



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 07:08 PM GMT

PDB ID : 1E72
Title : MYROSINASE FROM SINAPIS ALBA WITH BOUND GLUCO-HYDROX
IMOLACTAM AND SULFATE OR ASCORBATE
Authors : Burmeister, W.P.
Deposited on : 2000-08-23
Resolution : 1.60 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 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
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) : trunk26865

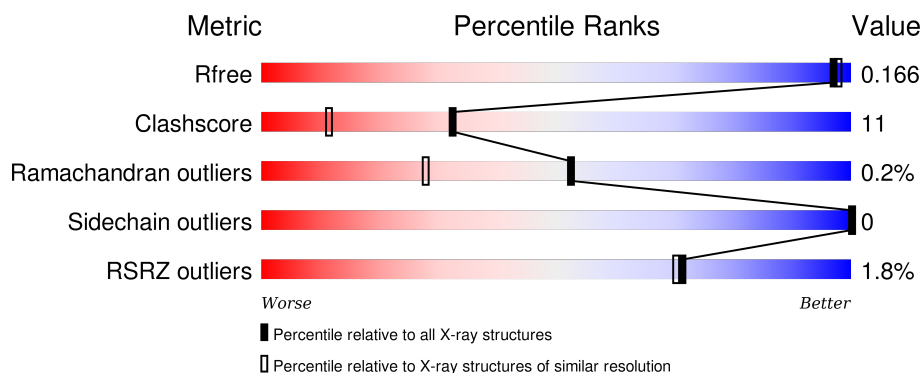
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.60 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}	91344	2475 (1.60-1.60)
Clashscore	102246	2732 (1.60-1.60)
Ramachandran outliers	100387	2654 (1.60-1.60)
Sidechain outliers	100360	2653 (1.60-1.60)
RSRZ outliers	91569	2479 (1.60-1.60)

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	M	501	<div> <div>2%</div> <div>87%</div> <div>11%</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
10	GOL	M	1510[A]	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
10	GOL	M	1510[B]	-	-	-	X
10	GOL	M	1511	-	-	-	X
10	GOL	M	1514	-	-	-	X
2	NAG	M	931	-	-	X	-
2	NAG	M	961	X	-	-	X
2	NAG	M	991	X	-	-	-
6	ASC	M	995	-	X	-	X
7	GOX	M	999	-	-	-	X
9	SO4	M	1503	-	-	-	X
9	SO4	M	1505[A]	-	-	-	X

2 Entry composition [i](#)

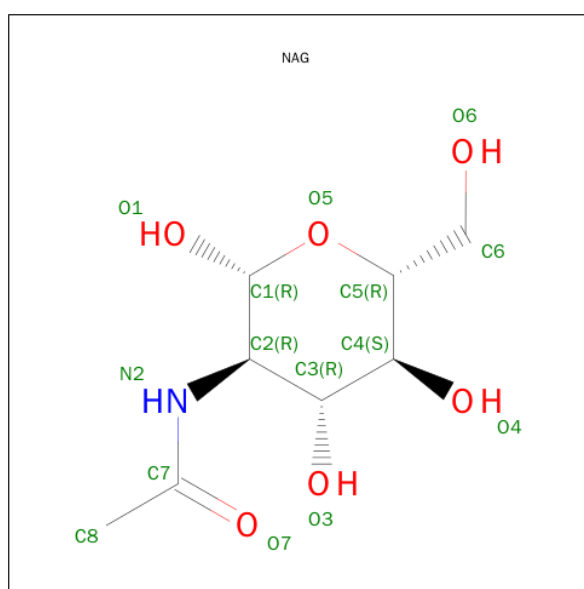
There are 11 unique types of molecules in this entry. The entry contains 5204 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 MYROSINASE MA1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	M	499	Total	C	N	O	S	0	21	0
			4082	2618	660	788	16			

- Molecule 2 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	M	1	Total	C	N	O	0	0
			14	8	1	5		
2	M	1	Total	C	N	O	0	0
			14	8	1	5		
2	M	1	Total	C	N	O	0	0
			14	8	1	5		
2	M	1	Total	C	N	O	0	0
			14	8	1	5		
2	M	1	Total	C	N	O	0	0
			14	8	1	5		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	M	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 3 is a polymer of unknown type called SUGAR (2-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	M	2	Total	C	N	O	0	0
			28	16	2	10		

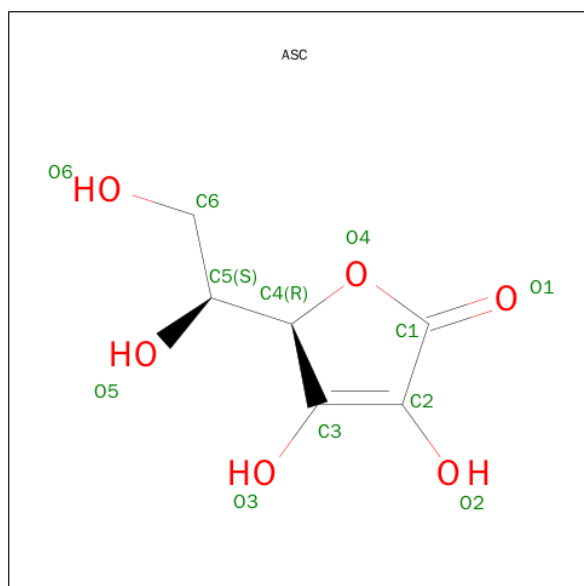
- Molecule 4 is a polymer of unknown type called SUGAR (5-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	M	5	Total	C	N	O	0	0
			58	33	2	23		

- Molecule 5 is a polymer of unknown type called SUGAR (6-MER).

