



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 01:18 am GMT

PDB ID : 2WCO
Title : STRUCTURES OF THE STREPTOMYCES COELICOLOR A3(2)
HYALURONAN LYASE IN COMPLEX WITH OLIGOSACCHARIDE
SUBSTRATES AND AN INHIBITOR
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Deposited on : 2009-03-12
Resolution : 1.94 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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : recalc28949

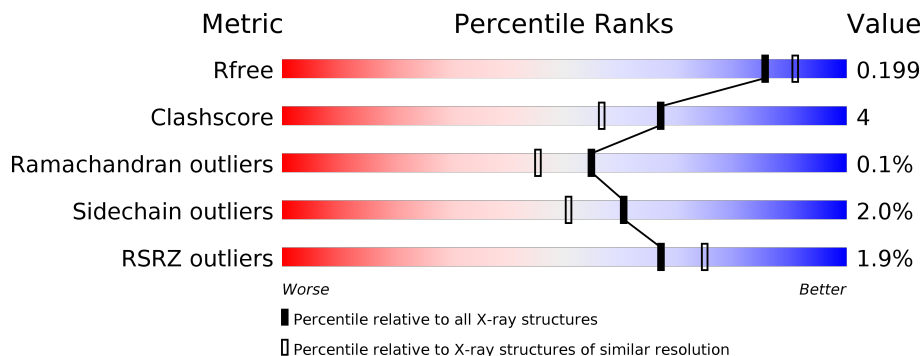
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.94 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}	100719	3233 (1.96-1.92)
Clashscore	112137	3430 (1.96-1.92)
Ramachandran outliers	110173	3395 (1.96-1.92)
Sidechain outliers	110143	3395 (1.96-1.92)
RSRZ outliers	101464	3250 (1.96-1.92)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	765	<div> <div>2%</div> <div>84%</div> <div>12%</div> <div>..</div> </div>

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	FMT	A	1767	-	-	X	-
3	FMT	A	1768	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6547 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 PUTATIVE SECRETED LYASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	742	5741	3612	1060	1050	19	0	12	0

There are 21 discrepancies between the modelled and reference sequences:

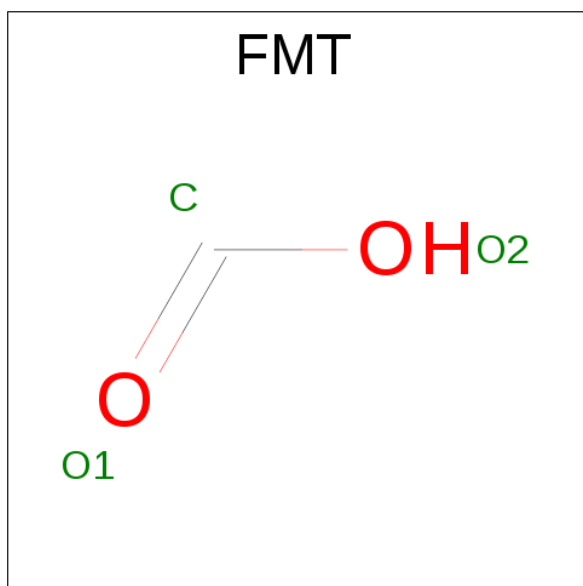
Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	EXPRESSION TAG	UNP O86516
A	2	GLY	-	EXPRESSION TAG	UNP O86516
A	3	SER	-	EXPRESSION TAG	UNP O86516
A	4	SER	-	EXPRESSION TAG	UNP O86516
A	5	HIS	-	EXPRESSION TAG	UNP O86516
A	6	HIS	-	EXPRESSION TAG	UNP O86516
A	7	HIS	-	EXPRESSION TAG	UNP O86516
A	8	HIS	-	EXPRESSION TAG	UNP O86516
A	9	HIS	-	EXPRESSION TAG	UNP O86516
A	10	HIS	-	EXPRESSION TAG	UNP O86516
A	11	SER	-	EXPRESSION TAG	UNP O86516
A	12	SER	-	EXPRESSION TAG	UNP O86516
A	13	GLY	-	EXPRESSION TAG	UNP O86516
A	14	LEU	-	EXPRESSION TAG	UNP O86516
A	15	VAL	-	EXPRESSION TAG	UNP O86516
A	16	PRO	-	EXPRESSION TAG	UNP O86516
A	17	ARG	-	EXPRESSION TAG	UNP O86516
A	18	GLY	-	EXPRESSION TAG	UNP O86516
A	19	SER	-	EXPRESSION TAG	UNP O86516
A	20	HIS	-	EXPRESSION TAG	UNP O86516
A	21	MET	-	EXPRESSION TAG	UNP O86516

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



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

- Molecule 3 is FORMIC ACID (three-letter code: FMT) (formula: CH_2O_2).



