



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:34 AM GMT

PDB ID : 2HOX
Title : alliinase from allium sativum (garlic)
Authors : Shimon, L.J.W.; Rabinkov, A.; Wilcheck, M.; Mirelman, D.; Frolow, F.
Deposited on : 2006-07-17
Resolution : 1.40 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

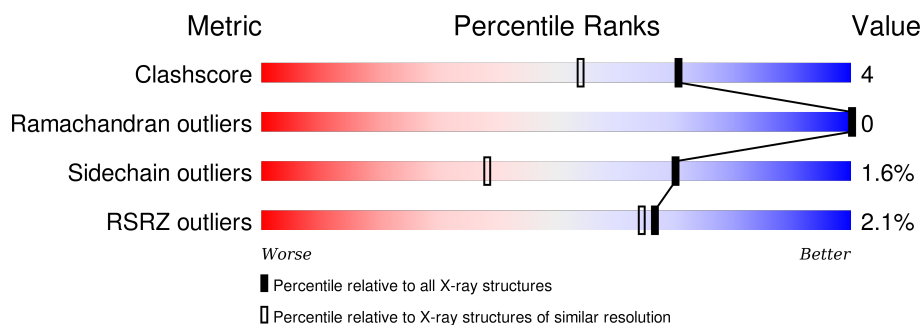
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.40 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))
Clashscore	102246	1295 (1.40-1.40)
Ramachandran outliers	100387	1259 (1.40-1.40)
Sidechain outliers	100360	1258 (1.40-1.40)
RSRZ outliers	91569	1198 (1.40-1.40)

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	427	<div> <div>2%</div> <div> <div></div> <div>91%</div> <div>8%</div> </div> </div>
1	B	427	<div> <div>2%</div> <div> <div></div> <div>93%</div> <div>6%</div> </div> </div>
1	C	427	<div> <div>3%</div> <div> <div></div> <div>91%</div> <div>7%</div> </div> </div>
1	D	427	<div> <div>2%</div> <div> <div></div> <div>91%</div> <div>9%</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
2	NAG	A	500	-	-	-	X
3	NAG	A	4500	X	-	-	X
3	NAG	B	502	-	-	-	X
7	NAG	D	503	-	-	-	X

2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 16852 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 Alliin lyase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	425	Total	C	N	O	S	0	1	0
			3436	2190	577	646	23			
1	B	427	Total	C	N	O	S	0	0	0
			3447	2197	582	645	23			
1	C	425	Total	C	N	O	S	0	1	0
			3445	2200	579	643	23			
1	D	427	Total	C	N	O	S	0	0	0
			3447	2197	582	645	23			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	176	ASP	ASN	SEE REMARK 999	UNP Q01594
B	176	ASP	ASN	SEE REMARK 999	UNP Q01594
C	176	ASP	ASN	SEE REMARK 999	UNP Q01594
D	176	ASP	ASN	SEE REMARK 999	UNP Q01594

- Molecule 2 is a polymer of unknown type called SUGAR (6-MER).

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

- Molecule 3 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: C₈H₁₅NO₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	B	1	Total	C	N	O	0	0
			14	8	1	5		
3	B	1	Total	C	N	O	0	0
			14	8	1	5		
3	B	1	Total	C	N	O	0	0
			14	8	1	5		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	C	1	Total	C	N	O	0	0
			14	8	1	5		
3	C	1	Total	C	N	O	0	0
			14	8	1	5		

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

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	B	2	Total	C	N	O	0	0
			24	14	1	9		

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

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	C	4	Total	C	N	O	0	0
			49	28	2	19		

- Molecule 6 is a polymer of unknown type called SUGAR (3-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	D	3	Total	C	N	O	0	0
			38	22	2	14		

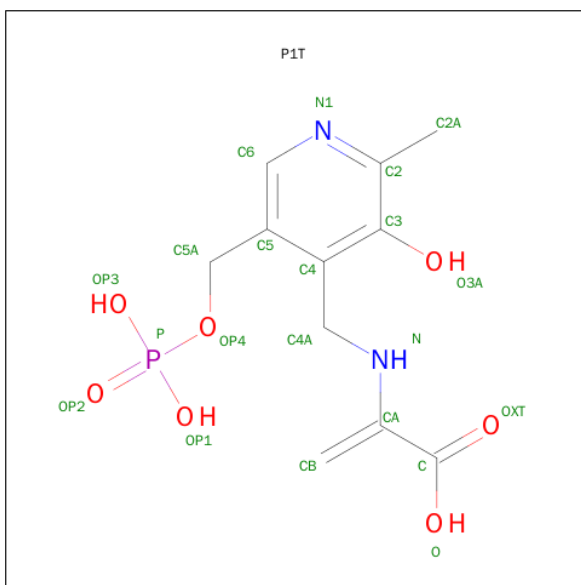
- Molecule 7 is a polymer of unknown type called SUGAR (2-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	D	2	Total	C	N	O	0	0
			28	16	2	10		

