



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2014 – 03:04 AM GMT

PDB ID : 1OJP
Title : SPECIFICITY AND MECHANISM OF STREPTOCOCCUS PNEUMONIAE HYALURONATE LYASE: COMPLEX WITH 6-SULPHATED CHONDROITIN DISACCHARIDE
Authors : Rigden, D.J.; Jedrzejewski, M.J.
Deposited on : 2003-07-11
Resolution : 1.90 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

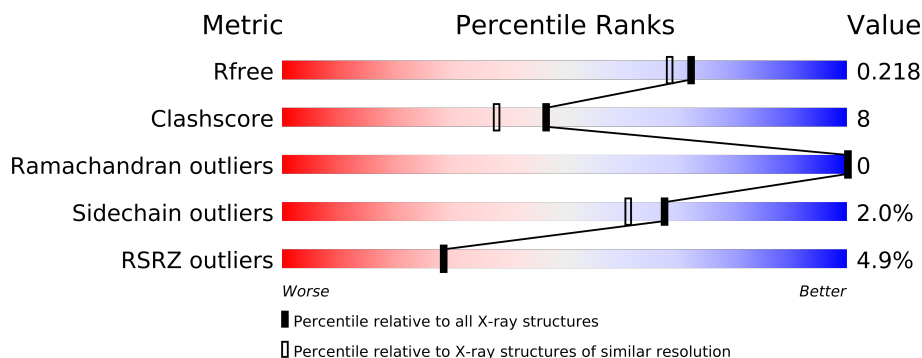
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 1.90 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}	66092	3684 (1.90-1.90)
Clashscore	79885	4465 (1.90-1.90)
Ramachandran outliers	78287	4413 (1.90-1.90)
Sidechain outliers	78261	4414 (1.90-1.90)
RSRZ outliers	66119	3686 (1.90-1.90)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	731	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6559 atoms, of which 0 are hydrogen and 0 are deuterium.

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 HYALURONATE LYASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	723	Total	C	N	O	S	0	12	1
			5905	3716	991	1176	22			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	893	HIS	-	EXPRESSION TAG	UNP Q54873
A	894	HIS	-	EXPRESSION TAG	UNP Q54873
A	895	HIS	-	EXPRESSION TAG	UNP Q54873
A	896	HIS	-	EXPRESSION TAG	UNP Q54873
A	897	HIS	-	EXPRESSION TAG	UNP Q54873
A	898	HIS	-	EXPRESSION TAG	UNP Q54873
A	173	THR	ALA	CONFLICT	UNP Q54873
A	196	ASP	GLU	CONFLICT	UNP Q54873
A	223	ILE	THR	CONFLICT	UNP Q54873
A	496	ARG	CYS	CONFLICT	UNP Q54873
A	541	THR	PRO	CONFLICT	UNP Q54873
A	704	SER	GLY	CONFLICT	UNP Q54873
A	736	SER	PHE	CONFLICT	UNP Q54873
A	790	GLY	ARG	CONFLICT	UNP Q54873

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	2	Total	C	N	O	S	0	0
			30	14	1	14	1		

There are 14 discrepancies between the modelled and reference sequences:

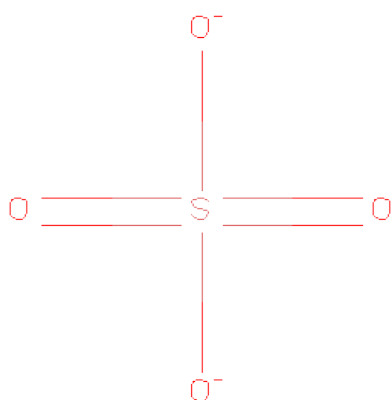
Chain	Residue	Modelled	Actual	Comment	Reference
A	893	HIS	-	EXPRESSION TAG	UNP Q54873
A	894	HIS	-	EXPRESSION TAG	UNP Q54873

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Chain	Residue	Modelled	Actual	Comment	Reference
A	895	HIS	-	EXPRESSION TAG	UNP Q54873
A	896	HIS	-	EXPRESSION TAG	UNP Q54873
A	897	HIS	-	EXPRESSION TAG	UNP Q54873
A	898	HIS	-	EXPRESSION TAG	UNP Q54873
A	173	THR	ALA	CONFLICT	UNP Q54873
A	196	ASP	GLU	CONFLICT	UNP Q54873
A	223	ILE	THR	CONFLICT	UNP Q54873
A	496	ARG	CYS	CONFLICT	UNP Q54873
A	541	THR	PRO	CONFLICT	UNP Q54873
A	704	SER	GLY	CONFLICT	UNP Q54873
A	736	SER	PHE	CONFLICT	UNP Q54873
A	790	GLY	ARG	CONFLICT	UNP Q54873

