



wwPDB X-ray Structure Validation Summary Report ⓘ

Jan 31, 2016 – 08:07 PM GMT

PDB ID : 1IVE
Title : STRUCTURES OF AROMATIC INHIBITORS OF INFLUENZA VIRUS
NEURAMINIDASE
Authors : Jedrzejewski, M.J.; Luo, M.
Deposited on : 1994-12-12
Resolution : 2.40 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

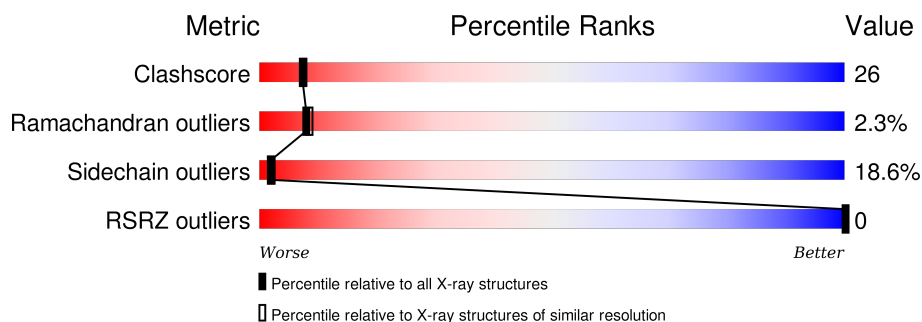
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION



The reported resolution of this entry is 2.40 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	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)
RSRZ outliers	91569	2928 (2.40-2.40)

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	388	
1	B	388	

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	472	-	-	-	X
2	NAG	B	472	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	BMA	A	476	-	-	X	-
3	FUL	A	477	-	-	X	-
4	NAG	A	478	-	-	X	-
4	NAG	A	479	-	-	X	X
4	NAG	B	479	-	-	-	X
5	MAN	B	481	-	-	X	-
5	MAN	B	482	X	-	-	X
5	MAN	B	485	X	-	-	-
8	FUC	B	477	X	-	-	-

2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 8216 atoms, of which 1788 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 INFLUENZA A SUBTYPE N2 NEURAMINIDASE.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	388	Total	C	H	N	O	S	0	0	0
			3745	1866	723	545	588	23			
1	B	388	Total	C	H	N	O	S	0	0	0
			3745	1866	723	545	588	23			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	339	ASP	ASN	CONFLICT	UNP P06820
B	339	ASP	ASN	CONFLICT	UNP P06820

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

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	2	Total	C	H	N	O		0	0
			55	16	27	2	10			
2	B	2	Total	C	H	N	O		0	0
			55	16	27	2	10			

- Molecule 3 is a polymer of unknown type called SUGAR (4-MER).

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	4	Total	C	H	N	O		0	0
			96	28	47	2	19			

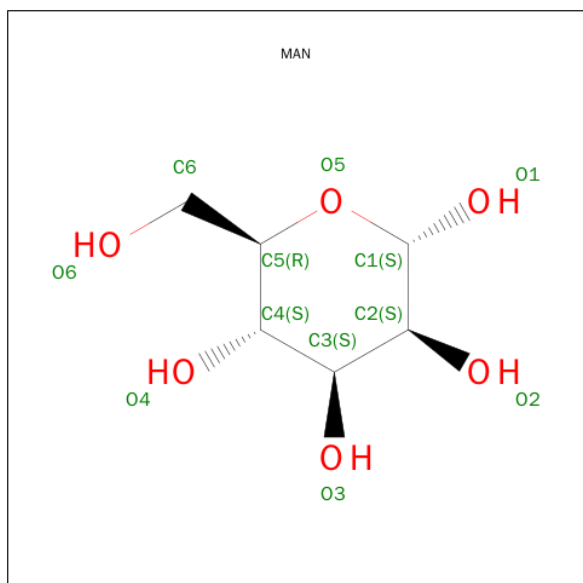
- Molecule 4 is a polymer of unknown type called SUGAR (3-MER).

