



Full wwPDB X-ray Structure Validation Report

Feb 26, 2014 – 09:41 PM GMT

PDB ID : 1LXK
Title : Streptococcus pneumoniae Hyaluronate Lyase in Complex with Tetrasaccharide Hyaluronan Substrate
Authors : Jedrzejewski, M.J.; Mello, L.V.; De Groot, B.L.; Li, S.
Deposited on : 2002-06-05
Resolution : 1.53 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>

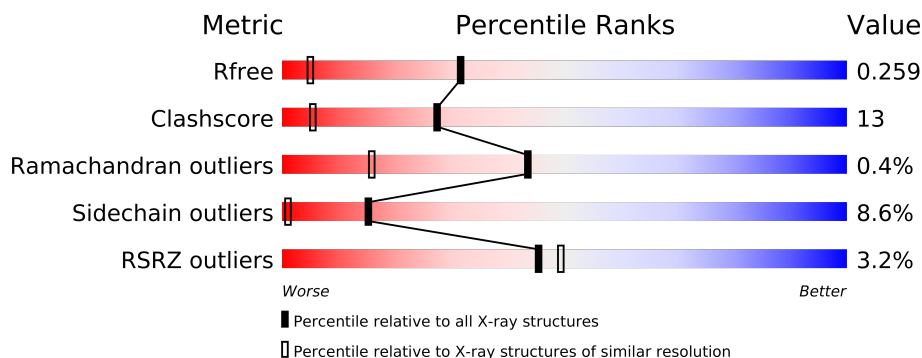
The following versions of software and data (see [references](#)) were used in the production of this report:

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 1.53 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	1031 (1.56-1.52)
Clashscore	79885	1155 (1.56-1.52)
Ramachandran outliers	78287	1127 (1.56-1.52)
Sidechain outliers	78261	1125 (1.56-1.52)
RSRZ outliers	66119	1031 (1.56-1.52)

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	721	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6512 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	719	Total	C	N	O	S	0	0	0
			5774	3632	966	1154	22			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	408	PHE	TYR	ENGINEERED	GB 437705
A	731	VAL	GLY	SEE REMARK 999	GB 437705

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

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	4	Total	C	N	O	0	0
			53	28	2	23		

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	408	PHE	TYR	ENGINEERED	GB 437705
A	731	VAL	GLY	SEE REMARK 999	GB 437705

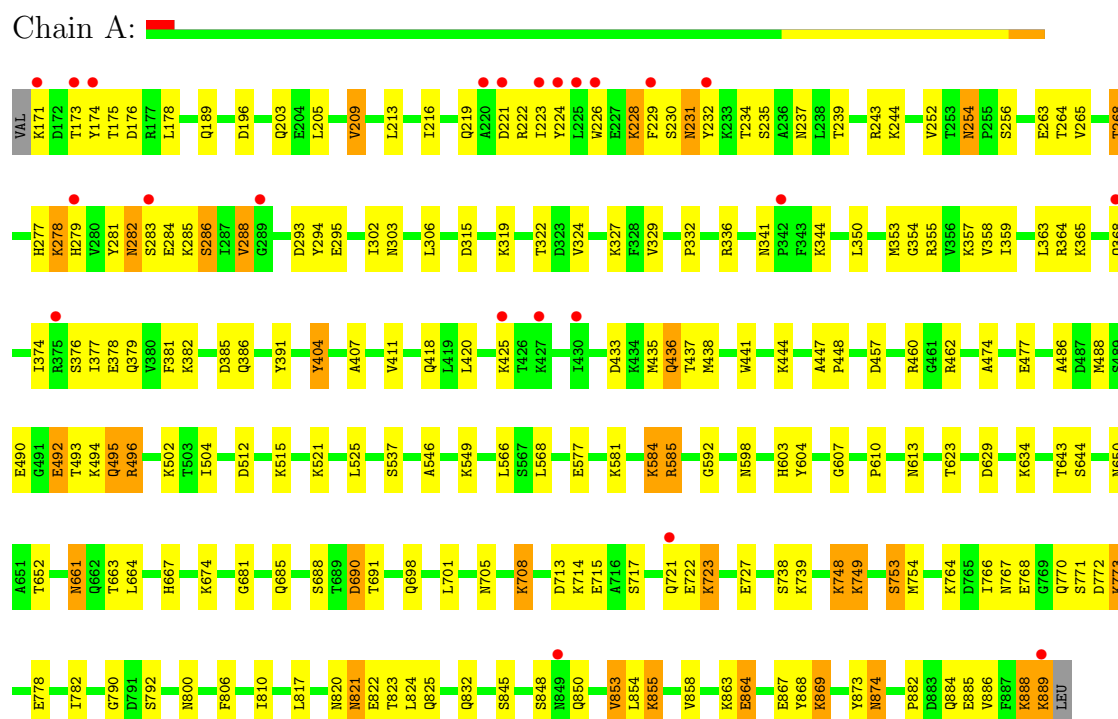
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	685	Total	O	0	0
			685	685		

