Full wwPDB X-ray Structure Validation Report ⓘ

Nov 5, 2026 – 06:14 PM JST

PDB ID	:	47VF / pdb_000047vf
Title	:	The structure of hemagglutinin T160A mutant (H3 numbering) from H9N2 influenza virus (A/chicken/Tianjin/CAS/2007) in complex with human receptor analog 6'SLNLN
Authors	:	Wang, M.; Peng, R.C.; Qi, J.X.; Shi, Y.; Gao, G.F.
Deposited on	:	2026-11-01
Resolution	:	2.10 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:

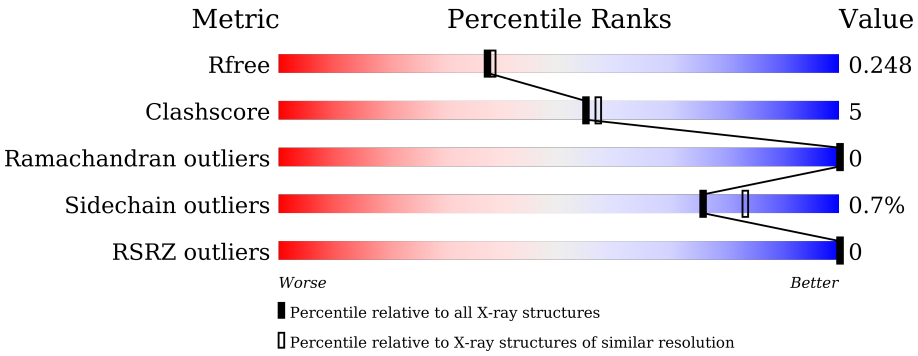
MolProbity	:	4-5-2 with Phenix2.2
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	2.2
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.51

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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}	180053	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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 <=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	316	86% 14% .
1	C	316	88% 12%
2	B	176	84% 9% 8%
2	D	176	88% 5% 8%
3	E	3	33% 33% 33%

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Mol	Chain	Length	Quality of chain
3	F	3	33% 33% 33%

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 8454 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 Hemagglutinin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	316	Total	C	N	O	S	0	0	0
			2475	1553	437	473	12			
1	C	316	Total	C	N	O	S	0	0	0
			2475	1553	437	473	12			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	86	VAL	THR	conflict	UNP A7YL49
C	86	VAL	THR	conflict	UNP A7YL49

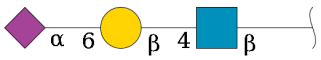
- Molecule 2 is a protein called Hemagglutinin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	162	Total	C	N	O	S	0	0	0
			1291	802	227	255	7			
2	D	162	Total	C	N	O	S	0	0	0
			1291	802	227	255	7			

There are 4 discrepancies between the modelled and reference sequences:

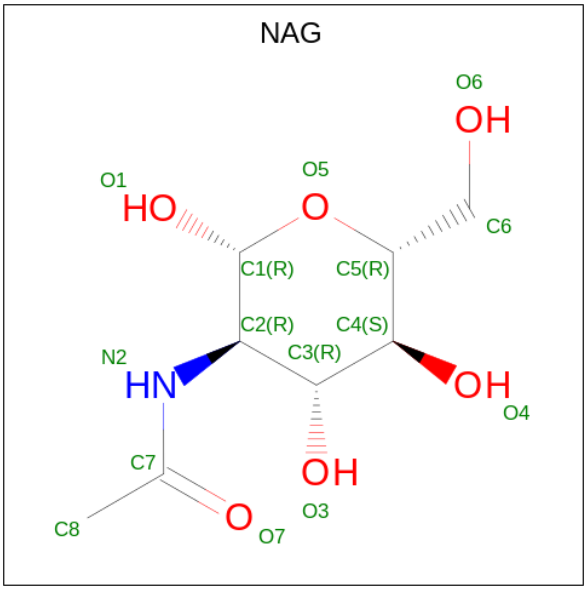
Chain	Residue	Modelled	Actual	Comment	Reference
B	385	VAL	ILE	conflict	UNP B5U6G4
B	487	ARG	LYS	conflict	UNP B5U6G4
D	385	VAL	ILE	conflict	UNP B5U6G4
D	487	ARG	LYS	conflict	UNP B5U6G4