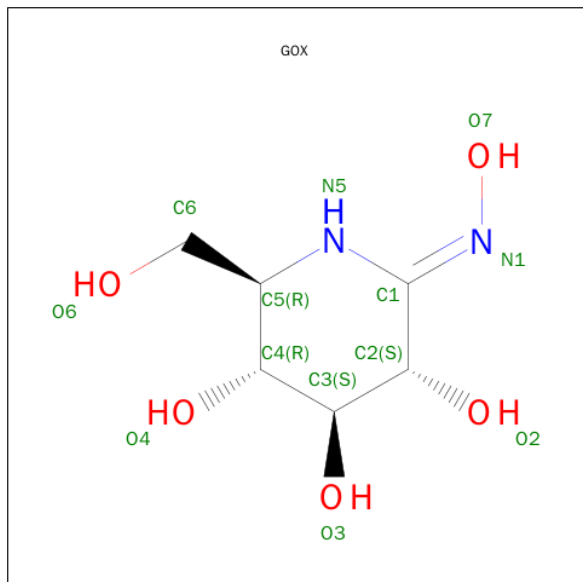
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	M	6	Total	C	N	O	0	0
			69	39	2	28		

- Molecule 6 is ASCORBIC ACID (three-letter code: ASC) (formula: $C_6H_8O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	M	1	Total	C	O	0	0
			12	6	6		

- Molecule 7 is (2S,3S,4R,5R)-6-(HYDROXYAMINO)-2-(HYDROXYMETHYL)-2,3,4,5-TETRAHYDROPYRIDINE-3,4,5-TRIOL (three-letter code: GOX) (formula: $C_6H_{12}N_2O_5$).

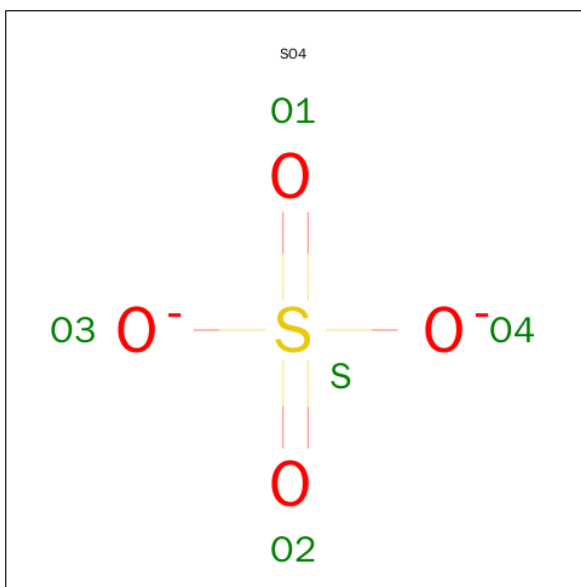


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	M	1	Total	C	N	O	0	0
			13	6	2	5		

- Molecule 8 is ZINC ION (three-letter code: ZN) (formula: Zn).

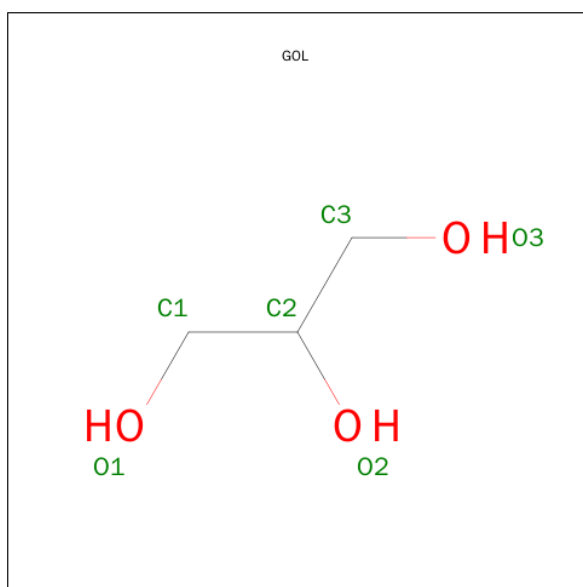
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	M	1	Total	Zn	0	0
			1	1		

- Molecule 9 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	M	1	Total	O	S	0	0
			5	4	1		
9	M	1	Total	O	S	0	0
			5	4	1		
9	M	1	Total	O	S	0	1
			10	8	2		
9	M	1	Total	O	S	0	0
			5	4	1		
9	M	1	Total	O	S	0	0
			5	4	1		
9	M	1	Total	O	S	0	0
			5	4	1		
9	M	1	Total	O	S	0	0
			5	4	1		

- Molecule 10 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
10	M	1	Total	C	O	0	1
			7	3	4		
10	M	1	Total	C	O	0	0
			6	3	3		
10	M	1	Total	C	O	0	0
			6	3	3		
10	M	1	Total	C	O	0	0
			6	3	3		

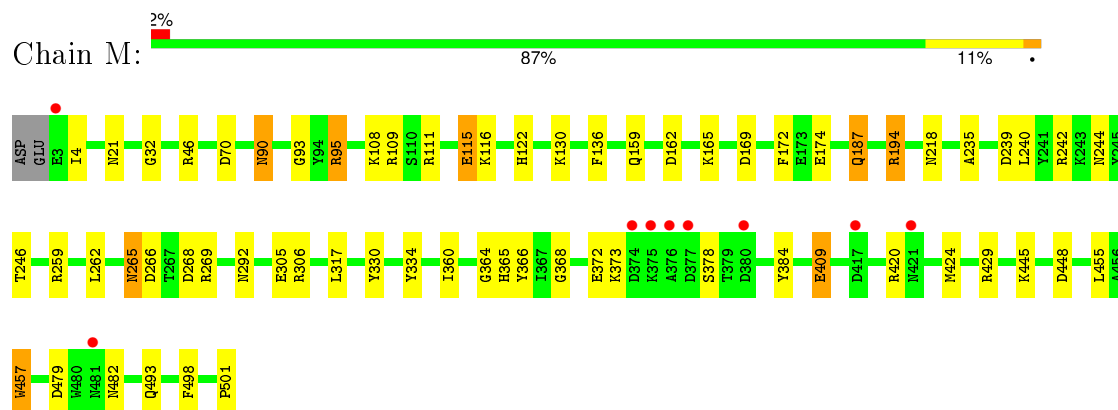
- Molecule 11 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	M	787	Total	O	0	0
			787	787		

