098c in complex with meso-tartrate
Authors : Mechaly, A.E.; Haouz, A.; Miras, I.; Weber, P.; Shepard, W.; Cole, S.; Alzari, P.M.; Bellinzoni, M.
Deposited on : 2011-12-27
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

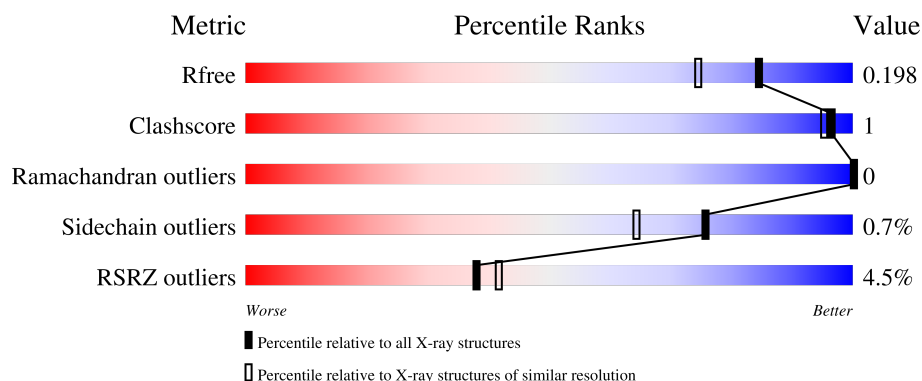
1 Overall quality at a glance

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	495	2% 87% 9%
1	B	495	2% 86% 9%
1	C	495	6% 88% 8%
1	D	495	6% 87% 9%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	SRT	C	1466	-	X	-	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 15056 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 FUMARATE HYDRATASE CLASS II.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	450	Total	C	N	O	S	0	2	0
			3348	2088	607	642	11			
1	B	449	Total	C	N	O	S	0	0	0
			3339	2080	607	641	11			
1	C	456	Total	C	N	O	S	0	1	0
			3334	2077	605	640	12			
1	D	451	Total	C	N	O	S	0	1	0
			3314	2071	599	632	12			

There are 88 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-21	MET	-	expression tag	UNP O53446
A	-20	SER	-	expression tag	UNP O53446
A	-19	TYR	-	expression tag	UNP O53446
A	-18	TYR	-	expression tag	UNP O53446
A	-17	HIS	-	expression tag	UNP O53446
A	-16	HIS	-	expression tag	UNP O53446
A	-15	HIS	-	expression tag	UNP O53446
A	-14	HIS	-	expression tag	UNP O53446
A	-13	HIS	-	expression tag	UNP O53446
A	-12	HIS	-	expression tag	UNP O53446
A	-11	LEU	-	expression tag	UNP O53446
A	-10	GLU	-	expression tag	UNP O53446
A	-9	SER	-	expression tag	UNP O53446
A	-8	THR	-	expression tag	UNP O53446
A	-7	SER	-	expression tag	UNP O53446
A	-6	LEU	-	expression tag	UNP O53446
A	-5	TYR	-	expression tag	UNP O53446
A	-4	LYS	-	expression tag	UNP O53446
A	-3	LYS	-	expression tag	UNP O53446
A	-2	ALA	-	expression tag	UNP O53446
A	-1	GLY	-	expression tag	UNP O53446