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	604	Total 604	O 604	0	0

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: HYALURONATE LYASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	83.70Å 103.01Å 101.16Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	64.49 – 1.90 64.49 – 1.90	Depositor EDS
% Data completeness (in resolution range)	91.9 (64.49-1.90) 92.0 (64.49-1.90)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.27 (at 1.90Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.196 , 0.221 0.193 , 0.218	Depositor DCC
R_{free} test set	3179 reflections (5.24%)	DCC
Wilson B-factor (Å ²)	22.5	Xtriage
Anisotropy	0.689	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 44.5	EDS
Estimated twinning fraction	0.024 for -h,l,k	Xtriage
L-test for twinning	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 63868 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6559	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 4.05% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NG6, GCD, SO4

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	3.72	8/6026 (0.1%)	1.61	12/8135 (0.1%)

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	2

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	203[A]	GLN	CD-NE2	131.87	4.62	1.32
1	A	203[B]	GLN	CD-NE2	131.87	4.62	1.32
1	A	862[A]	ARG	NE-CZ	104.07	2.68	1.33
1	A	862[B]	ARG	NE-CZ	104.07	2.68	1.33
1	A	319[A]	LYS	CD-CE	82.69	3.58	1.51
1	A	319[B]	LYS	CD-CE	82.69	3.58	1.51
1	A	319[A]	LYS	CE-NZ	79.62	3.48	1.49
1	A	319[B]	LYS	CE-NZ	79.62	3.48	1.49

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	862[A]	ARG	NE-CZ-NH1	-62.69	88.96	120.30
1	A	862[B]	ARG	NE-CZ-NH1	-62.69	88.96	120.30
1	A	862[A]	ARG	CD-NE-CZ	-45.68	59.65	123.60
1	A	862[B]	ARG	CD-NE-CZ	-45.68	59.65	123.60
1	A	319[A]	LYS	CD-CE-NZ	-30.58	41.37	111.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	319[B]	LYS	CD-CE-NZ	-30.58	41.37	111.70
1	A	862[A]	ARG	NE-CZ-NH2	29.24	134.92	120.30
1	A	862[B]	ARG	NE-CZ-NH2	29.24	134.92	120.30
1	A	203[A]	GLN	CG-CD-NE2	-25.87	54.61	116.70
1	A	203[B]	GLN	CG-CD-NE2	-25.87	54.61	116.70
1	A	203[A]	GLN	OE1-CD-NE2	-23.82	67.12	121.90
1	A	203[B]	GLN	OE1-CD-NE2	-23.82	67.12	121.90

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	203[A]	GLN	Sidechain
1	A	862[A]	ARG	Sidechain

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5905	0	5727	95	0
2	A	30	0	19	1	0
3	A	20	0	0	0	0
4	A	604	0	0	6	0
All	All	6559	0	5746	95	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 8.

All (95) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:862[A]:ARG:HD3	1:A:862[A]:ARG:CZ	1.60	1.28
1:A:862[A]:ARG:CZ	1:A:862[A]:ARG:CD	2.31	1.07
1:A:319[B]:LYS:NZ	1:A:319[B]:LYS:HD3	1.74	1.02
1:A:613:ASN:H	1:A:698:GLN:HE22	1.08	0.92
1:A:319[B]:LYS:CD	1:A:319[B]:LYS:CE	2.50	0.90