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 ($\text{RSRZ} > 2$). 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: MYROSINASE MA1



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	135.30Å 137.20Å 80.60Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 1.60 9.99 – 1.60	Depositor EDS
% Data completeness (in resolution range)	97.5 (10.00-1.60) 97.5 (9.99-1.60)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.66 (at 1.60Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.129 , 0.157 0.143 , 0.166	Depositor DCC
R_{free} test set	4890 reflections (5.38%)	DCC
Wilson B-factor (Å ²)	14.2	Xtriage
Anisotropy	0.079	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.49 , 62.5	EDS
Estimated twinning fraction	0.016 for -k,-h,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 95834 reflections	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	5204	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.94% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: XYP, GOL, ZN, BMA, NAG, SO4, ASC, FUC, MAN, GOX

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	M	0.64	1/4290 (0.0%)	1.26	34/5833 (0.6%)

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	M	0	2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	M	501	PRO	N-CD	5.92	1.56	1.47

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	M	109	ARG	NE-CZ-NH2	-12.53	114.03	120.30
1	M	268	ASP	CB-CG-OD1	11.57	128.71	118.30
1	M	95	ARG	NE-CZ-NH2	-10.33	115.14	120.30
1	M	46	ARG	NE-CZ-NH2	-9.97	115.32	120.30
1	M	194	ARG	NE-CZ-NH2	9.32	124.96	120.30
1	M	269	ARG	NE-CZ-NH2	-9.15	115.72	120.30
1	M	242	ARG	NE-CZ-NH2	-9.02	115.79	120.30
1	M	115	GLU	OE1-CD-OE2	-8.61	112.97	123.30
1	M	269	ARG	NE-CZ-NH1	8.31	124.45	120.30
1	M	109	ARG	NH1-CZ-NH2	8.14	128.35	119.40
1	M	70	ASP	CB-CG-OD2	-8.14	110.98	118.30
1	M	162	ASP	CB-CG-OD1	7.28	124.85	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	M	482	ASN	CB-CG-OD1	6.94	135.48	121.60
1	M	457	TRP	O-C-N	-6.73	111.93	122.70
1	M	172	PHE	CB-CG-CD2	-6.35	116.35	120.80
1	M	194	ARG	NH1-CZ-NH2	-6.32	112.44	119.40
1	M	306	ARG	NE-CZ-NH1	5.91	123.25	120.30
1	M	479	ASP	CB-CG-OD1	5.90	123.61	118.30
1	M	457	TRP	CA-C-N	5.84	130.04	117.20
1	M	268	ASP	OD1-CG-OD2	-5.78	112.31	123.30
1	M	242	ARG	NE-CZ-NH1	5.68	123.14	120.30
1	M	330	TYR	CB-CG-CD2	-5.54	117.68	121.00
1	M	266	ASP	CB-CG-OD1	5.51	123.26	118.30
1	M	448	ASP	CB-CG-OD1	5.46	123.22	118.30
1	M	111	ARG	NE-CZ-NH2	-5.41	117.60	120.30
1	M	90	ASN	CB-CG-OD1	5.40	132.40	121.60
1	M	266	ASP	CB-CG-OD2	-5.40	113.44	118.30
1	M	109	ARG	NE-CZ-NH1	-5.38	117.61	120.30
1	M	409	GLU	OE1-CD-OE2	-5.38	116.84	123.30
1	M	482	ASN	CA-CB-CG	-5.36	101.60	113.40
1	M	429	ARG	NE-CZ-NH2	-5.34	117.63	120.30
1	M	259	ARG	NE-CZ-NH2	5.19	122.90	120.30
1	M	265	ASN	CB-CG-OD1	5.18	131.97	121.60
1	M	372	GLU	OE1-CD-OE2	-5.01	117.28	123.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	M	457	TRP	Mainchain,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	M	4082	0	3835	76	0
2	M	84	0	77	15	0
3	M	28	0	25	4	0
4	M	58	0	50	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	M	69	0	58	7	0
6	M	12	0	6	2	0
7	M	13	0	12	3	0
8	M	1	0	0	0	0
9	M	45	0	0	1	0
10	M	25	0	30	1	0
11	M	787	0	0	21	0
All	All	5204	0	4093	83	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 11.