- Molecule 8 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	B	1	Total	Cl	0	0
			1	1		
8	A	1	Total	Cl	0	0
			1	1		
8	D	1	Total	Cl	0	0
			1	1		
8	C	1	Total	Cl	0	0
			1	1		

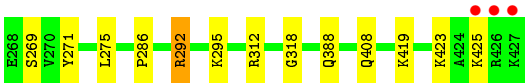
- Molecule 9 is 2-[(3-HYDROXY-2-METHYL-5-[(PHOSPHONOOXY)METHYL]PYRIDIN-4-YL)METHYL]AMINO]ACRYLIC ACID (three-letter code: P1T) (formula: C₁₁H₁₅N₂O₇P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
9	A	1	Total	C	N	O	P	0	0
			21	11	2	7	1		
9	B	1	Total	C	N	O	P	0	0
			21	11	2	7	1		
9	C	1	Total	C	N	O	P	0	0
			21	11	2	7	1		
9	D	1	Total	C	N	O	P	0	0
			21	11	2	7	1		

- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	752	Total	O	0	0
			752	752		
10	B	685	Total	O	0	0
			685	685		
10	C	774	Total	O	0	0
			774	774		
10	D	458	Total	O	0	2
			458	458		



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	67.67Å 126.89Å 102.66Å 90.00° 97.30° 90.00°	Depositor
Resolution (Å)	46.10 – 1.40 46.11 – 1.40	Depositor EDS
% Data completeness (in resolution range)	96.5 (46.10-1.40) 96.4 (46.11-1.40)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.50 (at 1.39Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.167 , 0.205 0.177 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	16.1	Xtriage
Anisotropy	0.132	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 52.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	1 of 327315 reflections (0.000%)	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	16852	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.08% 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, BMA, NAG, CL, P1T, FUC

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	0.87	2/3525 (0.1%)	0.86	2/4772 (0.0%)
1	B	0.87	2/3536 (0.1%)	0.87	2/4785 (0.0%)
1	C	0.87	3/3543 (0.1%)	0.84	2/4793 (0.0%)
1	D	0.87	2/3536 (0.1%)	0.84	1/4785 (0.0%)
All	All	0.87	9/14140 (0.1%)	0.85	7/19135 (0.0%)

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	C	0	1

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	224	TYR	CE2-CZ	-6.82	1.29	1.38
1	C	277	TYR	CE1-CZ	-6.80	1.29	1.38
1	A	277	TYR	CE1-CZ	-6.28	1.30	1.38
1	C	277	TYR	CD2-CE2	5.96	1.48	1.39
1	A	288	GLU	CB-CG	-5.71	1.41	1.52
1	D	292	ARG	CG-CD	-5.56	1.38	1.51
1	B	277	TYR	CE2-CZ	-5.39	1.31	1.38
1	C	288	GLU	CB-CG	-5.19	1.42	1.52
1	B	352	TYR	CG-CD2	5.01	1.45	1.39