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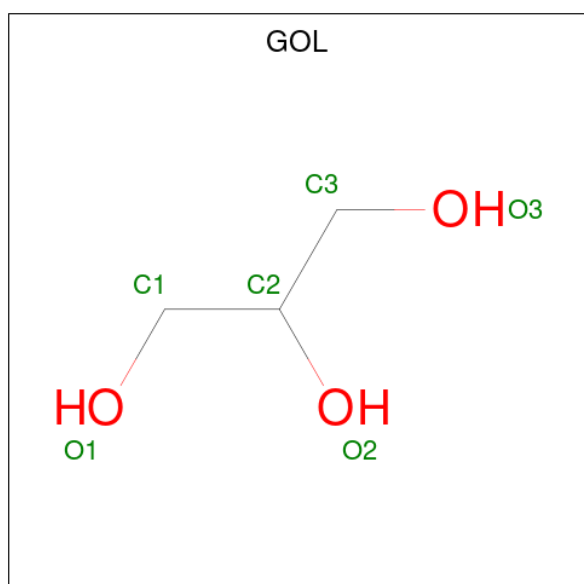
Chain	Residue	Modelled	Actual	Comment	Reference
A	0	SER	-	expression tag	UNP O53446
B	-21	MET	-	expression tag	UNP O53446
B	-20	SER	-	expression tag	UNP O53446
B	-19	TYR	-	expression tag	UNP O53446
B	-18	TYR	-	expression tag	UNP O53446
B	-17	HIS	-	expression tag	UNP O53446
B	-16	HIS	-	expression tag	UNP O53446
B	-15	HIS	-	expression tag	UNP O53446
B	-14	HIS	-	expression tag	UNP O53446
B	-13	HIS	-	expression tag	UNP O53446
B	-12	HIS	-	expression tag	UNP O53446
B	-11	LEU	-	expression tag	UNP O53446
B	-10	GLU	-	expression tag	UNP O53446
B	-9	SER	-	expression tag	UNP O53446
B	-8	THR	-	expression tag	UNP O53446
B	-7	SER	-	expression tag	UNP O53446
B	-6	LEU	-	expression tag	UNP O53446
B	-5	TYR	-	expression tag	UNP O53446
B	-4	LYS	-	expression tag	UNP O53446
B	-3	LYS	-	expression tag	UNP O53446
B	-2	ALA	-	expression tag	UNP O53446
B	-1	GLY	-	expression tag	UNP O53446
B	0	SER	-	expression tag	UNP O53446
C	-21	MET	-	expression tag	UNP O53446
C	-20	SER	-	expression tag	UNP O53446
C	-19	TYR	-	expression tag	UNP O53446
C	-18	TYR	-	expression tag	UNP O53446
C	-17	HIS	-	expression tag	UNP O53446
C	-16	HIS	-	expression tag	UNP O53446
C	-15	HIS	-	expression tag	UNP O53446
C	-14	HIS	-	expression tag	UNP O53446
C	-13	HIS	-	expression tag	UNP O53446
C	-12	HIS	-	expression tag	UNP O53446
C	-11	LEU	-	expression tag	UNP O53446
C	-10	GLU	-	expression tag	UNP O53446
C	-9	SER	-	expression tag	UNP O53446
C	-8	THR	-	expression tag	UNP O53446
C	-7	SER	-	expression tag	UNP O53446
C	-6	LEU	-	expression tag	UNP O53446
C	-5	TYR	-	expression tag	UNP O53446
C	-4	LYS	-	expression tag	UNP O53446
C	-3	LYS	-	expression tag	UNP O53446

