



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 8, 2024 – 01:47 am GMT

PDB ID : 6EZX
Title : Crystal structure of GH20 Exo beta-N-Acetylglucosaminidase from *Vibrio harveyi* in complex with N-acetylglucosamine
Authors : Porfetye, A.T.; Meekrathok, P.; Burger, M.; Vetter, I.R.; Suginta, W.
Deposited on : 2017-11-16
Resolution : 2.50 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4.02b-467
Mogul : 1.8.4, CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.36
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36

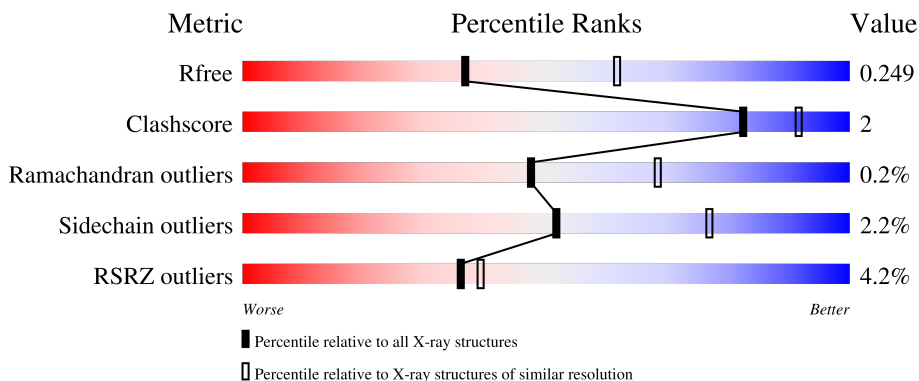
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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}	130704	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	652	 4% 89% 9%
1	B	652	 4% 92% 6%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	MLI	A	702	-	-	X	-

2 Entry composition i

There are 4 unique types of molecules in this entry. The entry contains 11254 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 Beta-N-acetylglucosaminidase Nag2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	639	5161	3285	880	974	22	0	3	0
1	B	639	5164	3287	882	973	22	0	3	0

There are 20 discrepancies between the modelled and reference sequences:

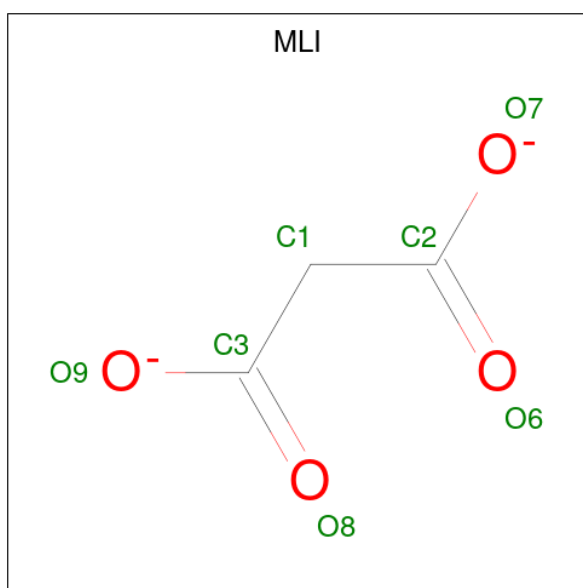
Chain	Residue	Modelled	Actual	Comment	Reference
A	643	ARG	-	expression tag	UNP D9ISE0
A	644	SER	-	expression tag	UNP D9ISE0
A	645	ARG	-	expression tag	UNP D9ISE0
A	646	SER	-	expression tag	UNP D9ISE0
A	647	HIS	-	expression tag	UNP D9ISE0
A	648	HIS	-	expression tag	UNP D9ISE0
A	649	HIS	-	expression tag	UNP D9ISE0
A	650	HIS	-	expression tag	UNP D9ISE0
A	651	HIS	-	expression tag	UNP D9ISE0
A	652	HIS	-	expression tag	UNP D9ISE0
B	643	ARG	-	expression tag	UNP D9ISE0
B	644	SER	-	expression tag	UNP D9ISE0
B	645	ARG	-	expression tag	UNP D9ISE0
B	646	SER	-	expression tag	UNP D9ISE0
B	647	HIS	-	expression tag	UNP D9ISE0
B	648	HIS	-	expression tag	UNP D9ISE0
B	649	HIS	-	expression tag	UNP D9ISE0
B	650	HIS	-	expression tag	UNP D9ISE0
B	651	HIS	-	expression tag	UNP D9ISE0
B	652	HIS	-	expression tag	UNP D9ISE0

