



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:19 PM UTC

PDB ID : 4U3E / pdb_00004u3e
Title : Anaerobic ribonucleotide reductase
Authors : Funk, M.A.; Drennan, C.L.
Deposited on : 2014-07-20
Resolution : 1.64 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

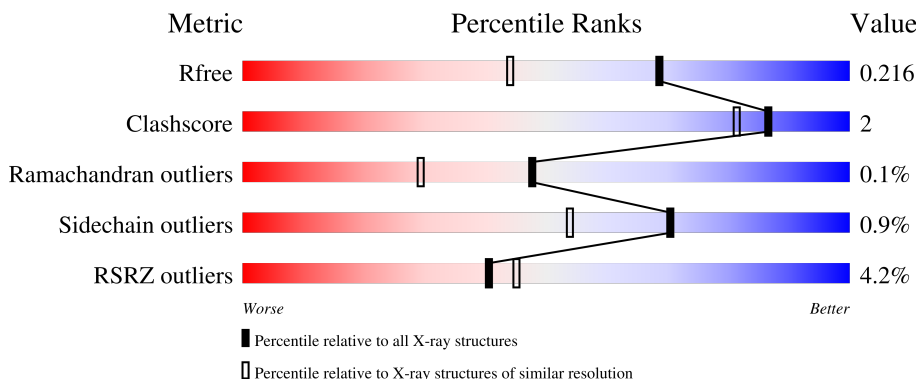
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.64 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	1141 (1.64-1.64)
Clashscore	190562	1171 (1.64-1.64)
Ramachandran outliers	187476	1151 (1.64-1.64)
Sidechain outliers	187428	1150 (1.64-1.64)
RSRZ outliers	180081	1141 (1.64-1.64)

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	671	4% 89% 6% 5%
1	B	671	3% 87% 5% 8%

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 11270 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 Ribonucleoside triphosphate reductase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	637	Total	C	N	O	S	Se	0	8	0
			5217	3362	866	962	5	22			
1	B	619	Total	C	N	O	S	Se	0	6	0
			5114	3302	846	941	4	21			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MSE	-	expression tag	UNP Q9WYL6
A	-18	GLY	-	expression tag	UNP Q9WYL6
A	-17	SER	-	expression tag	UNP Q9WYL6
A	-16	SER	-	expression tag	UNP Q9WYL6
A	-15	HIS	-	expression tag	UNP Q9WYL6
A	-14	HIS	-	expression tag	UNP Q9WYL6
A	-13	HIS	-	expression tag	UNP Q9WYL6
A	-12	HIS	-	expression tag	UNP Q9WYL6
A	-11	HIS	-	expression tag	UNP Q9WYL6
A	-10	HIS	-	expression tag	UNP Q9WYL6
A	-9	SER	-	expression tag	UNP Q9WYL6
A	-8	SER	-	expression tag	UNP Q9WYL6
A	-7	GLY	-	expression tag	UNP Q9WYL6
A	-6	LEU	-	expression tag	UNP Q9WYL6
A	-5	VAL	-	expression tag	UNP Q9WYL6
A	-4	PRO	-	expression tag	UNP Q9WYL6
A	-3	ARG	-	expression tag	UNP Q9WYL6
A	-2	GLY	-	expression tag	UNP Q9WYL6
A	-1	SER	-	expression tag	UNP Q9WYL6
A	0	HIS	-	expression tag	UNP Q9WYL6
B	-19	MSE	-	expression tag	UNP Q9WYL6
B	-18	GLY	-	expression tag	UNP Q9WYL6
B	-17	SER	-	expression tag	UNP Q9WYL6
B	-16	SER	-	expression tag	UNP Q9WYL6
B	-15	HIS	-	expression tag	UNP Q9WYL6