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	83.74Å 103.82Å 101.35Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 1.53 64.55 – 1.53	Depositor EDS
% Data completeness (in resolution range)	(Not available) (20.00-1.53) 93.2 (64.55-1.53)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.08 (at 1.52Å)	Xtriage
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.190 , 0.239 0.226 , 0.259	Depositor DCC
R_{free} test set	1222 reflections (0.98%)	DCC
Wilson B-factor (Å ²)	15.9	Xtriage
Anisotropy	0.531	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 75.2	EDS
Estimated twinning fraction	0.013 for -h,l,k	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	2 of 124794 reflections (0.002%)	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6512	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.87% 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: NAG, BDP

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	0.35	0/5893	0.58	0/7958

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	5774	0	5591	148	0
2	A	53	0	38	1	0
3	A	685	0	0	27	0
All	All	6512	0	5629	148	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 13.

All (148) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:613:ASN:H	1:A:698:GLN:HE22	1.08	0.92

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:264:THR:O	1:A:268:THR:HG22	1.70	0.91
1:A:354:GLY:HA3	1:A:377:ILE:HD11	1.58	0.86
1:A:821:ASN:ND2	1:A:824:LEU:H	1.76	0.82
1:A:821:ASN:HD21	1:A:824:LEU:H	1.32	0.74
1:A:171:LYS:HB2	1:A:175:THR:HG21	1.73	0.70
1:A:644:SER:HB3	3:A:1281:HOH:O	1.93	0.67
1:A:650:ASN:HD21	1:A:832:GLN:HE22	1.41	0.67
1:A:754:MET:HG2	1:A:782:ILE:HG12	1.75	0.67
1:A:295:GLU:HB3	1:A:329:VAL:HG22	1.77	0.66
1:A:171:LYS:HE3	1:A:315:ASP:OD2	1.95	0.66
1:A:889:LYS:HG2	3:A:1688:HOH:O	1.95	0.66
1:A:708:LYS:HD3	1:A:715:GLU:HG3	1.78	0.66
1:A:546:ALA:HB1	3:A:1058:HOH:O	1.98	0.64
1:A:821:ASN:ND2	1:A:823:THR:H	1.95	0.63
1:A:288:VAL:HG13	1:A:294:TYR:OH	1.99	0.63
1:A:178:LEU:HD11	1:A:363:LEU:HD23	1.81	0.62
1:A:705:ASN:HB3	3:A:1240:HOH:O	2.00	0.61
1:A:336:ARG:HH12	2:A:1000:NAG:H2	1.64	0.61
1:A:874:ASN:HD22	1:A:874:ASN:C	2.04	0.61
1:A:820:ASN:HD22	1:A:825:GLN:HG2	1.65	0.61
1:A:603:HIS:ND1	3:A:1444:HOH:O	2.31	0.60
1:A:189:GLN:HG2	3:A:1546:HOH:O	2.03	0.59
1:A:821:ASN:C	1:A:821:ASN:HD22	2.06	0.59
1:A:286:SER:O	1:A:288:VAL:HG12	2.03	0.59
1:A:663:THR:HB	1:A:688:SER:HB3	1.85	0.58
1:A:821:ASN:HD22	1:A:823:THR:H	1.50	0.58
1:A:492:GLU:HG3	1:A:493:THR:N	2.18	0.58
1:A:855:LYS:HB2	1:A:886:VAL:HG12	1.84	0.58
1:A:652:THR:HB	3:A:1281:HOH:O	2.03	0.57
1:A:764:LYS:HD2	1:A:772:ASP:HB3	1.86	0.57
1:A:607:GLY:C	1:A:610:PRO:HD2	2.25	0.57
1:A:205:LEU:O	1:A:209:VAL:HG13	2.05	0.57
1:A:411:VAL:HG13	3:A:1623:HOH:O	2.03	0.57
1:A:254:ASN:ND2	1:A:256:SER:H	2.03	0.57
1:A:717:SER:O	1:A:723:LYS:HE3	2.06	0.56
1:A:585:ARG:HG3	1:A:766:ILE:HG22	1.88	0.56
1:A:488:MET:HE2	3:A:1602:HOH:O	2.05	0.55
1:A:845:SER:HB2	1:A:853:VAL:HG23	1.88	0.55
1:A:888:LYS:O	1:A:889:LYS:HB2	2.06	0.55
1:A:355:ARG:HH11	1:A:418:GLN:NE2	2.06	0.54
1:A:376:SER:O	1:A:379:GLN:HG2	2.08	0.53
1:A:889:LYS:HD3	3:A:1688:HOH:O	2.07	0.53