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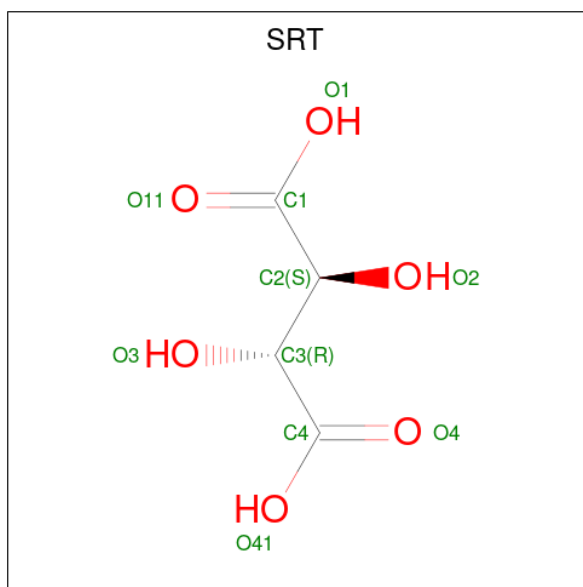
Chain	Residue	Modelled	Actual	Comment	Reference
C	-2	ALA	-	expression tag	UNP O53446
C	-1	GLY	-	expression tag	UNP O53446
C	0	SER	-	expression tag	UNP O53446
D	-21	MET	-	expression tag	UNP O53446
D	-20	SER	-	expression tag	UNP O53446
D	-19	TYR	-	expression tag	UNP O53446
D	-18	TYR	-	expression tag	UNP O53446
D	-17	HIS	-	expression tag	UNP O53446
D	-16	HIS	-	expression tag	UNP O53446
D	-15	HIS	-	expression tag	UNP O53446
D	-14	HIS	-	expression tag	UNP O53446
D	-13	HIS	-	expression tag	UNP O53446
D	-12	HIS	-	expression tag	UNP O53446
D	-11	LEU	-	expression tag	UNP O53446
D	-10	GLU	-	expression tag	UNP O53446
D	-9	SER	-	expression tag	UNP O53446
D	-8	THR	-	expression tag	UNP O53446
D	-7	SER	-	expression tag	UNP O53446
D	-6	LEU	-	expression tag	UNP O53446
D	-5	TYR	-	expression tag	UNP O53446
D	-4	LYS	-	expression tag	UNP O53446
D	-3	LYS	-	expression tag	UNP O53446
D	-2	ALA	-	expression tag	UNP O53446
D	-1	GLY	-	expression tag	UNP O53446
D	0	SER	-	expression tag	UNP O53446

- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

- Molecule 3 is S,R MESO-TARTARIC ACID (CCD ID: SRT) (formula: $C_4H_6O_6$).



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

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	517	Total	O	0	0
			517	517		
4	B	411	Total	O	0	0
			411	411		
4	C	392	Total	O	0	0
			392	392		
4	D	369	Total	O	0	0
			369	369		

- Molecule 1: FUMARATE HYDRATASE CLASS II





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	188.94Å 96.49Å 139.65Å 90.00° 112.44° 90.00°	Depositor
Resolution (Å)	24.16 – 1.65 24.16 – 1.65	Depositor EDS
% Data completeness (in resolution range)	99.5 (24.16-1.65) 99.7 (24.16-1.65)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.25 (at 1.65Å)	Xtriage
Refinement program	BUSTER 2.11.1	Depositor
R, R_{free}	0.163 , 0.185 0.176 , 0.198	Depositor DCC
R_{free} test set	14032 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	18.7	Xtriage
Anisotropy	0.087	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 52.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	15056	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.87% 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: SRT, GOL

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	0.81	0/3400	1.03	6/4623 (0.1%)
1	B	0.80	1/3385 (0.0%)	1.04	6/4601 (0.1%)
1	C	0.83	1/3384 (0.0%)	1.06	4/4607 (0.1%)
1	D	0.84	1/3364 (0.0%)	1.07	3/4577 (0.1%)
All	All	0.82	3/13533 (0.0%)	1.05	19/18408 (0.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	298	MET	SD-CE	-6.97	1.62	1.79
1	C	298	MET	SD-CE	-6.87	1.62	1.79
1	B	341	ILE	CG1-CD1	-6.49	1.26	1.51

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	141	ASP	CA-CB-CG	6.52	119.12	112.60
1	C	141	ASP	CA-CB-CG	6.48	119.08	112.60
1	B	141	ASP	CA-CB-CG	6.38	118.98	112.60
1	A	265	PHE	CA-CB-CG	-5.88	107.92	113.80
1	A	141	ASP	CA-CB-CG	5.76	118.36	112.60
1	C	265	PHE	CA-CB-CG	-5.72	108.08	113.80
1	B	265	PHE	CA-CB-CG	-5.56	108.24	113.80
1	A	383	PHE	CA-CB-CG	5.53	119.33	113.80
1	D	356	PHE	CA-CB-CG	-5.52	108.28	113.80
1	D	265	PHE	CA-CB-CG	-5.48	108.32	113.80
1	B	184	GLY	N-CA-C	-5.33	105.72	112.54
1	C	362	ILE	N-CA-CB	5.23	114.53	110.45
1	A	190	ASP	CA-CB-CG	5.21	117.81	112.60
1	A	184	GLY	N-CA-C	-5.18	105.91	112.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	313	ASP	CA-CB-CG	5.16	117.76	112.60
1	B	383	PHE	CA-CB-CG	5.15	118.95	113.80
1	B	111	ASN	CA-CB-CG	5.14	117.74	112.60
1	A	200	PHE	CA-CB-CG	-5.09	108.71	113.80
1	B	200	PHE	CA-CB-CG	-5.02	108.78	113.80