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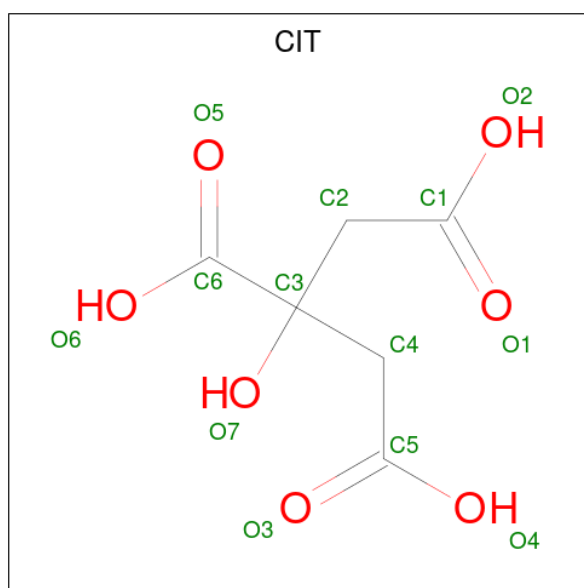
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Chain	Residue	Modelled	Actual	Comment	Reference
B	-14	HIS	-	expression tag	UNP Q9WYL6
B	-13	HIS	-	expression tag	UNP Q9WYL6
B	-12	HIS	-	expression tag	UNP Q9WYL6
B	-11	HIS	-	expression tag	UNP Q9WYL6
B	-10	HIS	-	expression tag	UNP Q9WYL6
B	-9	SER	-	expression tag	UNP Q9WYL6
B	-8	SER	-	expression tag	UNP Q9WYL6
B	-7	GLY	-	expression tag	UNP Q9WYL6
B	-6	LEU	-	expression tag	UNP Q9WYL6
B	-5	VAL	-	expression tag	UNP Q9WYL6
B	-4	PRO	-	expression tag	UNP Q9WYL6
B	-3	ARG	-	expression tag	UNP Q9WYL6
B	-2	GLY	-	expression tag	UNP Q9WYL6
B	-1	SER	-	expression tag	UNP Q9WYL6
B	0	HIS	-	expression tag	UNP Q9WYL6

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

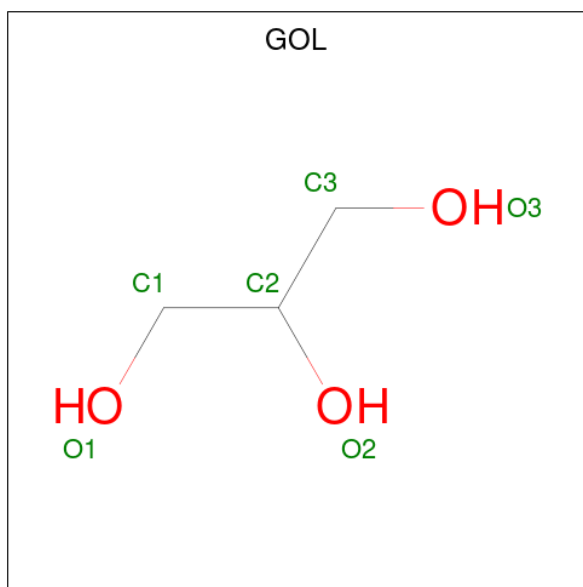
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Zn 1 1	0	0
2	B	1	Total Zn 1 1	0	0

- Molecule 3 is CITRIC ACID (CCD ID: CIT) (formula: C₆H₈O₇).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			13	6	7		
3	B	1	Total	C	O	0	0
			13	6	7		

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).

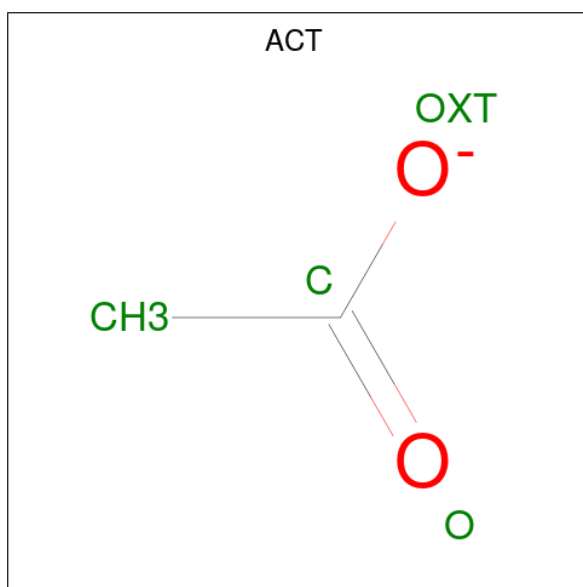


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 5 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Na	0	0
			1	1		

- Molecule 6 is ACETATE ION (CCD ID: ACT) (formula: $C_2H_3O_2$).