All (83) 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:M:90:ASN:HD21	2:M:911:NAG:C1	0.95	1.59
1:M:21:ASN:HD21	2:M:901:NAG:C1	0.95	1.58
1:M:244:ASN:HD21	2:M:931:NAG:C1	0.96	1.58
1:M:292:ASN:HD21	5:M:951:NAG:C1	0.95	1.55
1:M:265:ASN:HD21	4:M:942:NAG:C1	0.92	1.54
1:M:218:ASN:HD21	3:M:921:NAG:C1	0.93	1.54
1:M:360[B]:ILE:HD11	1:M:366[B]:TYR:CZ	1.45	1.50
1:M:360[B]:ILE:CD1	1:M:366[B]:TYR:CZ	2.05	1.37
1:M:360[B]:ILE:HD11	1:M:366[B]:TYR:CE1	1.68	1.28
1:M:360[B]:ILE:CG1	1:M:366[B]:TYR:CE1	2.21	1.24
1:M:360[B]:ILE:HD11	1:M:366[B]:TYR:OH	1.39	1.17
1:M:360[B]:ILE:HG12	1:M:366[B]:TYR:CE1	1.84	1.12
1:M:360[B]:ILE:HG12	1:M:366[B]:TYR:CD1	1.83	1.12
1:M:360[B]:ILE:CD1	1:M:366[B]:TYR:CE1	2.24	1.12
1:M:360[B]:ILE:CD1	1:M:366[B]:TYR:OH	2.08	0.95
5:M:954:BMA:C4	11:M:2739:HOH:O	2.14	0.94
1:M:246:THR:HG22	11:M:2390:HOH:O	1.67	0.93
6:M:995:ASC:H4	7:M:999:GOX:O7	1.72	0.89
1:M:165:LYS:NZ	2:M:931:NAG:H82	1.89	0.88
1:M:360[B]:ILE:HD13	1:M:366[B]:TYR:CZ	2.07	0.87
5:M:954:BMA:C5	11:M:2739:HOH:O	2.22	0.86
5:M:954:BMA:O4	11:M:2739:HOH:O	1.94	0.85
1:M:4:ILE:HD11	1:M:445:LYS:HD2	1.62	0.79
1:M:165:LYS:HZ1	2:M:931:NAG:H82	1.52	0.74
1:M:265:ASN:HD21	4:M:942:NAG:C2	1.97	0.73
1:M:130:LYS:HG3	11:M:2252:HOH:O	1.91	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:187:GLN:OE1	7:M:999:GOX:N1	2.24	0.68
1:M:218:ASN:HD21	3:M:921:NAG:C2	1.98	0.67
6:M:995:ASC:H5	11:M:2755:HOH:O	1.96	0.66
5:M:954:BMA:C6	11:M:2739:HOH:O	2.44	0.66
1:M:116:LYS:HG2	11:M:2232:HOH:O	1.97	0.65
1:M:360[B]:ILE:HG13	1:M:366[B]:TYR:CE1	2.29	0.64
1:M:169[B]:ASP:HB2	1:M:240:LEU:HD21	1.80	0.63
1:M:130:LYS:HB3	11:M:2265:HOH:O	2.00	0.62
1:M:360[B]:ILE:CG1	1:M:366[B]:TYR:CD1	2.62	0.60
1:M:360[B]:ILE:CG1	1:M:366[B]:TYR:CZ	2.71	0.60
1:M:365:HIS:HD2	11:M:2255:HOH:O	1.85	0.59
1:M:424:MET:HE3	11:M:2693:HOH:O	2.03	0.59
1:M:218:ASN:ND2	3:M:921:NAG:C2	2.64	0.58
1:M:21:ASN:HD21	2:M:901:NAG:C2	2.00	0.57
1:M:365:HIS:HE1	11:M:2550:HOH:O	1.86	0.57
1:M:108:LYS:HE2	11:M:2317:HOH:O	2.06	0.55
1:M:265:ASN:ND2	4:M:942:NAG:C2	2.64	0.55
1:M:244:ASN:ND2	2:M:931:NAG:O5	2.40	0.54
1:M:239:ASP:OD1	11:M:2385:HOH:O	2.19	0.53
1:M:90:ASN:ND2	2:M:911:NAG:C2	2.66	0.53
1:M:115:GLU:HG3	11:M:2229:HOH:O	2.09	0.52
1:M:240:LEU:HD13	2:M:931:NAG:H83	1.92	0.52
1:M:265:ASN:ND2	4:M:942:NAG:O5	2.37	0.52
1:M:4:ILE:HD11	1:M:445:LYS:CD	2.36	0.52
1:M:360[B]:ILE:HG12	1:M:366[B]:TYR:CZ	2.42	0.50
1:M:240:LEU:CD1	2:M:931:NAG:H83	2.42	0.49
10:M:1514:GOL:H11	11:M:2785:HOH:O	2.11	0.49
1:M:4:ILE:CD1	1:M:445:LYS:HD2	2.36	0.49
1:M:265:ASN:CG	4:M:942:NAG:C1	2.73	0.49
1:M:409:GLU:OE1	7:M:999:GOX:C1	2.60	0.49
1:M:360[A]:ILE:HG23	11:M:2538:HOH:O	2.13	0.47
1:M:218:ASN:ND2	3:M:921:NAG:O5	2.43	0.47
1:M:493:GLN:NE2	11:M:2693:HOH:O	2.47	0.47
1:M:95:ARG:HB2	1:M:455:LEU:HD13	1.97	0.47
1:M:360[A]:ILE:HD11	1:M:364:GLY:HA2	1.97	0.47
1:M:95:ARG:HA	1:M:136:PHE:O	2.15	0.47
1:M:194:ARG:HD2	9:M:1505[B]:SO4:O1	2.15	0.46
1:M:360[B]:ILE:HG12	1:M:366[B]:TYR:CG	2.44	0.46
1:M:244:ASN:HD21	2:M:931:NAG:C2	2.01	0.46
1:M:292:ASN:ND2	5:M:951:NAG:C2	2.68	0.45
1:M:445:LYS:HE2	11:M:2641:HOH:O	2.15	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:165:LYS:CE	2:M:931:NAG:H82	2.46	0.44
1:M:420:ARG:O	1:M:424:MET:HG2	2.17	0.44
1:M:165:LYS:HZ2	2:M:931:NAG:H82	1.78	0.44
1:M:21:ASN:ND2	2:M:901:NAG:C2	2.71	0.42
1:M:21:ASN:HA	1:M:498:PHE:CD2	2.55	0.42
1:M:159:GLN:NE2	11:M:2309:HOH:O	2.53	0.41
1:M:90:ASN:ND2	2:M:911:NAG:O5	2.45	0.41
1:M:373:LYS:NZ	1:M:378:SER:OG	2.44	0.41
1:M:122:HIS:HE1	1:M:174:GLU:O	2.02	0.41
1:M:235:ALA:HB1	1:M:317:LEU:HG	2.02	0.41
1:M:368:GLY:HA3	1:M:384:TYR:O	2.21	0.41
1:M:373:LYS:NZ	11:M:2559:HOH:O	2.52	0.41
1:M:32:GLY:HA3	1:M:93:GLY:O	2.21	0.41
1:M:262:LEU:O	1:M:334:TYR:HA	2.21	0.41
1:M:292:ASN:ND2	5:M:951:NAG:O5	2.42	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	M	518/501 (103%)	504 (97%)	13 (2%)	1 (0%)	52 28