- Molecule 3 is an oligosaccharide called N-acetyl-alpha-neuraminic acid-(2-6)-beta-D-galacto pyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf	Trace
3	E	3	Total C N O 46 25 2 19	0	0	0
3	F	3	Total C N O 46 25 2 19	0	0	0

- Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	242	Total O 242 242	0	0
5	B	106	Total O 106 106	0	0
5	C	253	Total O 253 253	0	0
5	D	159	Total O 159 159	0	0

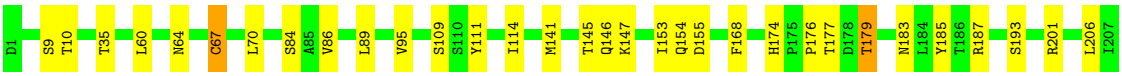
3 Residue-property plots (i)

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: Hemagglutinin

Chain A:





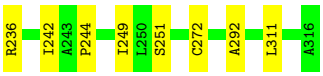


• Molecule 1: Hemagglutinin

Chain C:







• Molecule 2: Hemagglutinin

Chain B:

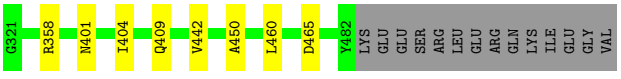




• Molecule 2: Hemagglutinin

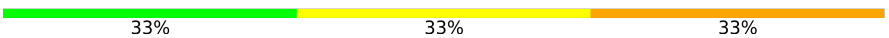
Chain D:





• Molecule 3: N-acetyl-alpha-neuraminic acid-(2-6)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

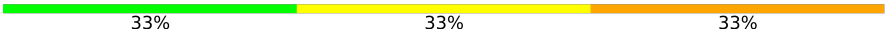
Chain E:





• Molecule 3: N-acetyl-alpha-neuraminic acid-(2-6)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain F:





4 Data and refinement statistics (i)

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α, β, γ	94.37Å 94.37Å 357.99Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	43.88 – 2.10 43.88 – 2.10	Depositor EDS
% Data completeness (in resolution range)	91.8 (43.88-2.10) 92.0 (43.88-2.10)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.61 (at 2.10Å)	Xtriage
Refinement program	PHENIX 2.1_6048	Depositor
R, R_{free}	0.229 , 0.248 0.228 , 0.248	Depositor DCC
R_{free} test set	3172 reflections (4.56%)	wwPDB-VP
Wilson B-factor (Å^2)	24.5	Xtriage
Anisotropy	0.088	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 24.3	EDS
L-test for twinning[2]	$< \|L\| > = 0.43$, $< L^2 > = 0.26$	Xtriage
Estimated twinning fraction	0.152 for -h-k,k,-l	Xtriage
F_o,F_c correlation	0.97	EDS
Total number of atoms	8454	wwPDB-VP
Average B, all atoms (Å^2)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.52% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 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: NAG, SIA, GAL

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	0.09	0/2533	0.27	0/3453
1	C	0.09	0/2533	0.28	0/3453
2	B	0.10	0/1312	0.25	0/1767
2	D	0.10	0/1312	0.24	0/1767
All	All	0.09	0/7690	0.27	0/10440

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	2475	0	2414	31	0
1	C	2475	0	2415	23	0
2	B	1291	0	1230	12	0
2	D	1291	0	1230	6	0
3	E	46	0	40	1	0
3	F	46	0	40	2	0
4	A	42	0	39	4	0
4	C	28	0	26	0	0
5	A	242	0	0	6	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	B	106	0	0	7	0
5	C	253	0	0	3	0
5	D	159	0	0	3	0
All	All	8454	0	7434	72	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 5.