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

- Molecule 4 is a polymer of unknown type called SUGAR (2-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	2	Total	C	N	O	0	0
			26	14	1	11		

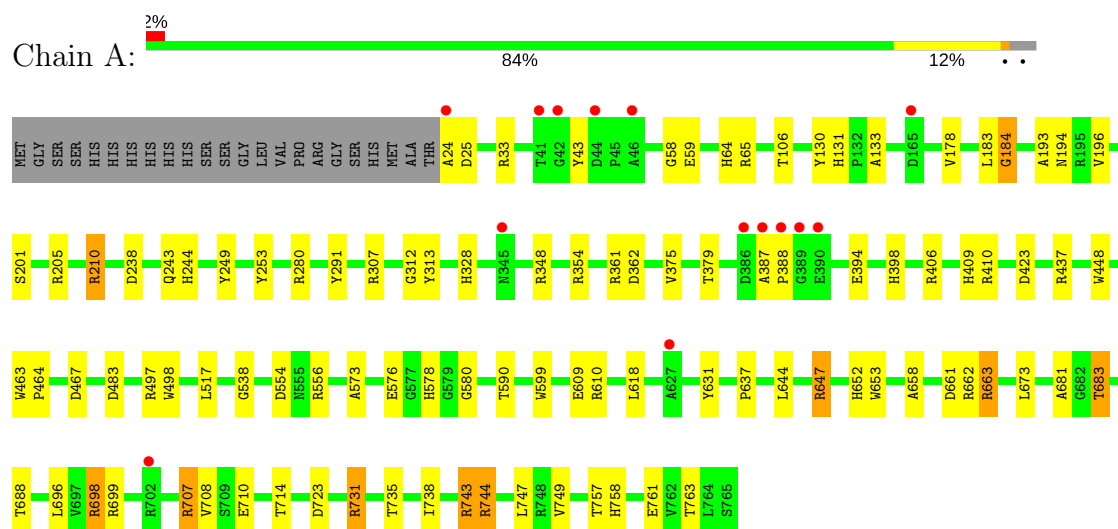
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	768	Total	O	0	0
			768	768		

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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: PUTATIVE SECRETED LYASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	140.11Å 140.11Å 100.65Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	100.00 – 1.94 19.73 – 1.94	Depositor EDS
% Data completeness (in resolution range)	89.4 (100.00-1.94) 89.5 (19.73-1.94)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.36 (at 1.93Å)	Xtriage
Refinement program	REFMAC 5.5.0082	Depositor
R, R_{free}	0.170 , 0.193 0.178 , 0.199	Depositor DCC
R_{free} test set	3368 reflections (5.32%)	DCC
Wilson B-factor (Å ²)	28.3	Xtriage
Anisotropy	0.029	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.42 , 53.5	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	6547	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.49% 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, FMT, NAG, GC4

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	1.22	15/5938 (0.3%)	1.06	26/8118 (0.3%)

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 (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	609	GLU	CB-CG	-6.57	1.39	1.52
1	A	576	GLU	CG-CD	6.52	1.61	1.51
1	A	253	TYR	CD1-CE1	6.39	1.49	1.39
1	A	681	ALA	CA-CB	5.93	1.65	1.52
1	A	59[A]	GLU	CG-CD	5.81	1.60	1.51
1	A	59[B]	GLU	CG-CD	5.81	1.60	1.51
1	A	763	THR	C-N	5.70	1.47	1.34
1	A	178	VAL	CB-CG2	5.61	1.64	1.52
1	A	743	ARG	CG-CD	5.49	1.65	1.51
1	A	313	TYR	CD1-CE1	-5.45	1.31	1.39
1	A	744	ARG	CG-CD	5.28	1.65	1.51
1	A	130	TYR	CD1-CE1	5.08	1.47	1.39
1	A	196	VAL	CB-CG2	5.04	1.63	1.52
1	A	291	TYR	CG-CD2	-5.01	1.32	1.39
1	A	201	SER	CB-OG	5.00	1.48	1.42