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	259	ARG	NE-CZ-NH2	7.51	124.06	120.30
1	B	89	ARG	NE-CZ-NH2	-7.07	116.77	120.30
1	C	89	ARG	NE-CZ-NH1	6.46	123.53	120.30
1	C	213	ARG	NE-CZ-NH2	-5.90	117.35	120.30
1	A	124	ARG	NE-CZ-NH2	-5.73	117.43	120.30
1	A	89	ARG	NE-CZ-NH2	-5.62	117.49	120.30
1	D	259	ARG	NE-CZ-NH1	-5.54	117.53	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	424	ALA	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	3436	0	3325	28	0
1	B	3447	0	3347	19	0
1	C	3445	0	3360	40	0
1	D	3447	0	3347	31	0
2	A	69	0	59	2	0
3	A	42	0	39	1	0
3	B	42	0	39	4	0
3	C	28	0	26	1	0
4	B	24	0	22	0	0
5	C	49	0	43	1	0
6	D	38	0	34	0	0
7	D	28	0	25	0	0
8	A	1	0	0	0	0
8	B	1	0	0	0	0
8	C	1	0	0	0	0
8	D	1	0	0	0	0
9	A	21	0	11	2	0
9	B	21	0	11	3	0
9	C	21	0	11	0	0
9	D	21	0	12	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
10	A	752	0	0	14	1
10	B	685	0	0	7	2
10	C	774	0	0	17	1
10	D	458	0	0	11	0
All	All	16852	0	13711	123	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 (123) 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:D:41:CYS:SG	10:D:6250:HOH:O	1.94	1.21
1:C:277:TYR:CE2	10:C:6486:HOH:O	1.98	1.12
1:C:277:TYR:CZ	10:C:6486:HOH:O	2.02	1.09
1:B:328:ASN:HD21	3:B:502:NAG:C1	1.78	0.95
1:B:137:HIS:HD2	10:B:6599:HOH:O	1.49	0.94
1:C:311:MET:CE	1:C:316:THR:HG22	2.03	0.87
1:A:137:HIS:HD2	10:A:6721:HOH:O	1.57	0.86
1:C:311:MET:HE1	1:C:316:THR:HG22	1.56	0.84
1:B:45:TYR:CE1	1:B:53:LYS:HD2	2.18	0.78
1:C:194:ASN:HB3	10:C:6468:HOH:O	1.84	0.78
1:D:190:VAL:HG23	1:D:191:ASN:HD22	1.49	0.77
3:B:3500:NAG:O3	10:B:6270:HOH:O	2.01	0.77
1:D:123:ASP:O	10:D:6449:HOH:O	2.01	0.76
1:B:55:GLN:O	10:B:6653:HOH:O	2.05	0.74
1:C:137:HIS:HE1	1:C:177:LYS:NZ	1.84	0.74
1:D:312:ARG:HD2	10:D:6417:HOH:O	1.88	0.73
10:C:6631:HOH:O	1:D:54:ILE:HD13	1.91	0.71
1:B:45:TYR:CZ	1:B:53:LYS:HD2	2.27	0.69
1:D:190:VAL:HG23	1:D:191:ASN:ND2	2.07	0.68
1:D:312:ARG:CD	10:D:6417:HOH:O	2.41	0.68
1:C:232:HIS:CG	10:C:6387:HOH:O	2.46	0.68
1:C:343:GLU:OE1	10:C:6286:HOH:O	2.12	0.67
1:C:141:ILE:HD11	1:C:177:LYS:HZ2	1.58	0.67
1:A:311:MET:HE1	1:A:320:LYS:HG3	1.76	0.66
1:C:19:ASN:CG	10:C:6742:HOH:O	2.33	0.65
1:C:137:HIS:HE1	1:C:177:LYS:HZ2	1.42	0.65
1:B:328:ASN:ND2	3:B:502:NAG:C1	2.57	0.65
1:C:317:PHE:CZ	1:C:321:LYS:HD2	2.32	0.65
1:C:317:PHE:CE1	1:C:321:LYS:HD2	2.31	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:49:ASP:OD2	10:A:6477:HOH:O	2.15	0.63
1:C:294:LEU:HD11	1:C:298:LYS:HE2	1.80	0.62
1:C:311:MET:HE1	1:C:320:LYS:HD2	1.82	0.62
1:A:311:MET:CE	1:A:320:LYS:HG3	2.29	0.61
1:A:170:GLU:HG2	10:A:6714:HOH:O	2.00	0.61
1:C:311:MET:HE2	1:C:311:MET:HA	1.83	0.61
1:A:419:LYS:HE3	10:A:6410:HOH:O	1.99	0.61
1:B:53:LYS:HE2	10:B:6424:HOH:O	2.01	0.60
1:D:232:HIS:ND1	1:D:318:GLY:HA3	2.16	0.60
1:C:311:MET:HE2	1:C:316:THR:HG22	1.84	0.60
1:C:425:LYS:HD2	10:C:6749:HOH:O	2.02	0.59
1:C:337:ASP:HB2	10:C:6329:HOH:O	2.04	0.57
1:B:45:TYR:CE1	1:B:53:LYS:CD	2.85	0.57
1:C:232:HIS:CE1	10:C:6387:HOH:O	2.58	0.57
1:C:232:HIS:ND1	1:C:318:GLY:HA3	2.19	0.57
1:A:311:MET:HE1	1:A:316:THR:HG22	1.86	0.57
1:C:15:VAL:HG21	1:C:28:LEU:HD23	1.88	0.56
1:A:212:LEU:HD12	1:A:237:LYS:HE2	1.87	0.56
1:A:106:GLU:HG3	1:A:126:ILE:HD12	1.88	0.55
1:A:232:HIS:ND1	1:A:318:GLY:HA3	2.21	0.55