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
4	A	3	Total	C	H	N	O		0	0
			74	22	35	2	15			
4	B	3	Total	C	H	N	O		0	0
			74	22	35	2	15			

- Molecule 5 is a polymer of unknown type called SUGAR (2-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	B	2	Total	C	H	O	0	0
			43	12	21	10		
5	B	2	Total	C	H	O	0	0
			43	12	21	10		

- Molecule 6 is SUGAR (ALPHA-D-MANNOSE) (three-letter code: MAN) (formula: C₆H₁₂O₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	B	1	Total	C	H	O	0	0
			22	6	11	5		
6	B	1	Total	C	H	O	0	0
			22	6	11	5		

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

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	A	2	Total	C	H	N	O	0
			55	16	27	2	10	
7	B	2	Total	C	H	N	O	0
			55	16	27	2	10	

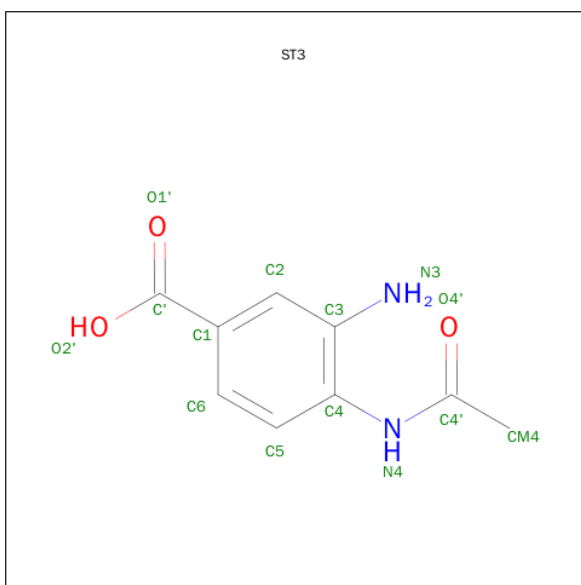
- Molecule 8 is a polymer of unknown type called SUGAR (4-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
8	B	4	Total	C	H	N	O	0	0
			96	28	47	2	19		

- Molecule 9 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	B	1	Total	Ca	0	0
			1	1		
9	A	1	Total	Ca	0	0
			1	1		

- Molecule 10 is 4-(ACETYLAMINO)-3-AMINO BENZOIC ACID (three-letter code: ST3) (formula: C₉H₁₀N₂O₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
10	A	1	Total	C	H	N	O	0	0
			17	9	3	2	3		
10	B	1	Total	C	H	N	O	0	0
			17	9	3	2	3		

4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	121.88Å 140.88Å 141.49Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	6.50 – 2.40 30.59 – 2.40	Depositor EDS
% Data completeness (in resolution range)	(Not available) (6.50-2.40) 54.4 (30.59-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.93 (at 2.42Å)	Xtriage
Refinement program	X-PLOR	Depositor
R, R_{free}	0.217 , (Not available) 0.223 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	18.4	Xtriage
Anisotropy	1.296	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 62.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Outliers	0 of 25898 reflections	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	8216	wwPDB-VP
Average B, all atoms (Å ²)	12.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.02% 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: BMA, NAG, CA, NDG, FUC, ST3, FUL, MAN

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	0.69	0/3092	0.77	1/4194 (0.0%)
1	B	0.69	0/3092	0.77	1/4194 (0.0%)
All	All	0.69	0/6184	0.77	2/8388 (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	A	0	18
1	B	0	18
5	B	2	0
8	B	1	0
All	All	3	36

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	103	ASP	CB-CG-OD1	5.67	123.40	118.30
1	B	103	ASP	CB-CG-OD1	5.67	123.40	118.30

All (3) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
5	B	482	MAN	C1
8	B	477	FUC	C1
5	B	485	MAN	C1

5 of 36 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	117	THR	Mainchain
1	A	118	ARG	Sidechain
1	A	121	TYR	Sidechain
1	A	122	VAL	Mainchain
1	A	84	TYR	Sidechain

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	3022	723	2851	145	3
1	B	3022	723	2852	158	4
2	A	28	27	25	4	0
2	B	28	27	25	5	0
3	A	49	47	43	4	16
4	A	39	35	32	24	6
4	B	39	35	32	6	1
5	B	44	42	38	2	12
6	B	22	22	20	0	0
7	A	28	27	24	2	0
7	B	28	27	24	2	0
8	B	49	47	43	4	0
9	A	1	0	0	0	0
9	B	1	0	0	0	0
10	A	14	3	9	2	0
10	B	14	3	9	1	0
All	All	6428	1788	6027	320	21

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 26.