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:862[B]:ARG:HD2	1:A:862[B]:ARG:CZ	2.06	0.85
1:A:319[A]:LYS:HD2	1:A:319[A]:LYS:NZ	1.97	0.80
1:A:862[B]:ARG:NE	1:A:862[B]:ARG:CZ	2.47	0.77
1:A:282:ASN:ND2	1:A:285:LYS:HG2	1.99	0.77
1:A:862[B]:ARG:NE	1:A:862[B]:ARG:HH11	1.81	0.76
1:A:319[A]:LYS:CD	1:A:319[A]:LYS:NZ	2.49	0.76
1:A:708:LYS:HE3	1:A:710:TYR:OH	1.85	0.76
1:A:862[B]:ARG:NE	1:A:862[B]:ARG:NH1	2.36	0.73
1:A:319[B]:LYS:NZ	1:A:319[B]:LYS:CD	2.49	0.73
1:A:862[B]:ARG:CD	1:A:862[B]:ARG:CZ	2.66	0.72
1:A:319[B]:LYS:HZ2	1:A:319[B]:LYS:HD3	1.55	0.70
1:A:315:ASP:O	1:A:319[A]:LYS:HG3	1.93	0.69
1:A:704[A]:SER:C	1:A:705:ASN:HD22	1.97	0.68
1:A:704[B]:SER:C	1:A:705:ASN:HD22	1.98	0.66
1:A:839:LYS:HD2	1:A:853:VAL:HG23	1.76	0.65
1:A:248:MET:O	1:A:252[B]:VAL:HG23	1.97	0.64
1:A:529:LEU:HD23	1:A:529:LEU:C	2.18	0.64
1:A:319[A]:LYS:HD2	1:A:319[A]:LYS:HZ2	1.60	0.62
1:A:650:ASN:HD21	1:A:832:GLN:HE22	1.47	0.60
1:A:758[B]:LEU:HD23	1:A:778:GLU:HB2	1.85	0.59
1:A:422:VAL:HA	4:A:2009:HOH:O	2.03	0.59
1:A:874:ASN:HD22	1:A:874:ASN:C	2.06	0.59
1:A:254:ASN:HD22	1:A:254:ASN:C	2.07	0.58
1:A:319[B]:LYS:CE	1:A:319[B]:LYS:CG	2.82	0.57
1:A:862[A]:ARG:NE	1:A:862[A]:ARG:CZ	2.68	0.57
1:A:282:ASN:HD22	1:A:282:ASN:C	2.09	0.55
1:A:282:ASN:HD21	1:A:285:LYS:HG2	1.70	0.55
1:A:613:ASN:N	1:A:698:GLN:HE22	1.92	0.54
1:A:458:MET:HE2	1:A:567:SER:HB2	1.89	0.54
1:A:254:ASN:ND2	1:A:256:SER:H	2.05	0.53
1:A:808[B]:GLN:HG3	4:A:2536:HOH:O	2.08	0.53
1:A:664:LEU:C	1:A:664:LEU:HD23	2.28	0.53
1:A:354:GLY:HA3	1:A:377:ILE:HD11	1.90	0.53
1:A:286:SER:O	1:A:288:VAL:HG23	2.10	0.52
1:A:298:THR:HB	1:A:299:PRO:HD3	1.92	0.52
1:A:374:ILE:HD13	1:A:423:ILE:HG23	1.93	0.51
1:A:822:GLU:CD	1:A:822:GLU:H	2.14	0.51
1:A:581:LYS:HB3	1:A:768:GLU:HB2	1.93	0.50
1:A:248:MET:O	1:A:252[A]:VAL:HG13	2.11	0.50
1:A:252[A]:VAL:HG23	1:A:253:THR:HG23	1.94	0.50
1:A:223:ILE:HG13	1:A:224:TYR:CD2	2.47	0.50
1:A:584:LYS:NZ	1:A:768:GLU:HG2	2.27	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:243:ARG:O	1:A:247:GLU:HG3	2.11	0.49
1:A:327:LYS:HD3	1:A:327:LYS:C	2.32	0.49
1:A:705:ASN:N	1:A:705:ASN:HD22	2.08	0.49
1:A:502:LYS:HB2	1:A:529:LEU:HD21	1.94	0.49
1:A:170:VAL:HG23	1:A:315:ASP:OD2	2.13	0.49
1:A:292:TRP:CD2	2:A:901:NG6:H5	2.46	0.49
1:A:417:SER:HB2	1:A:484:ARG:HB2	1.95	0.49
1:A:862[A]:ARG:NH1	1:A:862[A]:ARG:CD	2.74	0.48
1:A:664:LEU:HD23	1:A:665:THR:N	2.29	0.48
1:A:274:GLU:OE1	1:A:274:GLU:HA	2.14	0.47
1:A:404:TYR:CE1	1:A:461:GLY:HA3	2.49	0.47
1:A:202:ASN:ND2	1:A:251:GLN:HE22	2.12	0.47
1:A:331:ASP:OD2	1:A:334:HIS:ND1	2.43	0.47
1:A:870:ILE:HD11	1:A:887:PHE:CD2	2.50	0.46
1:A:203[A]:GLN:HG2	4:A:2031:HOH:O	2.14	0.46
1:A:314:SER:OG	1:A:317:GLU:HG3	2.16	0.46
1:A:744[B]:TYR:CE2	1:A:797:LEU:HD13	2.51	0.46
1:A:502:LYS:CB	1:A:529:LEU:HD21	2.46	0.46
1:A:254:ASN:ND2	1:A:254:ASN:C	2.69	0.46
1:A:374:ILE:O	1:A:378:GLU:HG3	2.16	0.46
1:A:839:LYS:HD2	1:A:853:VAL:CG2	2.45	0.45
1:A:377:ILE:O	1:A:380:VAL:HG12	2.17	0.45
1:A:483:HIS:NE2	1:A:529:LEU:HD12	2.31	0.45
1:A:462:ARG:HD2	1:A:582:GLU:HB2	1.99	0.45
1:A:640:PHE:CD1	1:A:875:PRO:HG2	2.52	0.45
1:A:820:ASN:HD22	1:A:825:GLN:HG2	1.83	0.44
1:A:458:MET:HE1	1:A:566:LEU:C	2.37	0.44
1:A:671:PHE:HB2	1:A:678:ALA:HB3	1.99	0.43
1:A:845:SER:HB2	1:A:853:VAL:HG23	2.01	0.43
1:A:869:LYS:NZ	4:A:2568:HOH:O	2.51	0.43
1:A:667:HIS:HD2	4:A:2546:HOH:O	2.02	0.43
1:A:570:SER:HA	1:A:636:LEU:HB3	2.01	0.43
1:A:529:LEU:HD23	1:A:530:LEU:N	2.34	0.43
1:A:874:ASN:ND2	1:A:874:ASN:C	2.72	0.42
1:A:495:GLN:NE2	4:A:2258:HOH:O	2.52	0.42
1:A:354:GLY:O	1:A:358:VAL:HB	2.20	0.42
1:A:208:LYS:HD3	1:A:247:GLU:OE2	2.19	0.42
1:A:882:PRO:HB2	1:A:884:GLN:NE2	2.34	0.42
1:A:407:ALA:O	1:A:410:ASN:HB2	2.20	0.42
1:A:404:TYR:CD1	1:A:407:ALA:HB3	2.54	0.42
1:A:581:LYS:HD3	1:A:768:GLU:HG3	2.01	0.41
1:A:331:ASP:HA	1:A:332:PRO:HD2	1.97	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:882:PRO:HB2	1:A:884:GLN:HE22	1.85	0.41
1:A:661:ASN:H	1:A:661:ASN:HD22	1.67	0.41
1:A:661:ASN:H	1:A:661:ASN:ND2	2.19	0.41
1:A:669:SER:OG	1:A:825:GLN:NE2	2.54	0.40
1:A:492:GLU:OE2	1:A:496:ARG:HD3	2.21	0.40
1:A:788:GLN:HE21	1:A:788:GLN:HA	1.87	0.40