All (72) 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:A:89:LEU:HD11	1:A:95:VAL:HG23	1.76	0.68
1:A:208:GLY:C	5:A:501:HOH:O	2.36	0.68
1:C:154:GLN:O	1:C:236:ARG:HA	1.96	0.66
2:B:374:ASN:ND2	5:B:502:HOH:O	2.30	0.64
1:A:206:LEU:O	1:A:210:ARG:NH2	2.30	0.63
1:A:155:ASP:H	4:A:401:NAG:H81	1.65	0.61
1:A:154:GLN:O	1:A:236:ARG:HA	2.01	0.61
2:D:450:ALA:HB2	2:D:460:LEU:HD23	1.83	0.61
2:D:358:ARG:HD3	5:D:519:HOH:O	2.04	0.57
1:C:89:LEU:HD11	1:C:95:VAL:HG23	1.88	0.55
1:C:185:TYR:O	1:C:187:ARG:N	2.38	0.55
1:C:209:PRO:HD3	5:C:542:HOH:O	2.07	0.55
2:B:358:ARG:NH2	5:B:509:HOH:O	2.40	0.55
1:C:60:LEU:HD12	1:C:70:LEU:HD13	1.89	0.54
1:C:5:ILE:HD11	2:D:442:VAL:HG21	1.89	0.54
2:B:453:ASP:OD2	2:B:457:CYS:HB2	2.08	0.54
1:A:179:THR:O	1:A:183:ASN:ND2	2.41	0.54
1:A:60:LEU:HD12	1:A:70:LEU:HD13	1.91	0.53
1:A:210:ARG:HD3	1:A:219:ARG:HG2	1.89	0.53
1:C:311:LEU:H	1:C:311:LEU:HD23	1.74	0.53
2:B:348:ASN:HB2	5:B:518:HOH:O	2.08	0.52
1:C:67:CYS:HB3	1:C:70:LEU:HD12	1.92	0.52
1:C:145:THR:HG21	3:F:3:SIA:H111	1.91	0.52
2:B:382:GLN:NE2	2:B:384:GLU:OE2	2.37	0.52
1:A:64:ASN:HB3	1:A:67:CYS:SG	2.49	0.52
1:A:174:HIS:NE2	1:A:221:ASP:OD2	2.41	0.52
1:A:154:GLN:HA	4:A:401:NAG:H81	1.90	0.52
1:C:111:TYR:HA	1:C:249:ILE:O	2.10	0.51
1:C:174:HIS:NE2	1:C:221:ASP:OD2	2.43	0.51
1:A:67:CYS:HB3	1:A:70:LEU:HD12	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:402:NAG:H83	4:A:402:NAG:H3	1.93	0.50
2:B:374:ASN:CG	5:B:502:HOH:O	2.54	0.50
2:B:401:ASN:HA	2:B:404:ILE:HD12	1.92	0.50
2:D:409:GLN:NE2	5:D:501:HOH:O	2.24	0.50
1:C:64:ASN:HB3	1:C:67:CYS:SG	2.51	0.50
1:A:145:THR:HG21	3:E:3:SIA:H111	1.95	0.49
2:B:450:ALA:HB2	2:B:460:LEU:HD23	1.94	0.49
1:C:168:PHE:CZ	1:C:225:SER:HB3	2.47	0.48
1:C:164:LYS:HD3	1:C:251:SER:HB2	1.95	0.48
1:C:192:THR:O	1:C:202:THR:HA	2.13	0.48
1:A:201:ARG:NE	5:A:512:HOH:O	2.39	0.47
1:A:311:LEU:HD23	1:A:311:LEU:H	1.79	0.47
2:B:479:ARG:HD2	5:B:508:HOH:O	2.15	0.47
2:D:401:ASN:HA	2:D:404:ILE:HD12	1.96	0.47
1:A:10:THR:HG22	5:A:537:HOH:O	2.15	0.47
1:A:176:PRO:HA	5:A:501:HOH:O	2.15	0.46
1:A:114:ILE:HD12	1:A:249:ILE:HD11	1.96	0.46
1:C:170:TRP:HB3	1:C:244:PRO:HG3	1.96	0.46
1:C:208:GLY:O	1:C:210:ARG:NH1	2.48	0.46
1:A:185:TYR:O	1:A:187:ARG:N	2.39	0.46
1:C:272:CYS:SG	1:C:292:ALA:HB1	2.56	0.45
1:A:168:PHE:CZ	1:A:225:SER:HB3	2.52	0.45
1:C:157:GLN:NE2	5:C:501:HOH:O	2.28	0.45
1:A:208:GLY:O	5:A:501:HOH:O	2.20	0.45
1:A:9:SER:O	2:B:335:SER:HA	2.17	0.44
2:D:465:ASP:HB3	5:D:635:HOH:O	2.17	0.44
2:B:366:ASP:HB2	5:B:504:HOH:O	2.16	0.44
1:A:272:CYS:SG	1:A:292:ALA:HB1	2.57	0.44
1:C:114:ILE:HD12	1:C:249:ILE:HD11	2.01	0.42
1:A:141:MET:HB3	1:A:242:ILE:HG22	2.01	0.42
1:A:193:SER:O	1:A:235:VAL:HA	2.19	0.42
1:A:84:SER:O	1:A:86:VAL:HG13	2.19	0.42
1:A:35:THR:HG21	4:A:403:NAG:H62	2.02	0.42
1:C:141:MET:HB3	1:C:242:ILE:HG22	2.01	0.41
1:A:109:SER:OG	1:A:252:GLY:HA3	2.21	0.41
1:A:177:THR:HB	5:A:583:HOH:O	2.19	0.41
5:C:751:HOH:O	3:F:1:NAG:H2	2.20	0.41
2:B:480:ARG:NH2	5:B:515:HOH:O	2.53	0.41
1:C:193:SER:O	1:C:235:VAL:HA	2.20	0.41
1:A:60:LEU:HD21	1:A:111:TYR:CZ	2.56	0.40
1:A:153:ILE:HD13	1:A:236:ARG:NH2	2.36	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:50:LEU:HB2	1:C:76:TRP:HB3	2.02	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	314/316 (99%)	299 (95%)	15 (5%)	0	100	100
1	C	314/316 (99%)	299 (95%)	15 (5%)	0	100	100
2	B	160/176 (91%)	158 (99%)	2 (1%)	0	100	100
2	D	160/176 (91%)	155 (97%)	5 (3%)	0	100	100
All	All	948/984 (96%)	911 (96%)	37 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	278/278 (100%)	274 (99%)	4 (1%)	59	67
1	C	278/278 (100%)	276 (99%)	2 (1%)	76	83
2	B	137/150 (91%)	137 (100%)	0	100	100
2	D	137/150 (91%)	137 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	830/856 (97%)	824 (99%)	6 (1%)	76 83