The worst 5 of 320 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)
4:A:479:NAG:H82	1:B:455:THR:CG2	1.74	1.18

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:478:NAG:O5	1:B:455:THR:HB	1.49	1.11
4:A:479:NAG:H82	1:B:455:THR:OG1	1.59	1.02
4:A:479:NAG:H82	1:B:455:THR:HG21	1.46	0.97
1:A:437:TRP:H	1:A:469:ILE:HG21	1.33	0.93

The worst 5 of 21 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)
1:A:163:LEU:O	1:B:172:ARG:NH1[4_555]	1.32	0.88
1:A:163:LEU:O	1:B:172:ARG:HH12[4_555]	0.74	0.86
3:A:476:BMA:H61	4:A:479:NAG:HO6[3_654]	0.99	0.61
3:A:476:BMA:H4	4:A:479:NAG:H62[3_654]	1.00	0.60
3:A:477:FUL:H5	5:B:481:MAN:H62[3_654]	1.01	0.59

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	386/388 (100%)	326 (84%)	51 (13%)	9 (2%)	8	8
1	B	386/388 (100%)	326 (84%)	51 (13%)	9 (2%)	8	8
All	All	772/776 (100%)	652 (84%)	102 (13%)	18 (2%)	8	8

5 of 18 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	222	ILE
1	A	284	TYR
1	A	347	GLN
1	A	387	ASN
1	B	222	ILE

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	338/338 (100%)	275 (81%)	63 (19%)	2	2
1	B	338/338 (100%)	275 (81%)	63 (19%)	2	2
All	All	676/676 (100%)	550 (81%)	126 (19%)	2	2

5 of 126 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	430	ARG
1	B	140	LEU
1	B	420	ARG
1	A	435	ARG
1	B	82	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 25 such sidechains are listed below:

Mol	Chain	Res	Type
1	A	419	ASN
1	B	131	GLN
1	B	402	ASN
1	B	104	ASN
1	B	142	ASN

5.3.3 RNA ⓘ

There are no RNA molecules 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 ⓘ

26 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	472	1,2	14,14,15	1.83	2 (14%)	15,19,21	4.08	9 (60%)
2	NAG	A	473	2	14,14,15	2.54	6 (42%)	15,19,21	3.33	7 (46%)
3	NAG	A	474	1,3	14,14,15	1.52	4 (28%)	15,19,21	3.87	7 (46%)
3	NAG	A	475	3	14,14,15	3.02	10 (71%)	15,19,21	2.28	6 (40%)
3	BMA	A	476	3,4	11,11,12	5.08	9 (81%)	14,15,17	2.40	7 (50%)
3	FUL	A	477	3	10,10,11	2.41	5 (50%)	14,14,16	2.87	8 (57%)
4	NAG	A	478	1,4	14,14,15	2.22	4 (28%)	15,19,21	4.72	9 (60%)
4	NAG	A	479	3,4	14,14,15	3.77	9 (64%)	15,19,21	3.50	7 (46%)
4	BMA	A	480	5,4,6	11,11,12	4.61	9 (81%)	14,15,17	3.31	6 (42%)
7	NAG	A	484	1,7	14,14,15	2.05	6 (42%)	15,19,21	3.14	4 (26%)
7	NDG	A	485	7	14,14,15	2.33	5 (35%)	15,19,21	3.78	7 (46%)
2	NAG	B	472	1,2	14,14,15	1.83	2 (14%)	15,19,21	4.08	9 (60%)
2	NAG	B	473	2	14,14,15	2.54	6 (42%)	15,19,21	3.33	7 (46%)
8	NAG	B	474	1,8	14,14,15	1.52	4 (28%)	15,19,21	3.87	7 (46%)
8	NAG	B	475	8	14,14,15	3.02	10 (71%)	15,19,21	2.28	6 (40%)
8	BMA	B	476	8	11,11,12	5.08	9 (81%)	14,15,17	2.40	7 (50%)
8	FUC	B	477	8	10,10,11	2.41	5 (50%)	14,14,16	2.87	8 (57%)
4	NAG	B	478	1,4	14,14,15	2.22	4 (28%)	15,19,21	4.72	9 (60%)
4	NAG	B	479	4	14,14,15	3.77	9 (64%)	15,19,21	3.50	7 (46%)
4	BMA	B	480	5,4,6	11,11,12	4.61	9 (81%)	14,15,17	3.31	6 (42%)
5	MAN	B	481	5,4	11,11,12	3.68	6 (54%)	14,15,17	2.26	8 (57%)
5	MAN	B	482	5	11,11,12	5.69	9 (81%)	14,15,17	2.61	5 (35%)
5	MAN	B	484	5,4	11,11,12	3.68	6 (54%)	14,15,17	2.26	8 (57%)
5	MAN	B	485	5	11,11,12	5.69	9 (81%)	14,15,17	2.61	5 (35%)
7	NAG	B	487	1,7	14,14,15	2.05	6 (42%)	15,19,21	3.14	4 (26%)
7	NDG	B	488	7	14,14,15	2.33	5 (35%)	15,19,21	3.78	7 (46%)