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	M	187	GLN

5.3.2 Protein sidechains [i](#)

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	M	456/437 (104%)	456 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	M	90	ASN
1	M	218	ASN
1	M	244	ASN
1	M	265	ASN
1	M	292	ASN
1	M	365	HIS

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](#)

13 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
3	NAG	M	921	1,3	14,14,15	1.52	4 (28%)	15,19,21	2.99	6 (40%)
3	NAG	M	923	3	14,14,15	1.01	0	15,19,21	3.01	7 (46%)
4	FUC	M	941	4	10,10,11	1.35	1 (10%)	14,14,16	1.93	3 (21%)
4	NAG	M	942	1,4	14,14,15	0.99	1 (7%)	15,19,21	3.09	7 (46%)
4	NAG	M	943	4	14,14,15	1.13	1 (7%)	15,19,21	0.99	1 (6%)
4	BMA	M	944	4	11,11,12	1.96	2 (18%)	14,15,17	1.11	1 (7%)
4	XYP	M	945	4	9,9,10	0.91	0	12,12,14	1.90	3 (25%)
5	NAG	M	951	1,5	14,14,15	1.34	1 (7%)	15,19,21	2.48	5 (33%)
5	FUC	M	952	5	10,10,11	1.65	3 (30%)	14,14,16	2.63	6 (42%)
5	NAG	M	953	5	14,14,15	1.17	1 (7%)	15,19,21	2.00	4 (26%)
5	BMA	M	954	5	11,11,12	2.21	3 (27%)	14,15,17	5.29	12 (85%)
5	XYP	M	955	5	9,9,10	1.20	1 (11%)	12,12,14	2.26	4 (33%)
5	MAN	M	956	5	11,11,12	1.20	1 (9%)	14,15,17	1.91	3 (21%)

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	NAG	M	921	1,3	-	0/6/23/26	0/1/1/1
3	NAG	M	923	3	-	0/6/23/26	0/1/1/1
4	FUC	M	941	4	-	0/0/17/20	0/1/1/1
4	NAG	M	942	1,4	-	0/6/23/26	0/1/1/1
4	NAG	M	943	4	-	0/6/23/26	0/1/1/1
4	BMA	M	944	4	-	0/2/19/22	0/1/1/1
4	XYP	M	945	4	-	0/0/14/17	0/1/1/1
5	NAG	M	951	1,5	-	0/6/23/26	0/1/1/1
5	FUC	M	952	5	-	0/0/17/20	0/1/1/1
5	NAG	M	953	5	-	0/6/23/26	0/1/1/1
5	BMA	M	954	5	-	0/2/19/22	0/1/1/1
5	XYP	M	955	5	-	0/0/14/17	0/1/1/1
5	MAN	M	956	5	-	0/2/19/22	0/1/1/1

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	M	954	BMA	C2-C3	-5.22	1.45	1.52
4	M	944	BMA	C2-C3	-5.15	1.45	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	M	921	NAG	O7-C7	-2.61	1.17	1.23
5	M	951	NAG	C1-C2	-2.54	1.48	1.52
4	M	942	NAG	C3-C2	2.20	1.57	1.52
3	M	921	NAG	O5-C5	2.24	1.48	1.43
3	M	921	NAG	C3-C2	2.24	1.57	1.52
3	M	921	NAG	O4-C4	2.26	1.48	1.43
5	M	952	FUC	C1-C2	2.33	1.57	1.52
5	M	956	MAN	O5-C5	2.39	1.48	1.43
5	M	954	BMA	O3-C3	2.41	1.48	1.43
5	M	952	FUC	C6-C5	2.43	1.57	1.51
5	M	955	XYP	C2B-C3B	2.46	1.55	1.52
4	M	943	NAG	C1-C2	2.54	1.56	1.52
5	M	952	FUC	C2-C3	2.58	1.56	1.52
5	M	953	NAG	C1-C2	2.73	1.56	1.52
4	M	944	BMA	C4-C5	2.75	1.58	1.53
4	M	941	FUC	C2-C3	2.85	1.56	1.52
5	M	954	BMA	C4-C5	3.35	1.60	1.53