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

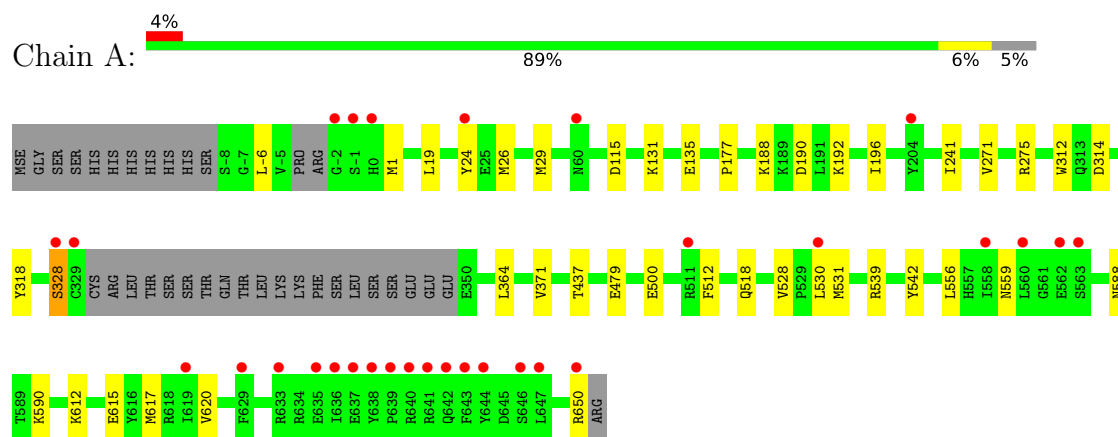
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	452	Total	O	0	1
			453	453		
7	B	440	Total	O	0	1
			441	441		

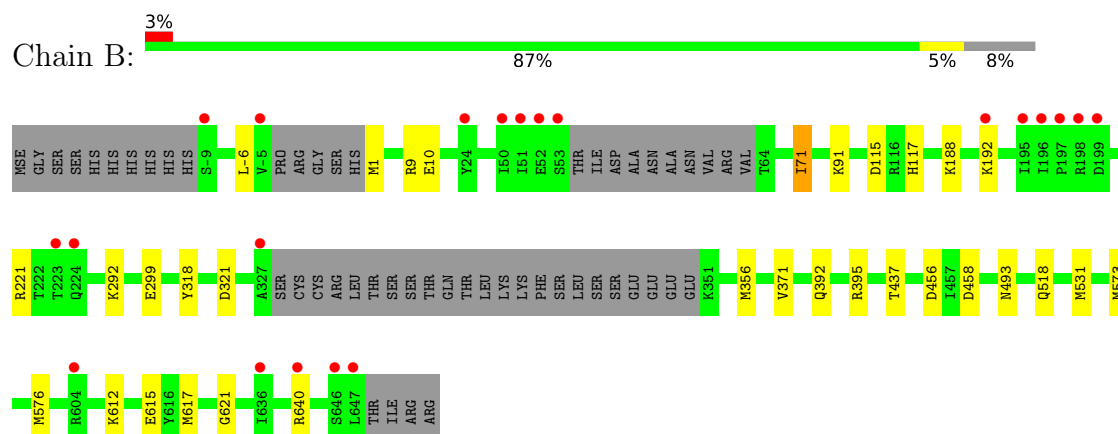
3 Residue-property plots [i](#)

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: Ribonucleoside triphosphate reductase



• Molecule 1: Ribonucleoside triphosphate reductase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	78.31Å 98.77Å 86.56Å 90.00° 111.72° 90.00°	Depositor
Resolution (Å)	49.38 – 1.64 49.38 – 1.64	Depositor EDS
% Data completeness (in resolution range)	95.1 (49.38-1.64) 89.0 (49.38-1.64)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.66 (at 1.64Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE: 1.9_1692)	Depositor
R, R_{free}	0.171 , 0.200 0.189 , 0.216	Depositor DCC
R_{free} test set	2000 reflections (1.36%)	wwPDB-VP
Wilson B-factor (Å ²)	15.6	Xtriage
Anisotropy	0.485	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 38.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	11270	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.40% 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: GOL, ACT, CIT, NA, ZN

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	0.36	0/5321	0.66	2/7150 (0.0%)
1	B	0.36	0/5217	0.67	0/7006
All	All	0.36	0/10538	0.67	2/14156 (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	556	LEU	CA-C-N	-5.36	114.48	122.74
1	A	556	LEU	C-N-CA	-5.36	114.48	122.74

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	5217	0	5129	25	0
1	B	5114	0	5030	21	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	13	0	4	0	0
3	B	13	0	5	0	0
4	A	6	0	8	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	6	0	8	0	0
5	A	1	0	0	0	0
6	A	4	0	3	0	0
7	A	453	0	0	3	0
7	B	441	0	0	0	0
All	All	11270	0	10187	43	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 2.