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:174:TYR:CD2	1:A:365:LYS:HG2	2.43	0.53
1:A:235:SER:HB2	1:A:293:ASP:HB2	1.90	0.53
1:A:288:VAL:HG22	1:A:288:VAL:O	2.09	0.53
1:A:713:ASP:C	1:A:714:LYS:HD2	2.30	0.52
1:A:386:GLN:NE2	3:A:1635:HOH:O	2.41	0.52
1:A:354:GLY:CA	1:A:377:ILE:HD11	2.37	0.52
1:A:495:GLN:CA	1:A:495:GLN:HE21	2.23	0.52
1:A:420:LEU:HD13	1:A:435:MET:CE	2.41	0.51
1:A:462:ARG:HD3	1:A:577:GLU:OE2	2.10	0.50
1:A:355:ARG:NH2	3:A:1623:HOH:O	2.44	0.50
1:A:173:THR:O	1:A:176:ASP:HB2	2.11	0.50
1:A:623:THR:HB	1:A:690:ASP:HB2	1.93	0.50
1:A:566:LEU:HD23	3:A:1062:HOH:O	2.10	0.50
1:A:213:LEU:HD11	1:A:265:VAL:HG22	1.93	0.50
1:A:230:SER:O	1:A:232:TYR:N	2.45	0.50
1:A:254:ASN:HD22	1:A:254:ASN:C	2.15	0.49
1:A:806:PHE:O	1:A:810:ILE:HG23	2.11	0.49
1:A:239:THR:HG22	1:A:243:ARG:HD2	1.94	0.49
1:A:322:THR:OG1	1:A:364:ARG:NH1	2.44	0.49
1:A:723:LYS:HD3	3:A:1203:HOH:O	2.12	0.49
1:A:216:ILE:HD12	1:A:226:TRP:CZ2	2.47	0.49
1:A:224:TYR:CD2	1:A:230:SER:HB3	2.47	0.49
1:A:882:PRO:HB2	1:A:884:GLN:NE2	2.27	0.49
1:A:604:TYR:HA	3:A:1444:HOH:O	2.13	0.48
1:A:882:PRO:HB2	1:A:884:GLN:HE22	1.78	0.48
1:A:228:LYS:HD2	1:A:229:PHE:CE1	2.48	0.48
1:A:598:ASN:HB2	1:A:739:LYS:O	2.14	0.48
1:A:303:ASN:HB3	1:A:359:ILE:HG21	1.96	0.47
1:A:303:ASN:HB3	1:A:359:ILE:CG2	2.44	0.47
1:A:701:LEU:HD12	1:A:778:GLU:HG2	1.96	0.47
1:A:374:ILE:O	1:A:378:GLU:HG3	2.15	0.47
1:A:634:LYS:HD2	3:A:1459:HOH:O	2.15	0.47
1:A:585:ARG:NH1	1:A:629:ASP:OD2	2.48	0.47
1:A:229:PHE:HE2	1:A:244:LYS:HE2	1.80	0.47
1:A:504:ILE:HG23	3:A:1016:HOH:O	2.15	0.47
1:A:354:GLY:O	1:A:358:VAL:HB	2.15	0.47
1:A:722:GLU:HG2	1:A:753:SER:OG	2.15	0.47
1:A:850:GLN:O	1:A:889:LYS:HD2	2.14	0.46
1:A:708:LYS:HD3	1:A:715:GLU:CG	2.45	0.46
1:A:254:ASN:ND2	1:A:254:ASN:C	2.69	0.46
1:A:888:LYS:O	1:A:889:LYS:CB	2.64	0.46
1:A:447:ALA:HB3	1:A:448:PRO:HD3	1.98	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:295:GLU:HB3	1:A:329:VAL:CG2	2.46	0.46
1:A:420:LEU:HD13	1:A:435:MET:HE3	1.96	0.46
1:A:764:LYS:HA	1:A:767:ASN:O	2.16	0.45
1:A:667:HIS:HD2	3:A:1142:HOH:O	1.99	0.45
1:A:329:VAL:HG12	1:A:357:LYS:HD2	1.98	0.45
1:A:381:PHE:O	1:A:437:THR:HG21	2.17	0.45
1:A:643:THR:HB	3:A:1058:HOH:O	2.16	0.45
1:A:277:HIS:CD2	3:A:1528:HOH:O	2.69	0.45
1:A:581:LYS:O	1:A:584:LYS:HD2	2.15	0.45
1:A:661:ASN:C	1:A:661:ASN:HD22	2.20	0.45
1:A:708:LYS:CD	1:A:715:GLU:HG3	2.46	0.45