All (62) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	M	923	NAG	C1-O5-C5	-9.29	100.45	112.25
3	M	921	NAG	C1-O5-C5	-9.00	100.83	112.25
5	M	954	BMA	C1-O5-C5	-8.49	101.47	112.25
5	M	954	BMA	O4-C4-C3	-8.38	91.47	110.34
5	M	951	NAG	C1-O5-C5	-7.30	102.98	112.25
5	M	954	BMA	O5-C5-C6	-7.20	91.76	107.35
5	M	954	BMA	C2-C3-C4	-6.84	99.43	111.04
4	M	942	NAG	C1-O5-C5	-6.45	104.07	112.25
4	M	942	NAG	C2-N2-C7	-5.93	115.42	123.04
5	M	952	FUC	C1-C2-C3	-5.91	102.55	109.54
5	M	953	NAG	C2-N2-C7	-4.71	116.98	123.04
5	M	955	XYP	C4B-C3B-C2B	-4.59	106.70	111.24
4	M	941	FUC	O3-C3-C2	-4.54	101.80	110.00
4	M	942	NAG	C8-C7-N2	-4.23	108.01	116.11
5	M	953	NAG	O4-C4-C3	-3.93	101.50	110.34
5	M	955	XYP	C5B-C4B-C3B	-3.82	105.02	109.54
3	M	921	NAG	O4-C4-C5	-3.74	99.33	109.24
4	M	945	XYP	O2B-C2B-C3B	-3.63	102.81	110.12
3	M	921	NAG	C2-N2-C7	-3.51	118.53	123.04
5	M	954	BMA	C1-C2-C3	-3.21	105.75	109.54
4	M	942	NAG	C4-C3-C2	-3.16	106.32	111.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	M	942	NAG	O3-C3-C2	-3.16	102.86	109.11
4	M	941	FUC	O2-C2-C3	-3.10	103.88	110.12
5	M	953	NAG	O4-C4-C5	-3.06	101.12	109.24
5	M	952	FUC	O2-C2-C1	-2.91	103.38	109.21
5	M	956	MAN	O2-C2-C3	-2.90	104.28	110.12
5	M	955	XYP	O2B-C2B-C3B	-2.88	104.32	110.12
5	M	956	MAN	C2-C3-C4	-2.83	106.24	111.04
3	M	923	NAG	C2-N2-C7	-2.81	119.43	123.04
5	M	952	FUC	O5-C5-C6	-2.80	101.49	106.13
5	M	954	BMA	O2-C2-C3	-2.71	104.66	110.12
4	M	941	FUC	O3-C3-C4	-2.69	104.29	110.34
3	M	923	NAG	C6-C5-C4	-2.67	106.43	113.02
5	M	952	FUC	C6-C5-C4	-2.65	107.86	113.08
3	M	921	NAG	C8-C7-N2	-2.65	111.04	116.11
3	M	921	NAG	O4-C4-C3	-2.57	104.56	110.34
3	M	923	NAG	C8-C7-N2	-2.56	111.20	116.11
3	M	923	NAG	O4-C4-C5	-2.50	102.62	109.24
5	M	953	NAG	C3-C2-N2	-2.47	104.65	110.56
3	M	923	NAG	O4-C4-C3	-2.41	104.92	110.34
5	M	951	NAG	C8-C7-N2	-2.36	111.58	116.11
5	M	951	NAG	C4-C3-C2	-2.32	107.63	111.23
5	M	952	FUC	C3-C4-C5	-2.31	105.83	109.72
5	M	951	NAG	O3-C3-C2	-2.25	104.65	109.11
4	M	945	XYP	C5B-C4B-C3B	-2.23	106.90	109.54
4	M	942	NAG	O4-C4-C3	-2.22	105.33	110.34
5	M	954	BMA	O3-C3-C4	-2.16	105.46	110.34
5	M	954	BMA	O3-C3-C2	-2.15	106.11	110.00
3	M	921	NAG	O3-C3-C2	-2.08	104.98	109.11
4	M	943	NAG	O3-C3-C2	-2.08	104.99	109.11
3	M	923	NAG	O6-C6-C5	-2.05	104.56	111.33
4	M	944	BMA	C2-C3-C4	2.14	114.68	111.04
5	M	954	BMA	O5-C1-C2	2.17	114.38	110.86
5	M	951	NAG	C3-C2-N2	2.35	116.19	110.56
5	M	955	XYP	O3B-C3B-C4B	2.99	115.39	110.00
4	M	942	NAG	O7-C7-N2	3.41	128.81	121.86
5	M	956	MAN	C1-C2-C3	3.93	114.19	109.54
4	M	945	XYP	C1B-C2B-C3B	4.31	114.64	109.54
5	M	952	FUC	C1-O5-C5	4.45	119.26	112.38
5	M	954	BMA	C3-C4-C5	4.62	118.25	110.20
5	M	954	BMA	O4-C4-C5	6.51	126.48	109.24
5	M	954	BMA	C6-C5-C4	7.33	131.09	113.02

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 16 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	M	921	NAG	4	0
4	M	942	NAG	5	0
5	M	951	NAG	3	0
5	M	954	BMA	4	0

5.6 Ligand geometry

Of 23 ligands modelled in this entry, 1 is monoatomic - leaving 22 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
9	SO4	M	1503	-	4,4,4	1.18	0	6,6,6	0.59	0
9	SO4	M	1504	-	4,4,4	1.01	0	6,6,6	0.49	0
9	SO4	M	1505[A]	-	4,4,4	1.21	0	6,6,6	1.20	1 (16%)
9	SO4	M	1505[B]	-	4,4,4	0.92	0	6,6,6	0.81	0
9	SO4	M	1506	-	4,4,4	0.59	0	6,6,6	0.51	0
9	SO4	M	1507	-	4,4,4	1.19	1 (25%)	6,6,6	0.79	0
9	SO4	M	1508	-	4,4,4	1.17	0	6,6,6	0.84	0
9	SO4	M	1509	-	4,4,4	0.93	0	6,6,6	0.12	0
10	GOL	M	1510[A]	-	5,5,5	0.85	0	5,5,5	1.30	1 (20%)
10	GOL	M	1510[B]	-	5,5,5	0.85	0	5,5,5	1.80	2 (40%)
10	GOL	M	1511	-	5,5,5	0.53	0	5,5,5	1.00	0
10	GOL	M	1512	-	5,5,5	0.26	0	5,5,5	0.75	0
9	SO4	M	1513	6	4,4,4	0.78	0	6,6,6	0.57	0
10	GOL	M	1514	-	5,5,5	0.25	0	5,5,5	0.61	0
2	NAG	M	901	1	14,14,15	1.26	1 (7%)	15,19,21	1.62	3 (20%)
2	NAG	M	911	1	14,14,15	1.19	1 (7%)	15,19,21	2.05	3 (20%)
2	NAG	M	931	1	14,14,15	1.63	3 (21%)	15,19,21	7.48	8 (53%)
2	NAG	M	961	1	14,14,15	1.28	1 (7%)	15,19,21	2.17	2 (13%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NAG	M	971	1	14,14,15	1.24	1 (7%)	15,19,21	0.88	1 (6%)
2	NAG	M	991	1	14,14,15	1.37	1 (7%)	15,19,21	2.07	5 (33%)
6	ASC	M	995	9	12,12,12	4.59	9 (75%)	17,17,17	4.25	12 (70%)
7	GOX	M	999	-	12,13,13	1.68	2 (16%)	9,18,18	3.27	5 (55%)