1:D:408:GLN:NE2	10:D:6393:HOH:O	2.38	0.55
1:A:176:ASP:CG	10:A:6408:HOH:O	2.44	0.55
1:C:137:HIS:CE1	1:C:177:LYS:HZ2	2.25	0.55
1:A:178:LYS:HG3	10:B:6636:HOH:O	2.07	0.54
1:B:77:LYS:NZ	10:B:6682:HOH:O	2.19	0.54
1:A:74:LYS:NZ	10:A:6326:HOH:O	2.40	0.54
1:A:74:LYS:NZ	10:A:6374:HOH:O	2.38	0.54
1:C:194:ASN:HB3	10:C:6766:HOH:O	2.09	0.53
1:D:419:LYS:O	1:D:423:LYS:HG2	2.09	0.52
1:B:328:ASN:HD21	3:B:502:NAG:C2	2.21	0.52
1:B:311:MET:SD	1:B:320:LYS:HD3	2.50	0.52
1:A:137:HIS:HE1	10:A:6720:HOH:O	1.93	0.51
10:C:6556:HOH:O	1:D:295:LYS:HD3	2.12	0.50
1:D:232:HIS:CG	10:D:6127:HOH:O	2.64	0.50
1:D:46:THR:OG1	10:D:6311:HOH:O	2.20	0.49
1:C:191:ASN:HA	10:C:6231:HOH:O	2.12	0.48
1:A:137:HIS:CE1	10:A:6720:HOH:O	2.65	0.48
1:C:311:MET:CE	1:C:320:LYS:HD2	2.43	0.48
10:A:6745:HOH:O	1:D:52:GLU:HG3	2.14	0.48
1:A:15:VAL:HG21	1:A:28:LEU:HD23	1.94	0.48
1:C:141:ILE:HD12	1:C:277:TYR:CE1	2.48	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:237:LYS:NZ	1:A:348:GLU:OE2	2.46	0.47
1:A:124:ARG:NH2	10:A:6286:HOH:O	2.47	0.47
1:C:317:PHE:CE1	1:C:321:LYS:CD	2.97	0.47
1:B:251:LYS:HZ3	9:B:6002:P1T:H4A2	1.79	0.46
1:A:19:ASN:ND2	1:A:21:SER:H	2.14	0.46
1:B:332:LEU:HD21	1:B:419:LYS:HG3	1.97	0.46
1:A:143:LEU:HD22	1:A:221:LYS:HD3	1.98	0.46
1:C:418:LEU:O	1:C:422:VAL:HG23	2.16	0.46
1:D:74:LYS:O	1:D:77:LYS:HG3	2.16	0.46
1:D:251:LYS:HZ3	9:D:6004:P1T:H4A2	1.81	0.46
1:D:388:GLN:NE2	10:D:6412:HOH:O	2.48	0.46
1:D:225:ASP:OD2	9:D:6004:P1T:N1	2.50	0.45
1:D:125:TYR:HB3	1:D:271:TYR:CD1	2.51	0.45
1:A:424:ALA:O	1:A:425:LYS:HB2	2.16	0.45
1:C:286:PRO:HG2	1:D:286:PRO:HG2	1.98	0.45
1:C:11:GLU:O	1:C:15:VAL:HG23	2.16	0.45
1:A:251:LYS:HZ3	9:A:6001:P1T:H4A2	1.81	0.45
1:C:370:TRP:CH2	1:C:425:LYS:HD3	2.52	0.45
1:B:199:ILE:HA	1:B:221:LYS:O	2.17	0.44
1:B:53:LYS:CE	10:B:6424:HOH:O	2.63	0.44
2:A:505:BMA:O2	10:A:6718:HOH:O	2.16	0.44
1:A:212:LEU:HD12	1:A:237:LYS:CE	2.47	0.44
5:C:500:NAG:H83	1:D:174:TYR:CE1	2.53	0.44
1:C:141:ILE:HD11	1:C:177:LYS:NZ	2.27	0.44
3:C:504:NAG:HO4	3:C:505:NAG:C1	2.31	0.44
1:C:298:LYS:CD	10:C:6507:HOH:O	2.65	0.44
1:D:232:HIS:CE1	10:D:6127:HOH:O	2.71	0.43
1:A:218:LYS:HB2	3:A:4500:NAG:H81	1.99	0.43
1:B:225:ASP:OD2	9:B:6002:P1T:N1	2.52	0.43
1:A:73:TRP:CH2	1:A:296:VAL:HG13	2.53	0.43
1:A:2:MET:N	10:A:6295:HOH:O	2.52	0.43
1:C:137:HIS:CE1	1:C:177:LYS:NZ	2.75	0.43
1:D:2:MET:CE	10:D:6307:HOH:O	2.66	0.42
1:B:110:LYS:O	1:B:114:GLU:HG2	2.20	0.42
1:D:312:ARG:NH2	10:D:6354:HOH:O	2.52	0.42
9:A:6001:P1T:H4A1	9:A:6001:P1T:H5A2	1.85	0.42
1:C:277:TYR:OH	10:C:6486:HOH:O	1.85	0.41
1:D:271:TYR:CZ	1:D:275:LEU:HD11	2.56	0.41
9:D:6004:P1T:H5A2	9:D:6004:P1T:H4A1	1.78	0.41
1:D:251:LYS:NZ	9:D:6004:P1T:HB1	2.36	0.41
1:D:267:ASP:OD2	1:D:269:SER:OG	2.33	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:157:LYS:NZ	1:B:197:GLN:OE1	2.49	0.41
9:B:6002:P1T:H5A2	9:B:6002:P1T:H4A1	1.89	0.41
1:A:199:ILE:HA	1:A:221:LYS:O	2.20	0.41
1:D:168:PHE:CE1	1:D:201:MET:HE2	2.55	0.41
1:C:311:MET:CE	1:C:316:THR:CG2	2.88	0.41
1:C:321:LYS:HG3	10:C:6509:HOH:O	2.21	0.41
2:A:505:BMA:H2	10:A:6718:HOH:O	2.20	0.41
1:C:114:GLU:HG3	10:C:6313:HOH:O	2.20	0.41
1:D:83:LEU:HD13	1:D:83:LEU:C	2.41	0.40
1:C:288:GLU:OE2	1:D:292:ARG:HD3	2.21	0.40
1:D:15:VAL:HG21	1:D:28:LEU:HD23	2.02	0.40
1:B:332:LEU:CD2	1:B:419:LYS:HG3	2.51	0.40