All (43) 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:612:LYS:NZ	1:A:615:GLU:OE2	2.13	0.80
1:A:650:ARG:HH22	1:B:9:ARG:H	1.30	0.78
1:B:188:LYS:HE3	1:B:192:LYS:NZ	2.09	0.68
1:A:188:LYS:HE2	1:A:192:LYS:HE3	1.78	0.65
1:A:650:ARG:HH22	1:B:9:ARG:N	1.94	0.65
1:B:392:GLN:HG2	1:B:395[B]:ARG:NH1	2.12	0.65
1:A:542:TYR:OH	7:A:983:HOH:O	2.13	0.64
1:A:559:ASN:HA	1:A:588:ASN:HB2	1.84	0.58
1:A:650:ARG:NH2	1:B:9:ARG:H	2.01	0.56
1:B:71:ILE:HD11	1:B:221:ARG:NH1	2.22	0.55
1:A:528:VAL:HG21	1:A:539:ARG:HG2	1.90	0.53
1:B:117[A]:HIS:CD2	1:B:117[A]:HIS:H	2.26	0.53
1:B:188:LYS:HE3	1:B:192:LYS:HZ2	1.74	0.53
1:B:10:GLU:OE1	1:B:91:LYS:HE3	2.11	0.51
1:A:590:LYS:HG2	1:A:617:MSE:SE	2.62	0.50
1:B:392:GLN:HG2	1:B:395[B]:ARG:HH12	1.74	0.49
1:A:559:ASN:ND2	7:A:1078:HOH:O	2.43	0.49
1:B:392:GLN:HA	1:B:395[B]:ARG:NH1	2.28	0.49
1:B:612:LYS:HE3	1:B:615:GLU:HG3	1.94	0.48
1:A:530:LEU:HG	1:A:531:MSE:HE2	1.95	0.48
1:A:500:GLU:HG2	1:A:620:VAL:HG11	1.95	0.48
1:A:612:LYS:HE3	1:A:615:GLU:HG3	1.97	0.47
1:A:177:PRO:HG2	1:A:364:LEU:HD13	1.96	0.46
1:B:-6:LEU:HG	1:B:1:MSE:HE3	1.99	0.45
1:A:-6:LEU:HD11	1:A:1:MSE:HG2	1.98	0.44
1:B:392:GLN:CG	1:B:395[B]:ARG:HH12	2.30	0.44
1:A:371:VAL:HG22	1:A:437:THR:HB	1.99	0.44
1:B:371:VAL:HG22	1:B:437:THR:HB	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:531:MSE:HE3	1:B:640:ARG:HB2	2.00	0.44
1:A:131:LYS:NZ	1:A:135:GLU:OE1	2.41	0.43
1:A:275:ARG:HD2	1:A:312:TRP:CD1	2.54	0.43
1:A:314:ASP:OD1	1:A:314:ASP:N	2.50	0.43
1:B:456:ASP:HB3	1:B:458:ASP:OD2	2.19	0.43
1:B:292:LYS:HE2	1:B:321:ASP:O	2.19	0.43
1:A:479:GLU:HG3	7:A:834:HOH:O	2.20	0.42
1:B:617:MSE:HB2	1:B:617:MSE:HE2	1.74	0.42
1:A:241:ILE:O	1:A:328:SER:HA	2.20	0.42
1:A:190:ASP:HB3	1:A:196:ILE:HG13	2.01	0.41
1:B:573:MSE:SE	1:B:576:MSE:HE3	2.70	0.41
1:A:29:MSE:HE1	1:A:512:PHE:CG	2.56	0.41
1:A:19:LEU:HD11	1:A:24:TYR:HA	2.03	0.41
1:A:26:MSE:HA	1:A:29:MSE:HE3	2.02	0.41
1:B:356[B]:MSE:SE	1:B:493:ASN:HD21	2.54	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	639/671 (95%)	628 (98%)	11 (2%)	0	100	100
1	B	617/671 (92%)	610 (99%)	6 (1%)	1 (0%)	43	27
All	All	1256/1342 (94%)	1238 (99%)	17 (1%)	1 (0%)	48	29

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	621	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	560/573 (98%)	555 (99%)	5 (1%)	70	54
1	B	550/573 (96%)	545 (99%)	5 (1%)	70	54
All	All	1110/1146 (97%)	1100 (99%)	10 (1%)	70	54