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	3348	0	3385	6	0
1	B	3339	0	3373	7	0
1	C	3334	0	3347	7	0
1	D	3314	0	3344	8	0
2	A	6	0	8	0	0
2	B	6	0	8	0	0
3	C	10	0	4	0	0
3	D	10	0	4	0	0
4	A	517	0	0	1	1
4	B	411	0	0	2	0
4	C	392	0	0	0	0
4	D	369	0	0	0	0
All	All	15056	0	13473	27	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 1.

All (27) 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:282:ARG:HG2	1:A:341:ILE:HD13	1.55	0.86
1:D:165:GLN:OE1	1:D:381:ARG:NH1	2.19	0.76
1:B:360:VAL:HG22	4:B:2318:HOH:O	1.87	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:360:VAL:HG22	4:A:2411:HOH:O	1.89	0.73
1:C:15:ASP:HB3	1:C:34:ARG:HH21	1.67	0.59
1:C:282:ARG:HG2	1:C:341:ILE:HD12	1.89	0.55
1:C:320:ILE:O	1:C:322:PRO:HD3	2.08	0.54
1:D:320:ILE:O	1:D:322:PRO:HD3	2.09	0.53
1:D:282:ARG:HG2	1:D:341:ILE:HD12	1.93	0.51
1:B:74:GLU:HG3	1:B:75:LYS:N	2.29	0.48
1:C:408:ILE:O	1:C:411:PRO:HD2	2.14	0.47
1:D:408:ILE:O	1:D:411:PRO:HD2	2.15	0.46
1:B:433:LYS:HD2	4:B:2372:HOH:O	2.17	0.45
1:C:465:LYS:HD3	1:D:233:THR:O	2.17	0.44
1:A:71:LEU:HD21	1:A:79:ILE:HD12	2.01	0.43
1:A:70:LEU:HD11	1:A:238:PRO:HG3	2.02	0.42
1:B:71:LEU:HD21	1:B:79:ILE:HD12	2.02	0.42
1:C:224:LEU:HD12	1:C:246:VAL:HG22	2.02	0.42
1:C:27:LEU:HD12	1:C:94:GLN:HG3	2.02	0.42
1:D:27:LEU:HD12	1:D:94:GLN:HG3	2.01	0.41
1:A:329:LEU:HD23	1:A:329:LEU:HA	1.91	0.41
1:B:70:LEU:HD11	1:B:238:PRO:HG3	2.02	0.41
1:B:308:GLU:HB3	1:B:398:LEU:HD21	2.03	0.41
1:A:364:MET:HE3	1:A:364:MET:HB3	1.99	0.41
1:D:238:PRO:HD2	1:D:241:PHE:HB2	2.03	0.41
1:B:97:ILE:HD13	1:B:107:SER:HB3	2.02	0.40
1:D:224:LEU:HD12	1:D:246:VAL:HG22	2.03	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)
4:A:2333:HOH:O	4:A:2333:HOH:O[2_656]	1.88	0.32