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)
10:B:6436:HOH:O	10:C:6361:HOH:O[1_455]	1.99	0.21
10:A:6664:HOH:O	10:B:6132:HOH:O[2_645]	2.11	0.09

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	424/427 (99%)	411 (97%)	13 (3%)	0	100	100
1	B	425/427 (100%)	413 (97%)	12 (3%)	0	100	100
1	C	426/427 (100%)	414 (97%)	12 (3%)	0	100	100
1	D	425/427 (100%)	412 (97%)	13 (3%)	0	100	100
All	All	1700/1708 (100%)	1650 (97%)	50 (3%)	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	374/375 (100%)	369 (99%)	5 (1%)	76	48
1	B	375/375 (100%)	366 (98%)	9 (2%)	57	19
1	C	376/375 (100%)	369 (98%)	7 (2%)	65	29
1	D	375/375 (100%)	371 (99%)	4 (1%)	80	56
All	All	1500/1500 (100%)	1475 (98%)	25 (2%)	70	35

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	LYS
1	A	25	ARG
1	A	176	ASP
1	A	268[A]	GLU
1	A	268[B]	GLU
1	B	1	LYS
1	B	6	MET
1	B	25	ARG
1	B	53	LYS
1	B	55	GLN
1	B	232	HIS
1	B	425	LYS
1	B	426	ARG
1	B	427	LYS
1	C	7	LYS
1	C	25	ARG
1	C	152	ASP
1	C	191	ASN
1	C	194	ASN
1	C	281	ASN
1	C	425	LYS
1	D	25	ARG
1	D	157	LYS
1	D	178	LYS
1	D	425	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	19	ASN
1	A	214	HIS
1	A	381	GLN
1	A	386	ASN
1	B	137	HIS
1	B	214	HIS
1	B	328	ASN
1	B	381	GLN
1	B	386	ASN
1	B	408	GLN
1	C	137	HIS
1	C	214	HIS
1	D	137	HIS
1	D	191	ASN
1	D	307	GLN
1	D	408	GLN

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](#)

17 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	500	1,2	14,14,15	0.64	0	15,19,21	1.34	2 (13%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	FUC	A	501	2	10,10,11	0.54	0	14,14,16	1.61	4 (28%)
2	NAG	A	502	2	14,14,15	0.37	0	15,19,21	1.11	2 (13%)
2	BMA	A	503	2	11,11,12	0.57	0	14,15,17	1.33	1 (7%)
2	XYP	A	504	2	9,9,10	0.80	0	12,12,14	0.82	0
2	BMA	A	505	2	11,11,12	0.90	1 (9%)	14,15,17	5.71	5 (35%)
4	NAG	B	500	1,4	14,14,15	0.47	0	15,19,21	2.08	1 (6%)
4	FUC	B	501	4	10,10,11	0.35	0	14,14,16	0.69	0
5	NAG	C	500	1,5	14,14,15	0.68	0	15,19,21	0.74	0
5	FUC	C	501	5	10,10,11	0.54	0	14,14,16	1.50	3 (21%)
5	NAG	C	502	5	14,14,15	0.66	0	15,19,21	1.56	4 (26%)
5	BMA	C	503	5	11,11,12	0.59	0	14,15,17	1.11	1 (7%)
6	NAG	D	500	1,6	14,14,15	0.51	0	15,19,21	1.40	1 (6%)
6	FUC	D	501	6	10,10,11	0.42	0	14,14,16	1.13	1 (7%)
6	NAG	D	502	6	14,14,15	0.67	1 (7%)	15,19,21	0.83	0
7	NAG	D	503	1,7	14,14,15	0.67	0	15,19,21	1.11	2 (13%)
7	NAG	D	504	7	14,14,15	0.58	0	15,19,21	1.25	2 (13%)