There are no symmetry-related clashes.

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	733/731 (100%)	704 (96%)	29 (4%)	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	653/649 (101%)	639 (98%)	14 (2%)	66	59

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	254	ASN
1	A	263	GLU
1	A	282	ASN
1	A	319[A]	LYS

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Mol	Chain	Res	Type
1	A	319[B]	LYS
1	A	368	GLN
1	A	404	TYR
1	A	433	ASP
1	A	529	LEU
1	A	661	ASN
1	A	705	ASN
1	A	773	LYS
1	A	822	GLU
1	A	874	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	202	ASN
1	A	231	ASN
1	A	237	ASN
1	A	254	ASN
1	A	261	GLN
1	A	277	HIS
1	A	282	ASN
1	A	368	GLN
1	A	386	GLN
1	A	418	GLN
1	A	661	ASN
1	A	667	HIS
1	A	683	ASN
1	A	698	GLN
1	A	788	GLN
1	A	820	ASN
1	A	825	GLN
1	A	832	GLN
1	A	874	ASN

5.3.3 RNA ⓘ

There are no RNA chains 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	$\# Z > 2$
2	GCD	A	900	2	10,11,12	5.32	4 (40%)	11,15,17	3.53	9 (81%)
2	NG6	A	901	2	19,19,19	2.35	8 (42%)	28,28,28	4.71	4 (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
2	GCD	A	900	2	-	0/4/17/20	0/1/1/1
2	NG6	A	901	2	-	0/10/30/30	0/1/1/1