1:A:607:GLY:O	1:A:610:PRO:HD2	2.16	0.45
1:A:502:LYS:HD3	1:A:537:SER:HB2	1.98	0.45
1:A:203:GLN:HG3	3:A:1374:HOH:O	2.17	0.45
1:A:768:GLU:HG3	3:A:1536:HOH:O	2.17	0.45
1:A:224:TYR:CG	1:A:230:SER:HB3	2.52	0.44
1:A:738:SER:O	1:A:800:ASN:HA	2.17	0.44
1:A:864:GLU:OE2	1:A:869:LYS:HD3	2.18	0.44
1:A:644:SER:HB2	1:A:873:TYR:HB3	1.98	0.44
1:A:568:LEU:HD23	1:A:592:GLY:HA2	2.00	0.44
1:A:420:LEU:HD12	1:A:488:MET:HE1	1.99	0.44
1:A:244:LYS:HE2	3:A:1366:HOH:O	2.16	0.44
1:A:585:ARG:HD3	1:A:629:ASP:OD1	2.17	0.44
1:A:231:ASN:O	1:A:237:ASN:ND2	2.49	0.44
1:A:771:SER:OG	1:A:773:LYS:HB2	2.17	0.44
1:A:661:ASN:H	1:A:661:ASN:HD22	1.66	0.44
1:A:436:GLN:O	1:A:436:GLN:HG2	2.17	0.44
1:A:858:VAL:HG12	3:A:1281:HOH:O	2.17	0.44
1:A:790:GLY:HA2	3:A:1132:HOH:O	2.17	0.43
1:A:863:LYS:HD2	1:A:868:TYR:OH	2.18	0.43
1:A:512:ASP:HB3	1:A:515:LYS:HG3	2.00	0.43
1:A:278:LYS:HG2	1:A:279:HIS:CD2	2.53	0.43
1:A:664:LEU:HA	1:A:685:GLN:O	2.18	0.43
1:A:623:THR:HA	1:A:691:THR:O	2.18	0.43
1:A:332:PRO:O	1:A:353:MET:HG2	2.18	0.43
1:A:749:LYS:HZ3	1:A:749:LYS:HG2	1.57	0.43
1:A:521:LYS:HD2	1:A:521:LYS:HA	1.80	0.43
1:A:681:GLY:O	1:A:792:SER:HB2	2.18	0.43
1:A:282:ASN:HD22	1:A:284:GLU:N	2.17	0.43
1:A:474:ALA:O	1:A:477:GLU:HB2	2.19	0.42
1:A:495:GLN:NE2	3:A:1005:HOH:O	2.51	0.42
1:A:486:ALA:O	1:A:494:LYS:HG3	2.20	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:882:PRO:O	1:A:885:GLU:HB2	2.19	0.42
1:A:223:ILE:HD13	1:A:223:ILE:N	2.34	0.42
1:A:496:ARG:HG3	1:A:496:ARG:HH11	1.85	0.42
1:A:457:ASP:HA	1:A:460:ARG:HG3	2.01	0.42
1:A:178:LEU:HA	1:A:178:LEU:HD23	1.90	0.41
1:A:302:ILE:O	1:A:306:LEU:HG	2.19	0.41
1:A:175:THR:HA	1:A:178:LEU:HB2	2.02	0.41
1:A:585:ARG:HH11	1:A:629:ASP:CG	2.23	0.41
1:A:350:LEU:O	1:A:353:MET:HB3	2.20	0.41
1:A:438:MET:HG2	1:A:441:TRP:CZ3	2.55	0.41
1:A:822:GLU:HG2	1:A:823:THR:HG23	2.03	0.41
1:A:767:ASN:HB3	1:A:770:GLN:CG	2.51	0.41
1:A:521:LYS:HE3	1:A:525:LEU:HG	2.02	0.41
1:A:282:ASN:HD22	1:A:284:GLU:H	1.67	0.41
1:A:821:ASN:HD21	1:A:824:LEU:N	2.07	0.41
1:A:748:LYS:O	1:A:749:LYS:C	2.59	0.40
1:A:869:LYS:HD2	1:A:869:LYS:N	2.36	0.40
1:A:404:TYR:CD1	1:A:407:ALA:HB3	2.57	0.40
1:A:281:TYR:CD2	1:A:324:VAL:HG11	2.57	0.40
1:A:234:THR:HG22	3:A:1533:HOH:O	2.20	0.40
1:A:391:TYR:CD2	1:A:549:LYS:HG3	2.57	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	717/721 (99%)	676 (94%)	38 (5%)	3 (0%)	43	16