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	500	1,2	-	0/6/23/26	0/1/1/1
2	FUC	A	501	2	-	0/0/17/20	0/1/1/1
2	NAG	A	502	2	-	0/6/23/26	0/1/1/1
2	BMA	A	503	2	-	0/2/19/22	0/1/1/1
2	XYP	A	504	2	-	0/0/14/17	0/1/1/1
2	BMA	A	505	2	-	0/2/19/22	0/1/1/1
4	NAG	B	500	1,4	-	0/6/23/26	0/1/1/1
4	FUC	B	501	4	-	0/0/17/20	0/1/1/1
5	NAG	C	500	1,5	-	0/6/23/26	0/1/1/1
5	FUC	C	501	5	-	0/0/17/20	0/1/1/1
5	NAG	C	502	5	-	0/6/23/26	0/1/1/1
5	BMA	C	503	5	-	0/2/19/22	0/1/1/1
6	NAG	D	500	1,6	-	0/6/23/26	0/1/1/1
6	FUC	D	501	6	-	0/0/17/20	0/1/1/1
6	NAG	D	502	6	-	0/6/23/26	0/1/1/1
7	NAG	D	503	1,7	-	0/6/23/26	0/1/1/1
7	NAG	D	504	7	-	0/6/23/26	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	505	BMA	O5-C1	-2.48	1.39	1.43
6	D	502	NAG	O5-C1	-2.11	1.40	1.43

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	505	BMA	C1-C2-C3	-15.50	91.20	109.54
2	A	505	BMA	C1-O5-C5	-13.37	95.28	112.25
2	A	503	BMA	O3-C3-C4	-3.86	101.64	110.34
2	A	500	NAG	O4-C4-C3	-2.70	104.26	110.34
2	A	502	NAG	O7-C7-C8	-2.68	117.15	122.06
2	A	500	NAG	C8-C7-N2	-2.45	111.42	116.11
2	A	501	FUC	C2-C3-C4	-2.43	106.91	111.04
5	C	503	BMA	O5-C1-C2	-2.31	107.11	110.86
7	D	503	NAG	C2-N2-C7	-2.23	120.18	123.04
5	C	502	NAG	O7-C7-C8	-2.12	118.17	122.06
5	C	501	FUC	C2-C3-C4	-2.01	107.62	111.04
5	C	501	FUC	C3-C4-C5	2.04	113.15	109.72
2	A	505	BMA	O5-C5-C6	2.04	111.76	107.35
2	A	502	NAG	C8-C7-N2	2.11	120.15	116.11
7	D	504	NAG	C3-C4-C5	2.22	114.06	110.20
7	D	503	NAG	C1-O5-C5	2.23	115.07	112.25
2	A	501	FUC	C3-C4-C5	2.34	113.66	109.72
5	C	502	NAG	C2-N2-C7	2.44	126.18	123.04
2	A	505	BMA	C3-C4-C5	2.47	114.51	110.20
5	C	502	NAG	C8-C7-N2	2.61	121.09	116.11
7	D	504	NAG	C4-C3-C2	2.61	115.28	111.23
2	A	501	FUC	C1-O5-C5	2.77	116.66	112.38
2	A	501	FUC	O5-C5-C4	2.91	114.58	109.53
5	C	502	NAG	C4-C3-C2	3.05	115.97	111.23
5	C	501	FUC	O5-C5-C6	3.33	111.64	106.13
6	D	501	FUC	O5-C5-C6	3.58	112.04	106.13
6	D	500	NAG	C1-O5-C5	3.92	117.22	112.25
2	A	505	BMA	O2-C2-C1	4.24	117.72	109.21
4	B	500	NAG	C1-O5-C5	7.22	121.41	112.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	505	BMA	2	0
5	C	500	NAG	1	0

5.6 Ligand geometry

Of 16 ligands modelled in this entry, 4 are monoatomic - leaving 12 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$
3	NAG	A	4500	1	14,14,15	0.38	0	15,19,21	1.82	1 (6%)
3	NAG	A	506	1	14,14,15	0.43	0	15,19,21	0.81	0
3	NAG	A	507	-	14,14,15	0.45	0	15,19,21	1.42	2 (13%)
9	P1T	A	6001	-	19,21,21	2.41	4 (21%)	22,30,30	2.38	9 (40%)
3	NAG	B	3500	1	14,14,15	0.54	0	15,19,21	0.84	0
3	NAG	B	502	-	14,14,15	0.52	0	15,19,21	1.51	2 (13%)
3	NAG	B	503	-	14,14,15	0.46	0	15,19,21	0.87	0
9	P1T	B	6002	-	19,21,21	2.61	5 (26%)	22,30,30	2.34	5 (22%)
3	NAG	C	504	1	14,14,15	0.50	0	15,19,21	0.86	0
3	NAG	C	505	-	14,14,15	0.67	0	15,19,21	1.25	1 (6%)
9	P1T	C	6003	-	19,21,21	2.80	5 (26%)	22,30,30	2.03	5 (22%)
9	P1T	D	6004	-	19,21,21	2.56	3 (15%)	22,30,30	2.18	8 (36%)