- Molecule 2 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: C₈H₁₅NO₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	
			Total	C	N			O
2	A	1	15	8	1	6	0	0
2	B	1	15	8	1	6	0	0

- Molecule 3 is MALONATE ION (three-letter code: MLI) (formula: $C_3H_2O_4$).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	C O		
3	A	1	7	3 4	0	0
3	B	1	7	3 4	0	0

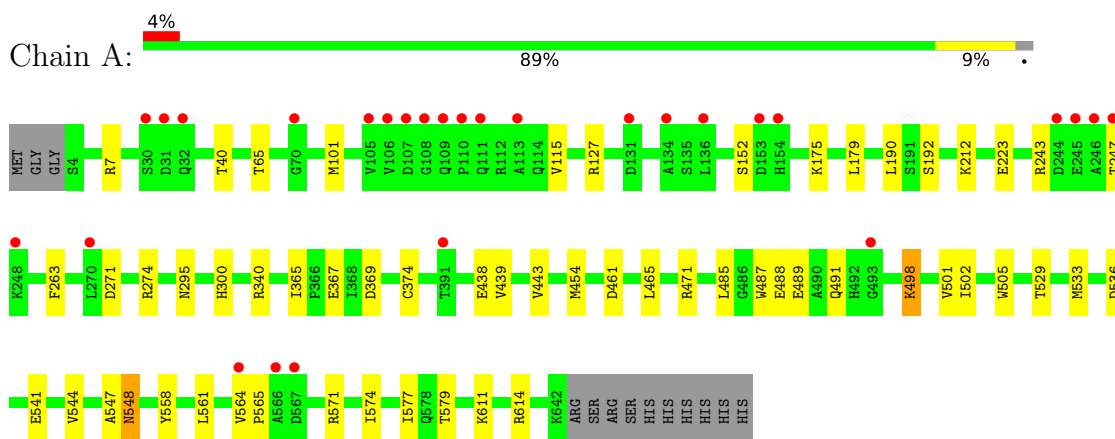
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	419	Total 419	O 419	0	0
4	B	466	Total 466	O 466	0	0

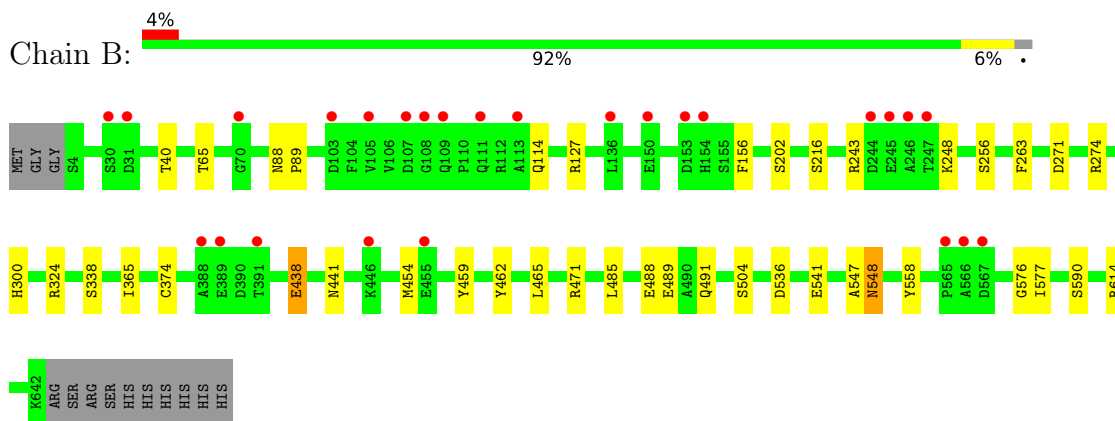
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: Beta-N-acetylglucosaminidase Nag2