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	406	ARG	NE-CZ-NH1	10.80	125.70	120.30
1	A	354	ARG	NE-CZ-NH2	-9.98	115.31	120.30
1	A	348	ARG	NE-CZ-NH2	-8.84	115.88	120.30
1	A	33	ARG	NE-CZ-NH1	8.39	124.50	120.30
1	A	410	ARG	NE-CZ-NH1	8.35	124.47	120.30
1	A	610	ARG	NE-CZ-NH2	-8.29	116.16	120.30
1	A	307	ARG	NE-CZ-NH2	-8.16	116.22	120.30
1	A	184	GLY	N-CA-C	7.93	132.94	113.10
1	A	33	ARG	NE-CZ-NH2	-7.62	116.49	120.30
1	A	699	ARG	NE-CZ-NH1	7.30	123.95	120.30
1	A	610	ARG	NE-CZ-NH1	6.83	123.72	120.30
1	A	483	ASP	CB-CG-OD1	6.42	124.08	118.30
1	A	437	ARG	NE-CZ-NH2	-6.21	117.19	120.30
1	A	698	ARG	NE-CZ-NH1	6.04	123.32	120.30
1	A	707	ARG	NE-CZ-NH1	6.03	123.32	120.30
1	A	467	ASP	CB-CG-OD1	6.00	123.70	118.30
1	A	647	ARG	NE-CZ-NH1	5.90	123.25	120.30
1	A	65	ARG	NE-CZ-NH1	5.89	123.24	120.30
1	A	238	ASP	CB-CG-OD1	5.70	123.43	118.30
1	A	423	ASP	CB-CG-OD2	-5.35	113.48	118.30
1	A	210	ARG	NE-CZ-NH1	-5.29	117.65	120.30
1	A	354	ARG	CG-CD-NE	-5.10	101.09	111.80
1	A	362	ASP	CB-CG-OD2	5.09	122.88	118.30
1	A	280	ARG	NE-CZ-NH2	-5.08	117.76	120.30
1	A	661	ASP	CB-CG-OD1	5.02	122.82	118.30
1	A	65	ARG	NE-CZ-NH2	-5.02	117.79	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	183	LEU	Peptide

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	5741	0	5590	49	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	6	0	8	2	0
3	A	6	0	2	2	0
4	A	26	0	21	0	0
5	A	768	0	0	8	1
All	All	6547	0	5621	51	2

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 4.

All (51) 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:652[B]:HIS:HD2	5:A:2665:HOH:O	1.39	1.06
1:A:673[A]:LEU:HD11	1:A:696:LEU:HD11	1.49	0.92
1:A:556:ARG:HE	1:A:578:HIS:HD2	1.27	0.79
1:A:24:ALA:N	1:A:210:ARG:HH12	1.80	0.78
1:A:673[A]:LEU:HD11	1:A:696:LEU:CD1	2.13	0.77
2:A:1766:GOL:H12	3:A:1767:FMT:H	1.74	0.68
1:A:556:ARG:HE	1:A:578:HIS:CD2	2.15	0.60
1:A:731:ARG:HB2	1:A:731:ARG:NH1	2.17	0.60
2:A:1766:GOL:H12	3:A:1767:FMT:C	2.31	0.59
1:A:652[B]:HIS:CD2	5:A:2665:HOH:O	2.28	0.57
1:A:662:ARG:C	1:A:663[A]:ARG:HG3	2.25	0.57
1:A:497:ARG:HH22	1:A:714:THR:HG22	1.69	0.57
1:A:573:ALA:O	1:A:644[A]:LEU:HD21	2.04	0.56
1:A:580:GLY:HA3	1:A:644[A]:LEU:HD22	1.88	0.56
1:A:723:ASP:HB2	5:A:2737:HOH:O	2.05	0.56
1:A:590:THR:HG22	1:A:618[A]:LEU:HD22	1.89	0.54
1:A:64:HIS:HE1	5:A:2094:HOH:O	1.89	0.54
1:A:193:ALA:H	1:A:243:GLN:HE22	1.55	0.54
1:A:708:VAL:HG11	1:A:749:VAL:HG21	1.91	0.53
1:A:312:GLY:HA3	1:A:599:TRP:CD2	2.44	0.52
1:A:328:HIS:HE1	5:A:2349:HOH:O	1.94	0.51
1:A:463:TRP:N	1:A:464:PRO:HD2	2.26	0.51
1:A:43:TYR:HB3	1:A:379[B]:THR:CG2	2.43	0.49
1:A:398:HIS:HD2	1:A:409:HIS:ND1	2.11	0.49
1:A:731:ARG:HB2	1:A:731:ARG:CZ	2.42	0.48
1:A:375:VAL:O	1:A:379[B]:THR:HG23	2.13	0.48
1:A:387:ALA:HB1	1:A:388:PRO:HD2	1.96	0.48
1:A:538:GLY:HA3	1:A:631:TYR:CE2	2.49	0.47
1:A:448:TRP:CE3	1:A:637:PRO:HB3	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:243:GLN:HG3	1:A:244:HIS:H	1.80	0.46
1:A:743:ARG:NH2	5:A:2736:HOH:O	2.41	0.46
1:A:517:LEU:HD23	1:A:517:LEU:C	2.35	0.46
1:A:43:TYR:HB3	1:A:379[B]:THR:HG22	1.97	0.46
1:A:64:HIS:HD2	5:A:2052:HOH:O	1.98	0.45
1:A:647:ARG:HD2	1:A:653:TRP:CD2	2.51	0.45
1:A:731:ARG:HG3	1:A:761:GLU:HB3	1.97	0.45
1:A:707:ARG:HB3	1:A:757:THR:CG2	2.46	0.45
1:A:243:GLN:HG3	1:A:244:HIS:N	2.31	0.45
1:A:683[B]:THR:HG22	1:A:688:THR:HG23	1.98	0.44
1:A:58:GLY:HA3	1:A:106:THR:O	2.18	0.44
1:A:738:ILE:HD13	1:A:747:LEU:HD22	2.00	0.44
1:A:463:TRP:N	1:A:464:PRO:CD	2.81	0.43
1:A:647:ARG:HD2	1:A:653:TRP:CE2	2.54	0.43
1:A:735:THR:OG1	1:A:758:HIS:CD2	2.73	0.42
1:A:663[A]:ARG:HG2	5:A:2691:HOH:O	2.19	0.42
1:A:658:ALA:CB	1:A:663[B]:ARG:HG3	2.49	0.42
1:A:361:ARG:NH2	1:A:394:GLU:HG3	2.35	0.41
1:A:131:HIS:CE1	1:A:133:ALA:HB3	2.54	0.41
1:A:24:ALA:CA	1:A:210:ARG:HH22	2.34	0.41