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	448/495 (90%)	435 (97%)	13 (3%)	0	100	100
1	B	445/495 (90%)	432 (97%)	13 (3%)	0	100	100
1	C	455/495 (92%)	446 (98%)	9 (2%)	0	100	100
1	D	448/495 (90%)	439 (98%)	9 (2%)	0	100	100
All	All	1796/1980 (91%)	1752 (98%)	44 (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	344/383 (90%)	341 (99%)	3 (1%)	70	56
1	B	343/383 (90%)	340 (99%)	3 (1%)	70	56
1	C	339/383 (88%)	338 (100%)	1 (0%)	86	80
1	D	338/383 (88%)	335 (99%)	3 (1%)	70	56
All	All	1364/1532 (89%)	1354 (99%)	10 (1%)	76	64

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	142[A]	THR
1	A	142[B]	THR
1	A	379	VAL
1	B	315	GLN
1	B	379	VAL
1	B	396	GLU
1	C	329	LEU
1	D	13	GLU
1	D	329	LEU
1	D	379	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	31	GLN
1	A	90	GLN
1	A	252	GLN
1	A	310	GLN
1	A	339	GLN
1	A	368	ASN
1	A	394	ASN
1	B	31	GLN
1	B	90	GLN
1	B	252	GLN
1	B	315	GLN
1	B	326	ASN
1	B	339	GLN
1	B	368	ASN
1	C	140	ASN
1	C	339	GLN
1	C	368	ASN
1	C	397	HIS
1	D	38	ASN
1	D	339	GLN
1	D	368	ASN

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](#)

4 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	GOL	A	1468	-	5,5,5	0.47	0	5,5,5	0.73	0
2	GOL	B	1467	-	5,5,5	0.44	0	5,5,5	1.55	1 (20%)
3	SRT	D	1466	-	9,9,9	1.52	1 (11%)	12,12,12	1.12	1 (8%)
3	SRT	C	1466	-	9,9,9	1.57	2 (22%)	12,12,12	1.49	2 (16%)

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	GOL	A	1468	-	-	4/4/4/4	-
2	GOL	B	1467	-	-	3/4/4/4	-
3	SRT	D	1466	-	-	10/12/12/12	-
3	SRT	C	1466	-	-	10/12/12/12	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	1466	SRT	C2-C1	3.05	1.57	1.52
3	D	1466	SRT	C2-C1	2.44	1.56	1.52
3	C	1466	SRT	O1-C1	-2.29	1.23	1.30

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	1466	SRT	O2-C2-C3	-3.37	103.31	110.17
2	B	1467	GOL	O2-C2-C3	2.80	120.78	109.18
3	D	1466	SRT	C2-C3-C4	2.56	115.50	109.82
3	C	1466	SRT	C2-C3-C4	2.25	114.81	109.82

There are no chirality outliers.

All (27) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1468	GOL	C1-C2-C3-O3
2	B	1467	GOL	O1-C1-C2-C3
3	C	1466	SRT	O2-C2-C3-O3
3	D	1466	SRT	O2-C2-C3-O3
3	C	1466	SRT	O3-C3-C4-O41
3	D	1466	SRT	O3-C3-C4-O41
3	C	1466	SRT	C2-C3-C4-O4
3	C	1466	SRT	C2-C3-C4-O41
3	D	1466	SRT	C2-C3-C4-O4
3	D	1466	SRT	C2-C3-C4-O41
3	C	1466	SRT	C1-C2-C3-O3
3	D	1466	SRT	C1-C2-C3-O3
3	C	1466	SRT	O1-C1-C2-O2
3	C	1466	SRT	O11-C1-C2-O2
3	D	1466	SRT	O1-C1-C2-O2
3	D	1466	SRT	O11-C1-C2-O2
3	D	1466	SRT	O2-C2-C3-C4
3	C	1466	SRT	O3-C3-C4-O4
3	D	1466	SRT	O3-C3-C4-O4
2	A	1468	GOL	O1-C1-C2-C3
2	A	1468	GOL	O2-C2-C3-O3
2	B	1467	GOL	O1-C1-C2-O2
3	C	1466	SRT	O2-C2-C3-C4
2	B	1467	GOL	O2-C2-C3-O3
2	A	1468	GOL	O1-C1-C2-O2
3	D	1466	SRT	C1-C2-C3-C4
3	C	1466	SRT	C1-C2-C3-C4