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

Mol	Chain	Res	Type
1	A	115	ASP
1	A	271	VAL
1	A	318	TYR
1	A	328	SER
1	A	518	GLN
1	B	71	ILE
1	B	115	ASP
1	B	299	GLU
1	B	318	TYR
1	B	518	GLN

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	229	ASN
1	A	526	GLN
1	B	151	HIS
1	B	256	HIS
1	B	313	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 8 ligands modelled in this entry, 3 are monoatomic - leaving 5 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	$\# Z > 2$
4	GOL	B	903	-	5,5,5	0.39	0	5,5,5	0.24	0
3	CIT	A	702	5	12,12,12	1.09	0	17,17,17	1.61	3 (17%)
3	CIT	B	902	-	12,12,12	1.08	0	17,17,17	1.34	1 (5%)
6	ACT	A	705	5	3,3,3	0.88	0	3,3,3	1.32	0
4	GOL	A	703	-	5,5,5	0.42	0	5,5,5	0.39	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
3	CIT	B	902	-	-	0/16/16/16	-
3	CIT	A	702	5	-	2/16/16/16	-
4	GOL	A	703	-	-	2/4/4/4	-
4	GOL	B	903	-	-	0/4/4/4	-

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	702	CIT	O6-C6-C3	4.24	121.27	113.14
3	B	902	CIT	O6-C6-C3	3.46	119.78	113.14
3	A	702	CIT	O2-C1-C2	2.65	122.76	114.35
3	A	702	CIT	O2-C1-O1	-2.07	118.01	123.33

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	703	GOL	C1-C2-C3-O3
4	A	703	GOL	O2-C2-C3-O3
3	A	702	CIT	O1-C1-C2-C3
3	A	702	CIT	O2-C1-C2-C3

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	617/671 (91%)	0.11	30 (4%) 35 40	8, 22, 42, 78	6 (0%)
1	B	599/671 (89%)	-0.02	21 (3%) 47 51	8, 21, 40, 86	5 (0%)
All	All	1216/1342 (90%)	0.05	51 (4%) 40 45	8, 21, 41, 86	11 (0%)

All (51) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	638	TYR	5.9
1	A	639	PRO	5.8
1	B	51	ILE	5.1
1	B	-9	SER	4.2
1	B	50	ILE	4.0
1	A	328	SER	3.9
1	A	0	HIS	3.9
1	A	640	ARG	3.7
1	A	60	ASN	3.7
1	B	647	LEU	3.7
1	A	643	PHE	3.5
1	A	641	ARG	3.4
1	A	-1	SER	3.3
1	B	197	PRO	3.2
1	A	636	ILE	3.2
1	A	329	CYS	3.2
1	B	53	SER	3.2
1	A	646	SER	3.1
1	A	637	GLU	3.1
1	A	530	LEU	3.0
1	A	647	LEU	3.0
1	A	629	PHE	3.0
1	A	644	TYR	2.9
1	B	52	GLU	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	650	ARG	2.8
1	B	198	ARG	2.7
1	A	562	GLU	2.7
1	A	-2	GLY	2.7
1	B	646	SER	2.6
1	B	640	ARG	2.6
1	B	192	LYS	2.6
1	B	196	ILE	2.6
1	B	24	TYR	2.5
1	A	619	ILE	2.5
1	A	511	ARG	2.5
1	A	558	ILE	2.5
1	B	199	ASP	2.4
1	A	560	LEU	2.4
1	B	327	ALA	2.4
1	A	642	GLN	2.4
1	B	-5	VAL	2.4
1	A	204	TYR	2.4
1	B	636	ILE	2.3
1	A	563	SER	2.3
1	A	24	TYR	2.3
1	A	633	ARG	2.2
1	B	224	GLN	2.2
1	B	223	THR	2.2
1	A	635	GLU	2.1
1	B	195	ILE	2.1
1	B	604	ARG	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
6	ACT	A	705	4/4	0.89	0.15	23,25,26,29	0
4	GOL	A	703	6/6	0.92	0.09	26,31,34,37	0
5	NA	A	704	1/1	0.93	0.09	31,31,31,31	0
4	GOL	B	903	6/6	0.93	0.08	27,32,34,37	0
3	CIT	B	902	13/13	0.94	0.09	19,26,51,53	0
3	CIT	A	702	13/13	0.95	0.07	18,21,30,31	0
2	ZN	B	901	1/1	0.97	0.05	28,28,28,28	0
2	ZN	A	701	1/1	0.98	0.04	27,27,27,27	0

6.5 Other polymers [i](#)

There are no such residues in this entry.