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	472	1,2	-	0/6/23/26	0/1/1/1
2	NAG	A	473	2	-	0/6/23/26	0/1/1/1
3	NAG	A	474	1,3	-	0/6/23/26	0/1/1/1
3	NAG	A	475	3	-	0/6/23/26	0/1/1/1
3	BMA	A	476	3,4	-	0/2/19/22	1/1/1/1
3	FUL	A	477	3	-	0/0/17/20	0/1/1/1
4	NAG	A	478	1,4	-	0/6/23/26	0/1/1/1
4	NAG	A	479	3,4	-	0/6/23/26	0/1/1/1
4	BMA	A	480	5,4,6	-	0/2/19/22	0/1/1/1
7	NAG	A	484	1,7	-	0/6/23/26	0/1/1/1
7	NDG	A	485	7	-	0/6/23/26	0/1/1/1
2	NAG	B	472	1,2	-	0/6/23/26	0/1/1/1
2	NAG	B	473	2	-	0/6/23/26	0/1/1/1
8	NAG	B	474	1,8	-	0/6/23/26	0/1/1/1
8	NAG	B	475	8	-	0/6/23/26	0/1/1/1
8	BMA	B	476	8	-	0/2/19/22	1/1/1/1
8	FUC	B	477	8	1/1/5/5	0/0/17/20	0/1/1/1
4	NAG	B	478	1,4	-	0/6/23/26	0/1/1/1
4	NAG	B	479	4	-	0/6/23/26	0/1/1/1
4	BMA	B	480	5,4,6	-	0/2/19/22	0/1/1/1
5	MAN	B	481	5,4	-	0/2/19/22	0/1/1/1
5	MAN	B	482	5	1/1/5/5	0/2/19/22	0/1/1/1
5	MAN	B	484	5,4	-	0/2/19/22	0/1/1/1
5	MAN	B	485	5	1/1/5/5	0/2/19/22	0/1/1/1
7	NAG	B	487	1,7	-	0/6/23/26	0/1/1/1
7	NDG	B	488	7	-	0/6/23/26	0/1/1/1

The worst 5 of 168 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	B	487	NAG	C4-C3	-3.58	1.43	1.52
7	A	484	NAG	C4-C3	-3.58	1.43	1.52
7	B	487	NAG	C4-C5	-2.95	1.46	1.53
7	A	484	NAG	C4-C5	-2.95	1.46	1.53
7	B	488	NDG	O-C1	-2.74	1.39	1.43

The worst 5 of 180 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	479	NAG	C4-C3-C2	-9.08	97.12	111.23
4	A	479	NAG	C4-C3-C2	-9.08	97.12	111.23
4	B	478	NAG	C8-C7-N2	-7.13	102.47	116.11
4	A	478	NAG	C8-C7-N2	-7.13	102.47	116.11
2	A	473	NAG	C4-C3-C2	-6.76	100.73	111.23

All (3) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
8	B	477	FUC	C1
5	B	482	MAN	C1
5	B	485	MAN	C1

There are no torsion outliers.

All (2) ring outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	476	BMA	C1-C2-C3-C4-C5-O5
8	B	476	BMA	C1-C2-C3-C4-C5-O5

22 monomers are involved in 68 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	472	NAG	2	0
2	A	473	NAG	2	0
3	A	475	NAG	4	0
3	A	476	BMA	1	7
3	A	477	FUL	3	9
4	A	478	NAG	10	0
4	A	479	NAG	16	6
4	A	480	BMA	2	0
7	A	485	NDG	2	0
2	B	472	NAG	2	0
2	B	473	NAG	3	0
8	B	475	NAG	4	0
8	B	476	BMA	1	0
8	B	477	FUC	3	0
4	B	478	NAG	4	0
4	B	479	NAG	4	1
4	B	480	BMA	2	0
5	B	481	MAN	1	8
5	B	482	MAN	1	4

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	484	MAN	1	0
5	B	485	MAN	1	0
7	B	488	NDG	2	0

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 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	$\# Z > 2$
10	ST3	A	471	-	11,14,14	0.83	0	15,19,19	2.23	2 (13%)
10	ST3	B	471	-	11,14,14	0.83	0	15,19,19	2.23	2 (13%)
6	MAN	B	483	4	11,11,12	4.42	5 (45%)	14,15,17	2.17	3 (21%)
6	MAN	B	486	4	11,11,12	4.42	5 (45%)	14,15,17	2.17	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
10	ST3	A	471	-	-	0/4/8/8	0/1/1/1
10	ST3	B	471	-	-	0/4/8/8	0/1/1/1
6	MAN	B	483	4	-	0/2/19/22	1/1/1/1
6	MAN	B	486	4	-	0/2/19/22	1/1/1/1