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	450/495 (90%)	-0.12	9 (2%) 65 70	12, 22, 43, 72	2 (0%)
1	B	449/495 (90%)	-0.00	11 (2%) 59 65	12, 24, 47, 77	0
1	C	456/495 (92%)	0.19	31 (6%) 23 26	12, 23, 59, 92	1 (0%)
1	D	451/495 (91%)	0.25	30 (6%) 24 27	13, 24, 59, 97	1 (0%)
All	All	1806/1980 (91%)	0.08	81 (4%) 38 42	12, 23, 52, 97	4 (0%)

All (81) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	12	ILE	6.9
1	D	447	ASP	6.7
1	C	447	ASP	5.8
1	C	17	MET	5.6
1	B	466	ALA	5.3
1	D	10	TYR	5.2
1	C	16	THR	5.1
1	C	446	GLY	5.1
1	D	418	TYR	4.9
1	C	10	TYR	4.8
1	C	448	ARG	4.8
1	C	14	HIS	4.8
1	D	14	HIS	4.8
1	D	139	SER	4.7
1	D	448	ARG	4.5
1	C	20	VAL	4.4
1	A	314	LEU	4.4
1	A	467	GLU	4.4
1	C	19	GLU	3.8
1	D	11	ARG	3.8
1	D	20	VAL	3.7

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Mol	Chain	Res	Type	RSRZ
1	D	13	GLU	3.7
1	C	139	SER	3.6
1	C	18	GLY	3.5
1	B	315	GLN	3.4
1	D	22	VAL	3.3
1	C	415	ALA	3.3
1	B	314	LEU	3.2
1	C	12	ILE	3.2
1	C	15	ASP	3.2
1	A	123	GLY	3.1
1	B	12	ILE	3.1
1	C	453	ASP	3.0
1	B	9	ASN	3.0
1	B	124	GLY	3.0
1	C	22	VAL	2.8
1	D	431	GLU	2.8
1	C	422	ALA	2.8
1	A	325	VAL	2.8
1	D	451	ILE	2.8
1	A	324	LYS	2.7
1	D	446	GLY	2.7
1	D	24	ALA	2.7
1	D	125	VAL	2.6
1	D	419	GLU	2.6
1	D	429	LEU	2.6
1	C	128	HIS	2.6
1	D	449	LEU	2.6
1	A	10	TYR	2.6
1	C	451	ILE	2.5
1	C	26	ALA	2.5
1	C	34	ARG	2.5
1	C	24	ALA	2.5
1	C	418	TYR	2.4
1	B	123	GLY	2.4
1	D	450	SER	2.4
1	C	13	GLU	2.4
1	D	445	ILE	2.4
1	A	12	ILE	2.3
1	B	325	VAL	2.3
1	B	44	ARG	2.3
1	C	11	ARG	2.3
1	C	449	LEU	2.3

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Mol	Chain	Res	Type	RSRZ
1	D	454	LEU	2.3
1	C	416	ILE	2.3
1	C	432	ARG	2.3
1	D	432	ARG	2.3
1	D	128	HIS	2.3
1	D	23	PRO	2.2
1	D	126	THR	2.2
1	B	88	ASP	2.2
1	D	31	GLN	2.2
1	D	444	LEU	2.2
1	C	44	ARG	2.2
1	D	409	VAL	2.1
1	C	414	SER	2.1
1	A	448	ARG	2.1
1	D	34	ARG	2.1
1	B	122	LYS	2.1
1	A	447	ASP	2.0
1	C	409	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($\text{\AA}^2$)	Q<0.9
2	GOL	A	1468	6/6	0.89	0.12	36,38,38,38	0
2	GOL	B	1467	6/6	0.92	0.10	30,31,32,33	0
3	SRT	D	1466	10/10	0.95	0.08	23,25,30,39	0
3	SRT	C	1466	10/10	0.96	0.07	20,22,26,34	0

6.5 Other polymers [i](#)

There are no such residues in this entry.