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	900	GCD	O5-C5	15.60	1.54	1.37
2	A	901	NG6	O5-C1	4.77	1.52	1.43
2	A	901	NG6	C2-N	4.57	1.53	1.45
2	A	900	GCD	C3-C4	4.47	1.56	1.50
2	A	901	NG6	O5-C5	4.04	1.54	1.44
2	A	901	NG6	C4-C3	2.99	1.60	1.52
2	A	901	NG6	O6-S	-2.72	1.48	1.57
2	A	901	NG6	C7-N	2.68	1.45	1.34
2	A	900	GCD	C4-C5	2.56	1.38	1.33
2	A	901	NG6	O7-C7	2.34	1.28	1.23
2	A	901	NG6	C1-C2	2.26	1.55	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	900	GCD	O6A-C6	2.02	1.28	1.22

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	901	NG6	C6-O6-S	23.29	146.71	116.76
2	A	901	NG6	O6-C6-C5	7.09	121.30	107.81
2	A	900	GCD	O3-C3-C2	-6.48	102.48	111.03
2	A	900	GCD	C2-C3-C4	4.45	119.86	111.68
2	A	900	GCD	O3-C3-C4	4.13	119.47	109.67
2	A	900	GCD	O6A-C6-C5	-3.87	111.03	120.19
2	A	900	GCD	O5-C5-C4	-3.66	118.30	122.91
2	A	900	GCD	O6B-C6-O6A	2.91	130.52	123.62
2	A	900	GCD	O5-C5-C6	2.69	116.00	111.75
2	A	900	GCD	C3-C4-C5	-2.58	117.45	121.82
2	A	901	NG6	O4-C4-C5	2.27	115.25	109.28
2	A	901	NG6	O4-C4-C3	-2.16	105.52	110.35
2	A	900	GCD	O2-C2-C3	-2.06	108.32	111.03

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.6 Ligand geometry ⓘ

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	$\# Z > 2$
3	SO4	A	1200	-	4,4,4	0.34	0	6,6,6	0.07	0
3	SO4	A	1201	-	4,4,4	0.21	0	6,6,6	0.08	0
3	SO4	A	1202	-	4,4,4	0.36	0	6,6,6	0.10	0
3	SO4	A	1203	-	4,4,4	0.39	0	6,6,6	0.07	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	SO4	A	1200	-	-	0/0/0/0	0/0/0/0
3	SO4	A	1201	-	-	0/0/0/0	0/0/0/0
3	SO4	A	1202	-	-	0/0/0/0	0/0/0/0
3	SO4	A	1203	-	-	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.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

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	723/731 (98%)	0.22	36 (4%)	28 28	14, 25, 47, 64	0

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	892	GLN	7.7
1	A	221	ASP	6.3
1	A	427	LYS	4.9
1	A	224	TYR	4.2
1	A	340	ASP	3.9
1	A	891	GLU	3.7
1	A	171	LYS	3.6
1	A	170	VAL	3.3
1	A	223	ILE	3.3
1	A	230	SER	3.3
1	A	232	TYR	3.2
1	A	375	ARG	3.2
1	A	174	TYR	3.1
1	A	368	GLN	2.9
1	A	491	GLY	2.8
1	A	704[A]	SER	2.7
1	A	335	PHE	2.7
1	A	428	ASN	2.6
1	A	289	GLY	2.6
1	A	338	THR	2.5
1	A	234	THR	2.4
1	A	425	LYS	2.4
1	A	370	ILE	2.4
1	A	173	THR	2.4
1	A	290	ASN	2.2
1	A	433	ASP	2.2
1	A	219	GLN	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	430	ILE	2.1
1	A	367	ASP	2.1
1	A	233	LYS	2.1
1	A	279	HIS	2.1
1	A	432[A]	LYS	2.1
1	A	228[A]	LYS	2.0
1	A	316	GLU	2.0
1	A	392	GLN	2.0
1	A	435	MET	2.0

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

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

6.3 Carbohydrates ⓘ

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	GCD	A	900	11/12	0.32	9.55	46,48,49,50	0
2	NG6	A	901	19/19	0.31	4.02	35,48,67,68	0

6.4 Ligands ⓘ

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	SO4	A	1201	5/5	0.10	0.72	45,45,48,49	0
3	SO4	A	1200	5/5	0.12	0.02	51,52,53,54	0
3	SO4	A	1202	5/5	0.12	-0.49	61,61,62,62	0
3	SO4	A	1203	5/5	0.17	-0.78	70,70,70,71	0

6.5 Other polymers ⓘ

There are no such residues in this entry.