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	A	4500	1	1/1/5/7	0/6/23/26	0/1/1/1
3	NAG	A	506	1	-	0/6/23/26	0/1/1/1
3	NAG	A	507	-	-	0/6/23/26	0/1/1/1
9	P1T	A	6001	-	-	0/10/15/15	0/1/1/1
3	NAG	B	3500	1	-	0/6/23/26	0/1/1/1
3	NAG	B	502	-	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	B	503	-	-	0/6/23/26	0/1/1/1
9	P1T	B	6002	-	-	0/10/15/15	0/1/1/1
3	NAG	C	504	1	-	0/6/23/26	0/1/1/1
3	NAG	C	505	-	-	0/6/23/26	0/1/1/1
9	P1T	C	6003	-	-	0/10/15/15	0/1/1/1
9	P1T	D	6004	-	-	0/10/15/15	0/1/1/1

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	C	6003	P1T	C4A-N	-8.42	1.29	1.46
9	A	6001	P1T	C4A-N	-8.41	1.29	1.46
9	D	6004	P1T	C4A-N	-8.15	1.29	1.46
9	B	6002	P1T	C4A-N	-7.80	1.30	1.46
9	B	6002	P1T	C-CA	-3.49	1.46	1.52
9	D	6004	P1T	C-CA	-3.40	1.46	1.52
9	C	6003	P1T	C-CA	-3.30	1.46	1.52
9	B	6002	P1T	P-OP1	-2.80	1.44	1.54
9	A	6001	P1T	C-CA	-2.69	1.47	1.52
9	C	6003	P1T	C4A-C4	-2.28	1.49	1.51
9	A	6001	P1T	P-OP3	-2.20	1.46	1.54
9	B	6002	P1T	C3-C4	-2.15	1.36	1.40
9	C	6003	P1T	C6-C5	2.23	1.42	1.37
9	A	6001	P1T	C3-C2	3.86	1.43	1.40
9	D	6004	P1T	C3-C2	4.98	1.44	1.40
9	B	6002	P1T	C3-C2	5.73	1.44	1.40
9	C	6003	P1T	C3-C2	6.17	1.45	1.40

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	A	6001	P1T	C3-C4-C5	-2.72	115.96	118.82
9	D	6004	P1T	C2A-C2-C3	-2.58	117.92	121.04
9	B	6002	P1T	C2A-C2-C3	-2.47	118.06	121.04
9	D	6004	P1T	OP3-P-OP4	-2.43	99.58	106.56
9	A	6001	P1T	OP3-P-OP4	-2.13	100.44	106.56
9	A	6001	P1T	C3-C2-N1	-2.12	117.68	120.61
9	C	6003	P1T	OP3-P-OP4	-2.02	100.76	106.56
9	A	6001	P1T	C2A-C2-N1	2.27	122.98	117.95
9	D	6004	P1T	C3-C2-N1	2.32	123.82	120.61
9	D	6004	P1T	O3A-C3-C2	2.42	121.86	117.66
3	A	507	NAG	C3-C4-C5	2.43	114.44	110.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	A	6001	P1T	OP1-P-OP4	2.76	114.52	106.56
3	B	502	NAG	C3-C4-C5	2.80	115.08	110.20
9	A	6001	P1T	C6-C5-C4	2.89	120.25	118.09
9	C	6003	P1T	C6-C5-C4	3.17	120.45	118.09
9	D	6004	P1T	C4-C4A-N	3.54	119.13	111.49
9	A	6001	P1T	C4-C4A-N	3.73	119.54	111.49
9	D	6004	P1T	OP4-C5A-C5	3.77	115.22	108.99
9	B	6002	P1T	C6-C5-C4	3.94	121.03	118.09
9	B	6002	P1T	C4-C4A-N	4.05	120.22	111.49
9	C	6003	P1T	C4-C4A-N	4.06	120.24	111.49
3	C	505	NAG	C3-C4-C5	4.10	117.34	110.20
9	D	6004	P1T	C4A-C4-C5	4.13	123.39	119.71
9	B	6002	P1T	OP4-C5A-C5	4.21	115.96	108.99
9	D	6004	P1T	C6-C5-C4	4.36	121.35	118.09
3	A	507	NAG	C1-O5-C5	4.42	117.86	112.25
9	C	6003	P1T	C4A-C4-C5	4.51	123.72	119.71
3	B	502	NAG	C1-O5-C5	4.61	118.10	112.25
9	C	6003	P1T	OP4-C5A-C5	4.69	116.75	108.99
9	A	6001	P1T	OP4-C5A-C5	4.88	117.06	108.99
9	A	6001	P1T	C4A-C4-C5	5.55	124.65	119.71
9	B	6002	P1T	C4A-C4-C5	6.59	125.59	119.71
3	A	4500	NAG	C1-O5-C5	6.62	120.65	112.25

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	A	4500	NAG	C1

There are no torsion outliers.

There are no ring outliers.