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
9	SO4	M	1503	-	-	0/0/0/0	0/0/0/0
9	SO4	M	1504	-	-	0/0/0/0	0/0/0/0
9	SO4	M	1505[A]	-	-	0/0/0/0	0/0/0/0
9	SO4	M	1505[B]	-	-	0/0/0/0	0/0/0/0
9	SO4	M	1506	-	-	0/0/0/0	0/0/0/0
9	SO4	M	1507	-	-	0/0/0/0	0/0/0/0
9	SO4	M	1508	-	-	0/0/0/0	0/0/0/0
9	SO4	M	1509	-	-	0/0/0/0	0/0/0/0
10	GOL	M	1510[A]	-	-	0/4/4/4	0/0/0/0
10	GOL	M	1510[B]	-	-	0/4/4/4	0/0/0/0
10	GOL	M	1511	-	-	0/4/4/4	0/0/0/0
10	GOL	M	1512	-	-	0/4/4/4	0/0/0/0
9	SO4	M	1513	6	-	0/0/0/0	0/0/0/0
10	GOL	M	1514	-	-	0/4/4/4	0/0/0/0
2	NAG	M	901	1	-	0/6/23/26	0/1/1/1
2	NAG	M	911	1	-	0/6/23/26	0/1/1/1
2	NAG	M	931	1	-	0/6/23/26	0/1/1/1
2	NAG	M	961	1	1/1/5/7	0/6/23/26	0/1/1/1
2	NAG	M	971	1	-	0/6/23/26	0/1/1/1
2	NAG	M	991	1	1/1/5/7	0/6/23/26	0/1/1/1
6	ASC	M	995	9	-	0/6/22/22	0/1/1/1
7	GOX	M	999	-	-	0/2/24/24	0/1/1/1

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	M	995	ASC	O1-C1	-8.58	1.03	1.21
2	M	911	NAG	O7-C7	-3.93	1.14	1.23
2	M	971	NAG	O7-C7	-3.80	1.14	1.23
2	M	931	NAG	O7-C7	-3.70	1.14	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	M	991	NAG	O7-C7	-3.61	1.14	1.23
2	M	961	NAG	O7-C7	-3.55	1.15	1.23
2	M	901	NAG	O7-C7	-3.28	1.15	1.23
2	M	931	NAG	C2-N2	-3.14	1.40	1.46
6	M	995	ASC	O4-C4	-2.75	1.40	1.45
6	M	995	ASC	O2-C2	-2.69	1.24	1.33
7	M	999	GOX	O7-N1	-2.65	1.33	1.40
2	M	931	NAG	C8-C7	-2.38	1.45	1.50
9	M	1507	SO4	O2-S	2.01	1.54	1.47
6	M	995	ASC	C5-C4	2.16	1.56	1.53
6	M	995	ASC	C2-C3	3.09	1.44	1.35
6	M	995	ASC	O3-C3	3.99	1.45	1.33
7	M	999	GOX	C4-C5	4.20	1.62	1.52
6	M	995	ASC	C4-C3	4.94	1.57	1.50
6	M	995	ASC	O4-C1	5.17	1.43	1.35
6	M	995	ASC	C2-C1	9.08	1.67	1.45

All (43) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	M	931	NAG	O7-C7-N2	-9.70	102.08	121.86
6	M	995	ASC	O4-C1-C2	-8.33	102.88	109.90
2	M	931	NAG	C1-O5-C5	-8.05	102.03	112.25
2	M	911	NAG	C1-O5-C5	-5.68	105.04	112.25
7	M	999	GOX	O3-C3-C2	-5.12	100.25	109.49
7	M	999	GOX	C3-C4-C5	-5.04	103.39	111.23
6	M	995	ASC	O4-C1-O1	-4.60	115.88	121.29
2	M	901	NAG	C1-O5-C5	-4.45	106.59	112.25
2	M	991	NAG	C4-C3-C2	-4.21	104.68	111.23
2	M	931	NAG	C4-C3-C2	-3.55	105.71	111.23
2	M	991	NAG	O3-C3-C2	-3.30	102.58	109.11
2	M	911	NAG	C4-C3-C2	-3.27	106.15	111.23
2	M	961	NAG	O3-C3-C2	-3.11	102.95	109.11
2	M	991	NAG	C2-N2-C7	-2.69	119.59	123.04
7	M	999	GOX	O4-C4-C5	-2.46	104.24	109.11
2	M	911	NAG	O3-C3-C2	-2.37	104.41	109.11
2	M	991	NAG	C6-C5-C4	-2.34	107.25	113.02
9	M	1505[A]	SO4	O2-S-O1	-2.30	102.20	109.50
2	M	901	NAG	C4-C3-C2	-2.30	107.66	111.23
6	M	995	ASC	O2-C2-C3	-2.26	120.83	128.87
6	M	995	ASC	C4-C3-C2	-2.19	104.98	109.24
2	M	901	NAG	O3-C3-C2	-2.13	104.89	109.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	M	971	NAG	O5-C5-C6	-2.04	102.92	107.35
2	M	931	NAG	O7-C7-C8	-2.04	118.33	122.06
10	M	1510[B]	GOL	O2-C2-C1	2.49	120.06	108.65
10	M	1510[A]	GOL	O2-C2-C1	2.49	120.06	108.65
6	M	995	ASC	O3-C3-C2	2.55	140.17	132.38
6	M	995	ASC	O4-C4-C5	2.66	114.19	109.71
2	M	931	NAG	C3-C4-C5	2.80	115.08	110.20
10	M	1510[B]	GOL	O1-C1-C2	2.86	124.04	110.18
7	M	999	GOX	C2-C1-N5	2.97	124.57	118.06
6	M	995	ASC	C5-C4-C3	3.28	118.78	114.14
6	M	995	ASC	O2-C2-C1	3.36	130.81	121.86
6	M	995	ASC	C4-O4-C1	3.90	113.33	109.29
2	M	991	NAG	C1-O5-C5	4.05	117.39	112.25
2	M	931	NAG	C3-C2-N2	4.07	120.30	110.56
6	M	995	ASC	C6-C5-C4	4.24	119.11	111.95
7	M	999	GOX	O2-C2-C3	5.30	120.75	109.95
2	M	961	NAG	C1-O5-C5	7.38	121.61	112.25
6	M	995	ASC	O4-C4-C3	7.50	110.05	104.06
6	M	995	ASC	O1-C1-C2	8.72	141.23	129.44
2	M	931	NAG	C8-C7-N2	12.26	139.57	116.11
2	M	931	NAG	C2-N2-C7	22.00	151.31	123.04