- Molecule 1: Beta-N-acetylglucosaminidase Nag2



4 Data and refinement statistics i

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	91.25Å 129.63Å 100.00Å 90.00° 114.36° 90.00°	Depositor
Resolution (Å)	48.02 – 2.50 48.02 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.6 (48.02-2.50) 99.6 (48.02-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.29 (at 2.51Å)	Xtrriage
Refinement program	REFMAC 5.8.0103	Depositor
R, R_{free}	0.200 , 0.248 0.205 , 0.249	Depositor DCC
R_{free} test set	3648 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	33.1	Xtrriage
Anisotropy	0.210	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 39.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.006 for h,-k,-h-l	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	11254	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 20.48 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 8.6025e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NAG, MLI

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	0.47	0/5307	0.67	1/7216 (0.0%)
1	B	0.48	0/5307	0.67	1/7215 (0.0%)
All	All	0.47	0/10614	0.67	2/14431 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	340	ARG	NE-CZ-NH2	5.15	122.87	120.30
1	B	324	ARG	NE-CZ-NH1	5.00	122.80	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	5161	0	5000	30	0
1	B	5164	0	5006	20	0
2	A	15	0	15	0	0
2	B	15	0	15	0	0
3	A	7	0	2	3	0
3	B	7	0	2	0	0
4	A	419	0	0	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	466	0	0	4	0
All	All	11254	0	10040	50	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 2.