8 monomers are involved in 15 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	4500	NAG	1	0
9	A	6001	P1T	2	0
3	B	3500	NAG	1	0
3	B	502	NAG	3	0
9	B	6002	P1T	3	0
3	C	504	NAG	1	0
3	C	505	NAG	1	0
9	D	6004	P1T	4	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 ⓘ

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	425/427 (99%)	-0.16	7 (1%) 74 73	10, 19, 37, 53	1 (0%)
1	B	427/427 (100%)	-0.15	8 (1%) 70 68	11, 18, 35, 65	1 (0%)
1	C	425/427 (99%)	-0.17	11 (2%) 59 57	10, 18, 33, 52	0
1	D	427/427 (100%)	-0.14	10 (2%) 64 62	11, 18, 36, 69	0
All	All	1704/1708 (99%)	-0.15	36 (2%) 67 64	10, 18, 36, 69	2 (0%)

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	427	LYS	7.6
1	D	427	LYS	7.5
1	A	1	LYS	6.9
1	D	152	ASP	5.9
1	B	425	LYS	4.8
1	C	416	TYR	4.3
1	C	1	LYS	4.2
1	B	426	ARG	4.0
1	B	152	ASP	3.6
1	A	22	GLU	3.6
1	B	148	THR	3.3
1	D	426	ARG	3.2
1	C	56	GLY	3.2
1	A	416	TYR	3.2
1	D	34	GLU	3.1
1	A	48	PRO	2.9
1	B	149	ALA	2.9
1	A	51	SER	2.9
1	C	191	ASN	2.8
1	C	55	GLN	2.7
1	D	425	LYS	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	55	GLN	2.6
1	C	22	GLU	2.6
1	D	123	ASP	2.5
1	D	193	SER	2.5
1	C	370	TRP	2.4
1	B	150	THR	2.4
1	D	19	ASN	2.3
1	C	337	ASP	2.3
1	C	51	SER	2.2
1	C	423	LYS	2.2
1	B	22	GLU	2.2
1	C	371	GLU	2.2
1	A	424	ALA	2.2
1	D	1	LYS	2.1
1	D	21	SER	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(Å ²)	Q<0.9
7	NAG	D	503	14/15	0.82	0.24	4.85	31,38,54,61	0
2	NAG	A	500	14/15	0.76	0.16	2.00	26,32,36,36	0
5	NAG	C	500	14/15	0.93	0.08	0.86	24,27,33,34	0
5	FUC	C	501	10/11	0.92	0.09	0.18	24,29,32,33	0
7	NAG	D	504	14/15	0.46	0.44	-	68,72,75,76	0
6	FUC	D	501	10/11	0.92	0.17	-	36,41,50,52	0
6	NAG	D	500	14/15	0.86	0.14	-	28,33,38,39	0
5	NAG	C	502	14/15	0.92	0.14	-	30,36,51,60	0
2	BMA	A	505	11/12	0.91	0.14	-	20,23,33,43	0
4	FUC	B	501	10/11	0.85	0.20	-	46,51,55,56	0
2	FUC	A	501	10/11	0.76	0.19	-	39,44,50,50	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
6	NAG	D	502	14/15	0.80	0.23	-	40,46,48,54	0
2	NAG	A	502	14/15	0.87	0.15	-	26,31,47,50	0
5	BMA	C	503	11/12	0.37	0.23	-	69,71,74,76	0
2	XYP	A	504	9/10	0.94	0.15	-	20,21,22,24	0
2	BMA	A	503	11/12	0.95	0.14	-	18,23,31,39	0
4	NAG	B	500	14/15	0.84	0.17	-	32,40,48,51	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	NAG	A	4500	14/15	0.48	0.36	13.64	64,70,79,80	0
3	NAG	B	502	14/15	0.67	0.21	3.09	48,59,69,69	0
3	NAG	C	504	14/15	0.91	0.14	1.80	27,33,44,44	0
3	NAG	A	506	14/15	0.88	0.15	1.62	32,42,55,61	0
9	P1T	D	6004	21/21	0.96	0.09	0.75	12,17,25,27	0
9	P1T	A	6001	21/21	0.97	0.09	0.65	13,17,29,29	0
3	NAG	B	3500	14/15	0.85	0.21	0.32	42,52,55,59	0
8	CL	C	5003	1/1	1.00	0.07	0.19	13,13,13,13	0
9	P1T	B	6002	21/21	0.97	0.09	0.02	12,17,24,25	0
9	P1T	C	6003	21/21	0.97	0.08	-0.29	13,17,25,28	0
8	CL	B	5002	1/1	0.99	0.07	-0.56	16,16,16,16	0
8	CL	A	5001	1/1	1.00	0.06	-0.97	12,12,12,12	0
8	CL	D	5004	1/1	0.99	0.06	-1.04	15,15,15,15	0
3	NAG	A	507	14/15	0.42	0.41	-	104,110,112,113	0
3	NAG	C	505	14/15	0.54	0.32	-	73,82,84,84	0
3	NAG	B	503	14/15	-0.04	0.60	-	128,132,132,133	0

6.5 Other polymers [i](#)

There are no such residues in this entry.