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	M	961	NAG	C1
2	M	991	NAG	C1

There are no torsion outliers.

There are no ring outliers.

7 monomers are involved in 21 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
9	M	1505[B]	SO4	1	0
10	M	1514	GOL	1	0
2	M	901	NAG	3	0
2	M	911	NAG	3	0
2	M	931	NAG	9	0
6	M	995	ASC	2	0
7	M	999	GOX	3	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	M	499/501 (99%)	-0.59	9 (1%) 71 70	11, 15, 27, 54	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	M	376	ALA	9.6
1	M	380	ASP	3.9
1	M	417	ASP	3.7
1	M	374	ASP	3.7
1	M	375	LYS	3.7
1	M	3	GLU	3.4
1	M	481	ASN	2.6
1	M	377	ASP	2.6
1	M	421	ASN	2.4

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(Å ²)	Q<0.9
3	NAG	M	923	14/15	0.55	0.30	1.49	38,43,47,48	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	NAG	M	921	14/15	0.93	0.08	1.24	20,24,29,32	0
4	NAG	M	942	14/15	0.93	0.08	0.30	18,21,23,25	0
4	XYP	M	945	9/10	0.63	0.38	-	42,43,45,46	0
4	FUC	M	941	10/11	0.87	0.20	-	29,33,34,37	0
5	NAG	M	953	14/15	0.93	0.10	-	20,24,29,32	0
5	MAN	M	956	11/12	0.57	0.31	-	26,37,41,41	0
5	BMA	M	954	11/12	0.86	0.14	-	24,28,33,35	0
5	XYP	M	955	9/10	0.54	0.27	-	37,40,42,45	0
4	BMA	M	944	11/12	0.75	0.32	-	38,41,43,44	0
5	NAG	M	951	14/15	0.93	0.09	-	18,19,23,25	0
5	FUC	M	952	10/11	0.91	0.10	-	21,23,25,26	0
4	NAG	M	943	14/15	0.92	0.09	-	26,29,34,36	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	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	NAG	M	961	14/15	0.61	0.45	60.94	36,40,46,47	0
10	GOL	M	1514	6/6	0.74	0.30	17.87	44,44,45,45	6
9	SO4	M	1505[A]	5/5	0.77	0.28	16.83	31,33,33,33	5
7	GOX	M	999	13/13	0.84	0.17	11.19	17,20,27,30	0
10	GOL	M	1510[B]	6/6	0.88	0.13	11.06	15,18,20,20	2
10	GOL	M	1510[A]	6/6	0.88	0.13	9.15	15,18,19,20	2
9	SO4	M	1503	5/5	0.98	0.12	5.99	24,26,28,30	5
6	ASC	M	995	12/12	0.88	0.15	3.49	13,23,23,23	12
10	GOL	M	1511	6/6	0.93	0.10	2.88	15,19,22,28	6
9	SO4	M	1513	5/5	0.98	0.07	1.80	20,21,21,22	5
2	NAG	M	901	14/15	0.72	0.18	1.57	31,34,38,39	0
9	SO4	M	1509	5/5	0.82	0.24	-	41,41,41,42	5
2	NAG	M	991	14/15	0.59	0.41	-	37,39,47,48	0
9	SO4	M	1504	5/5	0.94	0.18	-	26,29,31,32	5
8	ZN	M	1502	1/1	1.00	0.02	-	12,12,12,12	1
2	NAG	M	931	14/15	0.65	0.30	-	36,42,45,47	0
10	GOL	M	1512	6/6	0.76	0.21	-	34,34,35,37	6
9	SO4	M	1508	5/5	0.60	0.27	-	29,32,34,35	5
9	SO4	M	1506	5/5	0.99	0.05	-	20,20,24,25	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	NAG	M	971	14/15	0.57	0.52	-	56,63,65,67	0
9	SO4	M	1507	5/5	0.82	0.23	-	33,33,35,35	5
2	NAG	M	911	14/15	0.85	0.20	-	25,28,30,32	0
9	SO4	M	1505[B]	5/5	0.77	0.28	-	43,44,44,44	5

6.5 Other polymers [i](#)

There are no such residues in this entry.