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	231	ASN
1	A	848	SER

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	674	LYS

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	638/640 (100%)	583 (91%)	55 (9%)	15 1

All (55) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	196	ASP
1	A	209	VAL
1	A	219	GLN
1	A	221	ASP
1	A	222	ARG
1	A	228	LYS
1	A	252	VAL
1	A	254	ASN
1	A	263	GLU
1	A	268	THR
1	A	278	LYS
1	A	282	ASN
1	A	283	SER
1	A	285	LYS
1	A	286	SER
1	A	288	VAL
1	A	319	LYS
1	A	327	LYS
1	A	341	ASN
1	A	344	LYS
1	A	368	GLN
1	A	382	LYS
1	A	385	ASP
1	A	404	TYR
1	A	425	LYS
1	A	433	ASP
1	A	436	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	444	LYS
1	A	490	GLU
1	A	492	GLU
1	A	495	GLN
1	A	496	ARG
1	A	584	LYS
1	A	585	ARG
1	A	661	ASN
1	A	690	ASP
1	A	708	LYS
1	A	721	GLN
1	A	723	LYS
1	A	727	GLU
1	A	748	LYS
1	A	749	LYS
1	A	753	SER
1	A	773	LYS
1	A	817	LEU
1	A	821	ASN
1	A	853	VAL
1	A	854	LEU
1	A	855	LYS
1	A	864	GLU
1	A	867	GLU
1	A	869	LYS
1	A	874	ASN
1	A	888	LYS
1	A	889	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	197	GLN
1	A	202	ASN
1	A	254	ASN
1	A	277	HIS
1	A	282	ASN
1	A	349	ASN
1	A	368	GLN
1	A	379	GLN
1	A	418	GLN
1	A	436	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	495	GLN
1	A	580	ASN
1	A	661	ASN
1	A	667	HIS
1	A	683	ASN
1	A	698	GLN
1	A	705	ASN
1	A	729	GLN
1	A	759	GLN
1	A	820	ASN
1	A	821	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 ⓘ