All (50) 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:B:40:THR:HG22	1:B:65:THR:HG22	1.70	0.74
1:A:536:ASP:HB3	1:A:541:GLU:HG3	1.84	0.58
1:A:485:LEU:HD22	1:A:501:VAL:HB	1.86	0.57
1:A:7:ARG:HD2	4:A:1074:HOH:O	2.05	0.56
1:A:243:ARG:NH1	4:A:802:HOH:O	2.30	0.56
1:B:489:GLU:HG2	4:B:1032:HOH:O	2.07	0.55
1:B:536:ASP:HB3	1:B:541:GLU:HG3	1.90	0.53
1:A:558:TYR:O	1:A:614:ARG:NH2	2.42	0.52
1:A:40:THR:HG22	1:A:65:THR:HG22	1.92	0.52
1:A:547:ALA:O	1:A:548:ASN:HB2	2.10	0.52
1:B:547:ALA:O	1:B:548:ASN:HB2	2.11	0.50
1:B:454:MET:HE1	1:B:462:TYR:CG	2.47	0.50
1:B:271:ASP:HA	1:B:300:HIS:HB3	1.94	0.49
1:B:558:TYR:O	1:B:614:ARG:NH2	2.47	0.48
1:B:256:SER:HB2	4:B:1102:HOH:O	2.14	0.47
1:A:488:GLU:HA	1:A:502:ILE:HG23	1.96	0.47
1:B:576:GLY:C	1:B:577:ILE:HD12	2.35	0.46
1:B:488:GLU:CG	1:B:504:SER:HA	2.46	0.45
1:B:454:MET:HG3	1:B:459:TYR:HB2	1.98	0.45
1:A:488:GLU:HA	1:A:502:ILE:CG2	2.46	0.45
1:A:101:MET:HA	1:A:115:VAL:HG23	1.99	0.45
1:B:114:GLN:HE21	1:B:338:SER:HB3	1.83	0.44
1:B:438[B]:GLU:HB2	1:B:489:GLU:HG3	1.99	0.44
1:B:127:ARG:HD2	4:B:1180:HOH:O	2.18	0.44
1:A:471:ARG:NH2	4:A:814:HOH:O	2.51	0.43
1:A:439:VAL:HB	1:A:443:VAL:HG11	2.00	0.43
1:A:365:ILE:HG21	1:A:485:LEU:HD12	2.00	0.43
1:A:295:ASN:HD22	3:A:702:MLI:C1	2.32	0.43
1:A:498:LYS:HA	1:A:498:LYS:HD3	1.81	0.43
1:B:471:ARG:NH2	4:B:805:HOH:O	2.46	0.43
1:A:561:LEU:HD11	1:A:574:ILE:HG21	2.00	0.43
1:A:263:PHE:CD2	1:A:577:ILE:HD13	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:295:ASN:ND2	3:A:702:MLI:H12	2.34	0.42
1:B:156:PHE:HB2	1:B:216:SER:O	2.20	0.42
1:B:263:PHE:CD2	1:B:577:ILE:HD13	2.54	0.42
1:A:565:PRO:O	1:A:571:ARG:NH1	2.51	0.42
1:B:365:ILE:HG21	1:B:485:LEU:HD12	2.02	0.42
1:A:40:THR:HG23	4:A:874:HOH:O	2.20	0.41
1:A:529:THR:HB	1:A:579:THR:HA	2.02	0.41
1:A:40:THR:CG2	1:A:65:THR:HG22	2.50	0.41
1:A:295:ASN:HD22	3:A:702:MLI:H12	1.85	0.41
1:A:367:GLU:OE2	1:A:369:ASP:OD1	2.39	0.41
1:A:271:ASP:HA	1:A:300:HIS:HB3	2.02	0.41
1:A:127:ARG:HD2	4:A:1165:HOH:O	2.21	0.41
1:A:179:LEU:HD22	1:A:190:LEU:HB2	2.03	0.41
1:B:88:ASN:HB3	1:B:89:PRO:CD	2.50	0.41
1:A:533:MET:HA	1:A:544:VAL:O	2.21	0.41
1:B:454:MET:HE1	1:B:462:TYR:HA	2.04	0.40
1:A:212:LYS:HB2	1:A:223:GLU:HB2	2.03	0.40
1:A:487:TRP:HB3	1:A:505:TRP:HZ3	1.86	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	A	640/652 (98%)	631 (99%)	8 (1%)	1 (0%)	47 68
1	B	640/652 (98%)	630 (98%)	9 (1%)	1 (0%)	47 68
All	All	1280/1304 (98%)	1261 (98%)	17 (1%)	2 (0%)	47 68

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	548	ASN
1	A	548	ASN

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	A	559/567 (99%)	543 (97%)	16 (3%)	42 69
1	B	559/567 (99%)	548 (98%)	11 (2%)	55 79
All	All	1118/1134 (99%)	1091 (98%)	27 (2%)	52 74

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	152	SER
1	A	175	LYS
1	A	192	SER
1	A	247	THR
1	A	274	ARG
1	A	374	CYS
1	A	438[A]	GLU
1	A	438[B]	GLU
1	A	454	MET
1	A	461	ASP
1	A	465	LEU
1	A	489	GLU
1	A	491	GLN
1	A	498	LYS
1	A	564	VAL
1	A	611	LYS
1	B	202	SER
1	B	243	ARG
1	B	248	LYS
1	B	274	ARG
1	B	374	CYS
1	B	438[A]	GLU

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Mol	Chain	Res	Type
1	B	438[B]	GLU
1	B	441	ASN
1	B	465	LEU
1	B	491	GLN
1	B	590	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	114	GLN
1	B	441	ASN

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](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