The worst 5 of 10 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	486	MAN	O5-C1	2.63	1.48	1.43
6	B	483	MAN	O5-C1	2.63	1.48	1.43
6	B	486	MAN	C1-C2	4.64	1.63	1.52
6	B	483	MAN	C1-C2	4.64	1.63	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	486	MAN	C4-C5	4.85	1.63	1.53

The worst 5 of 10 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
10	B	471	ST3	C4-N4-C4'	-7.17	110.46	127.60
10	A	471	ST3	C4-N4-C4'	-7.17	110.46	127.60
6	B	486	MAN	C3-C4-C5	3.30	115.96	110.20
6	B	483	MAN	C3-C4-C5	3.30	115.96	110.20
10	B	471	ST3	CM4-C4'-N4	3.33	119.70	114.97

There are no chirality outliers.

There are no torsion outliers.

All (2) ring outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	B	486	MAN	C1-C2-C3-C4-C5-O5
6	B	483	MAN	C1-C2-C3-C4-C5-O5

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
10	A	471	ST3	2	0
10	B	471	ST3	1	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	A	388/388 (100%)	-0.49	0 100 100	2, 11, 19, 28	0
1	B	388/388 (100%)	-0.45	0 100 100	2, 11, 19, 28	0
All	All	776/776 (100%)	-0.47	0 100 100	2, 11, 19, 28	0

There are no RSRZ outliers to report.

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
5	MAN	B	482	11/12	0.54	0.50	48.22	15,15,31,33	0
4	NAG	B	479	14/15	0.76	0.56	15.53	15,15,27,28	0
4	NAG	A	479	14/15	0.85	0.47	13.24	15,15,27,28	0
2	NAG	B	472	14/15	0.92	0.29	3.31	15,15,26,27	0
2	NAG	A	472	14/15	0.88	0.30	3.08	15,15,26,27	0
4	NAG	B	478	14/15	0.78	0.21	1.81	11,15,22,24	0
4	NAG	A	478	14/15	0.75	0.18	1.68	11,15,22,24	0
7	NAG	B	487	14/15	0.87	0.17	1.22	15,15,25,26	0
7	NAG	A	484	14/15	0.90	0.19	0.88	15,15,25,26	0
5	MAN	B	485	11/12	0.53	0.53	-	15,15,31,33	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
8	FUC	B	477	10/11	0.91	0.24	-	15,15,25,27	0
5	MAN	B	484	11/12	0.74	0.54	-	15,15,29,29	0
2	NAG	A	473	14/15	0.83	0.37	-	15,15,30,31	0
4	BMA	A	480	11/12	0.72	0.39	-	15,26,30,31	0
7	NDG	A	485	14/15	0.85	0.36	-	15,15,35,37	0
5	MAN	B	481	11/12	0.64	0.40	-	15,15,29,29	0
3	NAG	A	474	14/15	0.92	0.17	-	15,16,24,29	0
8	NAG	B	474	14/15	0.87	0.19	-	15,16,24,29	0
3	FUL	A	477	10/11	0.82	0.39	-	15,15,25,27	0
3	NAG	A	475	14/15	0.80	0.41	-	15,15,32,34	0
8	NAG	B	475	14/15	0.82	0.30	-	15,15,32,34	0
4	BMA	B	480	11/12	0.77	0.48	-	15,26,30,31	0
7	NDG	B	488	14/15	0.78	0.42	-	15,15,35,37	0
3	BMA	A	476	11/12	0.77	0.49	-	15,15,32,33	0
8	BMA	B	476	11/12	0.65	0.66	-	15,15,32,33	0
2	NAG	B	473	14/15	0.92	0.35	-	15,15,30,31	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
10	ST3	A	471	14/14	0.91	0.17	0.73	12,18,22,24	0
10	ST3	B	471	14/14	0.93	0.14	0.50	12,18,22,24	0
9	CA	A	470	1/1	0.99	0.07	-1.56	14,14,14,14	0
9	CA	B	470	1/1	0.99	0.07	-1.68	14,14,14,14	0
6	MAN	B	483	11/12	0.86	0.36	-	15,15,29,32	0
6	MAN	B	486	11/12	0.86	0.52	-	15,15,29,32	0

6.5 Other polymers [i](#)

There are no such residues in this entry.