4 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	NAG	A	1000	2	15,15,15	2.26	6 (40%)	21,21,21	1.13	3 (14%)
2	BDP	A	1001	2	11,12,13	5.01	3 (27%)	13,17,19	3.72	5 (38%)
2	NAG	A	1002	2	12,14,15	1.95	3 (25%)	15,19,21	0.98	0
2	BDP	A	1003	2	11,12,13	2.81	2 (18%)	13,17,19	1.50	1 (7%)

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	NAG	A	1000	2	-	0/6/26/26	0/1/1/1
2	BDP	A	1001	2	-	0/4/21/24	0/1/1/1
2	NAG	A	1002	2	-	0/6/23/26	0/1/1/1
2	BDP	A	1003	2	-	0/4/21/24	0/1/1/1

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1001	BDP	C5-C6	-13.28	1.23	1.53
2	A	1003	BDP	C4-C5	-8.38	1.37	1.53
2	A	1001	BDP	C4-C5	8.07	1.68	1.53
2	A	1000	NAG	O5-C1	5.15	1.53	1.43
2	A	1001	BDP	O5-C5	5.06	1.61	1.45
2	A	1002	NAG	C2-N2	4.92	1.52	1.46
2	A	1000	NAG	C2-N2	3.65	1.52	1.45
2	A	1000	NAG	C1-C2	2.91	1.56	1.53
2	A	1003	BDP	O4-C4	2.88	1.49	1.43
2	A	1002	NAG	O5-C5	2.72	1.50	1.45
2	A	1000	NAG	C4-C5	2.66	1.58	1.53
2	A	1000	NAG	C4-C3	2.31	1.58	1.52
2	A	1002	NAG	C4-C5	2.26	1.58	1.53
2	A	1000	NAG	O5-C5	2.22	1.49	1.44

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1001	BDP	C4-C5-C6	9.28	127.14	110.84
2	A	1001	BDP	O5-C5-C4	-5.89	98.20	110.32
2	A	1001	BDP	C3-C4-C5	5.87	120.14	109.48
2	A	1001	BDP	O4-C4-C5	-3.21	101.98	109.81
2	A	1003	BDP	O6B-C6-O6A	3.03	130.93	124.07
2	A	1001	BDP	O5-C5-C6	2.74	117.44	110.57
2	A	1000	NAG	O5-C1-C2	2.33	111.97	109.61
2	A	1000	NAG	C8-C7-N2	-2.30	111.61	116.11
2	A	1000	NAG	O7-C7-N2	2.03	126.14	121.90

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

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	719/721 (99%)	0.30	23 (3%)	45 49	31, 41, 68, 98	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	427	LYS	4.9
1	A	220	ALA	4.2
1	A	174	TYR	3.7
1	A	221	ASP	3.5
1	A	224	TYR	3.4
1	A	232	TYR	3.1
1	A	223	ILE	3.1
1	A	375	ARG	3.0
1	A	430	ILE	2.9
1	A	368	GLN	2.6
1	A	342	PRO	2.6
1	A	225	LEU	2.5
1	A	229	PHE	2.4
1	A	889	LYS	2.4
1	A	226	TRP	2.4
1	A	849	ASN	2.3
1	A	171	LYS	2.2
1	A	721	GLN	2.2
1	A	283	SER	2.2
1	A	173	THR	2.2
1	A	425	LYS	2.1
1	A	289	GLY	2.1
1	A	279	HIS	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	BDP	A	1003	12/13	0.19	6.87	62,64,71,82	0
2	BDP	A	1001	12/13	0.18	2.83	68,70,81,84	0
2	NAG	A	1000	15/15	0.19	2.48	66,70,76,85	0
2	NAG	A	1002	14/15	0.14	1.79	40,60,63,65	0

6.4 Ligands ⓘ

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.