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$
2	NAG	A	701	-	15,15,15	0.80	1 (6%)	21,21,21	2.32	4 (19%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	MLI	B	702	-	6,6,6	1.08	0	7,7,7	0.73	0
2	NAG	B	701	-	15,15,15	0.61	0	21,21,21	1.25	4 (19%)
3	MLI	A	702	-	6,6,6	1.16	0	7,7,7	0.91	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
2	NAG	A	701	-	-	0/6/26/26	0/1/1/1
3	MLI	B	702	-	-	0/4/4/4	-
2	NAG	B	701	-	-	0/6/26/26	0/1/1/1
3	MLI	A	702	-	-	4/4/4/4	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	701	NAG	C1-C2	2.07	1.55	1.52

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	701	NAG	O5-C1-C2	-8.39	101.08	109.52
2	A	701	NAG	C1-C2-N2	-2.80	107.49	110.73
2	A	701	NAG	C1-C2-C3	2.77	114.32	110.54
2	B	701	NAG	C1-C2-N2	-2.52	107.81	110.73
2	B	701	NAG	C8-C7-N2	-2.27	112.26	116.10
2	B	701	NAG	O5-C1-C2	-2.25	107.26	109.52
2	B	701	NAG	C1-C2-C3	2.22	113.57	110.54
2	A	701	NAG	C8-C7-N2	-2.10	112.54	116.10

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	702	MLI	C3-C1-C2-O7
3	A	702	MLI	C3-C1-C2-O6
3	A	702	MLI	C2-C1-C3-O8
3	A	702	MLI	C2-C1-C3-O9

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	702	MLI	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

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	639/652 (98%)	0.18	28 (4%) 34 37	20, 34, 58, 97	0
1	B	639/652 (98%)	0.16	26 (4%) 37 40	20, 33, 58, 97	0
All	All	1278/1304 (98%)	0.17	54 (4%) 36 39	20, 34, 58, 97	0

All (54) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	246	ALA	9.2
1	A	107	ASP	6.0
1	B	245	GLU	5.6
1	A	108	GLY	5.3
1	B	246	ALA	5.2
1	B	389	GLU	4.2
1	B	108	GLY	4.0
1	A	245	GLU	3.9
1	A	30	SER	3.8
1	A	566	ALA	3.6
1	B	247	THR	3.5
1	A	109	GLN	3.4
1	B	153	ASP	3.4
1	B	109	GLN	3.3
1	A	136	LEU	3.2
1	B	107	ASP	3.2
1	B	566	ALA	3.0
1	A	110	PRO	3.0
1	A	105	VAL	2.9
1	A	154	HIS	2.9
1	A	153	ASP	2.9
1	B	31	ASP	2.8
1	A	113	ALA	2.8
1	A	31	ASP	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	30	SER	2.6
1	A	32	GLN	2.6
1	A	131	ASP	2.6
1	A	567	ASP	2.6
1	B	391	THR	2.5
1	A	106	VAL	2.5
1	A	70	GLY	2.5
1	A	111	GLN	2.5
1	B	113	ALA	2.5
1	A	248	LYS	2.5
1	B	244	ASP	2.4
1	B	105	VAL	2.4
1	B	455	GLU	2.3
1	B	565	PRO	2.3
1	A	270	LEU	2.3
1	B	70	GLY	2.2
1	B	446	LYS	2.2
1	B	150	GLU	2.2
1	A	134	ALA	2.2
1	A	493	GLY	2.2
1	B	111	GLN	2.2
1	B	103	ASP	2.1
1	B	567	ASP	2.1
1	A	564	VAL	2.1
1	B	388	ALA	2.1
1	A	247	THR	2.1
1	A	391	THR	2.1
1	B	136	LEU	2.1
1	B	154	HIS	2.1
1	A	244	ASP	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](#)

There are no monosaccharides in this entry.

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. 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	B-factors(Å ²)	Q<0.9
3	MLI	A	702	7/7	0.68	0.36	34,36,37,38	7
3	MLI	B	702	7/7	0.94	0.15	42,43,44,44	0
2	NAG	B	701	15/15	0.96	0.21	23,25,27,28	0
2	NAG	A	701	15/15	0.97	0.23	26,27,30,30	0

6.5 Other polymers [i](#)

There are no such residues in this entry.