All (2) 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)
5:A:2665:HOH:O	5:A:2665:HOH:O[8_555]	1.73	0.47
1:A:194:ASN:ND2	1:A:743:ARG:NH2[4_555]	1.98	0.22

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	752/765 (98%)	732 (97%)	19 (2%)	1 (0%)	55 45

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	184	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	574/582 (99%)	561 (98%)	13 (2%)	56 45

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

Mol	Chain	Res	Type
1	A	25	ASP
1	A	205	ARG
1	A	249	TYR
1	A	498	TRP
1	A	554	ASP
1	A	663[A]	ARG
1	A	663[B]	ARG
1	A	683[A]	THR
1	A	683[B]	THR
1	A	698	ARG
1	A	710	GLU
1	A	731	ARG
1	A	744	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	64	HIS
1	A	243	GLN

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Mol	Chain	Res	Type
1	A	328	HIS
1	A	398	HIS
1	A	427	HIS
1	A	571	HIS
1	A	578	HIS
1	A	758	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

2 carbohydrates 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	$\# Z > 2$
4	GC4	A	1769	4	8,11,12	0.83	0	9,15,17	2.00	3 (33%)
4	NAG	A	1770	4	15,15,15	1.06	0	21,21,21	1.76	3 (14%)

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
4	GC4	A	1769	4	-	0/0/17/20	0/1/1/1
4	NAG	A	1770	4	-	0/6/26/26	0/1/1/1

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1769	GC4	O5-C5-C4	2.41	112.41	109.60
4	A	1770	NAG	O3-C3-C2	2.49	114.66	109.61
4	A	1770	NAG	O5-C1-C2	3.28	112.81	109.52
4	A	1769	GC4	C3-C4-C5	3.39	116.11	109.75
4	A	1769	GC4	C4-C3-C2	3.64	115.02	110.34
4	A	1770	NAG	C1-C2-N2	5.07	116.60	110.73

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.6 Ligand geometry

3 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	1766	-	5,5,5	0.32	0	5,5,5	0.52	0
3	FMT	A	1767	-	0,2,2	0.00	-	0,1,1	0.00	-
3	FMT	A	1768	-	0,2,2	0.00	-	0,1,1	0.00	-

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	1766	-	-	0/4/4/4	0/0/0/0
3	FMT	A	1767	-	-	0/0/0/0	0/0/0/0
3	FMT	A	1768	-	-	0/0/0/0	0/0/0/0

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.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1766	GOL	2	0
3	A	1767	FMT	2	0

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	742/765 (96%)	-0.06	14 (1%) 67 75	21, 29, 43, 61	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	24	ALA	6.4
1	A	388	PRO	5.5
1	A	390	GLU	4.7
1	A	42	GLY	3.6
1	A	702	ARG	2.9
1	A	387	ALA	2.8
1	A	165	ASP	2.6
1	A	41	THR	2.6
1	A	345	ASN	2.6
1	A	389	GLY	2.3
1	A	44	ASP	2.2
1	A	46	ALA	2.1
1	A	386	ASP	2.1
1	A	627	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](#)

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(\AA^2)	Q<0.9
4	GC4	A	1769	11/12	0.97	0.10	0.74	28,32,39,39	0
4	NAG	A	1770	15/15	0.98	0.07	-0.46	23,27,31,33	0

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(\AA^2)	Q<0.9
3	FMT	A	1768	3/3	0.66	0.34	3.57	67,67,68,70	0
2	GOL	A	1766	6/6	0.97	0.11	0.11	29,34,35,36	0
3	FMT	A	1767	3/3	0.98	0.11	-0.08	37,37,37,37	0

6.5 Other polymers [i](#)

There are no such residues in this entry.