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	67	CYS
1	A	146	GLN
1	A	147	LYS
1	A	179	THR
1	C	67	CYS
1	C	206	LEU

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

Mol	Chain	Res	Type
2	D	362	GLN
2	D	436	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

6 monosaccharides 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	E	1	3	15,15,15	0.23	0	21,21,21	0.51	0
3	GAL	E	2	3	11,11,12	0.61	0	15,15,17	1.53	1 (6%)
3	SIA	E	3	3	20,20,21	1.59	3 (15%)	24,28,31	1.29	4 (16%)
3	NAG	F	1	3	15,15,15	0.11	0	21,21,21	0.12	0
3	GAL	F	2	3	11,11,12	0.53	0	15,15,17	0.95	0
3	SIA	F	3	3	20,20,21	1.97	2 (10%)	24,28,31	1.37	4 (16%)

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	E	1	3	-	2/6/26/26	0/1/1/1
3	GAL	E	2	3	-	0/2/19/22	0/1/1/1
3	SIA	E	3	3	-	0/18/34/38	0/1/1/1
3	NAG	F	1	3	-	0/6/26/26	0/1/1/1
3	GAL	F	2	3	-	0/2/19/22	0/1/1/1
3	SIA	F	3	3	-	2/18/34/38	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	F	3	SIA	C2-C1	7.56	1.59	1.52
3	E	3	SIA	C2-C1	5.50	1.57	1.52
3	E	3	SIA	O1B-C1	-2.90	1.21	1.30
3	F	3	SIA	O6-C2	2.81	1.47	1.43
3	E	3	SIA	O6-C2	2.76	1.47	1.43

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	E	2	GAL	C1-C2-C3	4.78	115.54	109.67
3	F	3	SIA	C6-O6-C2	3.49	118.80	111.34
3	E	3	SIA	C6-O6-C2	3.48	118.79	111.34
3	E	3	SIA	O1B-C1-C2	3.17	122.07	113.03
3	E	3	SIA	O1A-C1-C2	-2.41	116.89	122.57
3	F	3	SIA	O1B-C1-O1A	2.39	129.51	124.09
3	E	3	SIA	C4-C3-C2	2.34	114.00	109.81
3	F	3	SIA	C4-C3-C2	2.30	113.93	109.81
3	F	3	SIA	O1A-C1-C2	-2.12	117.56	122.57

There are no chirality outliers.

All (4) torsion outliers are listed below:

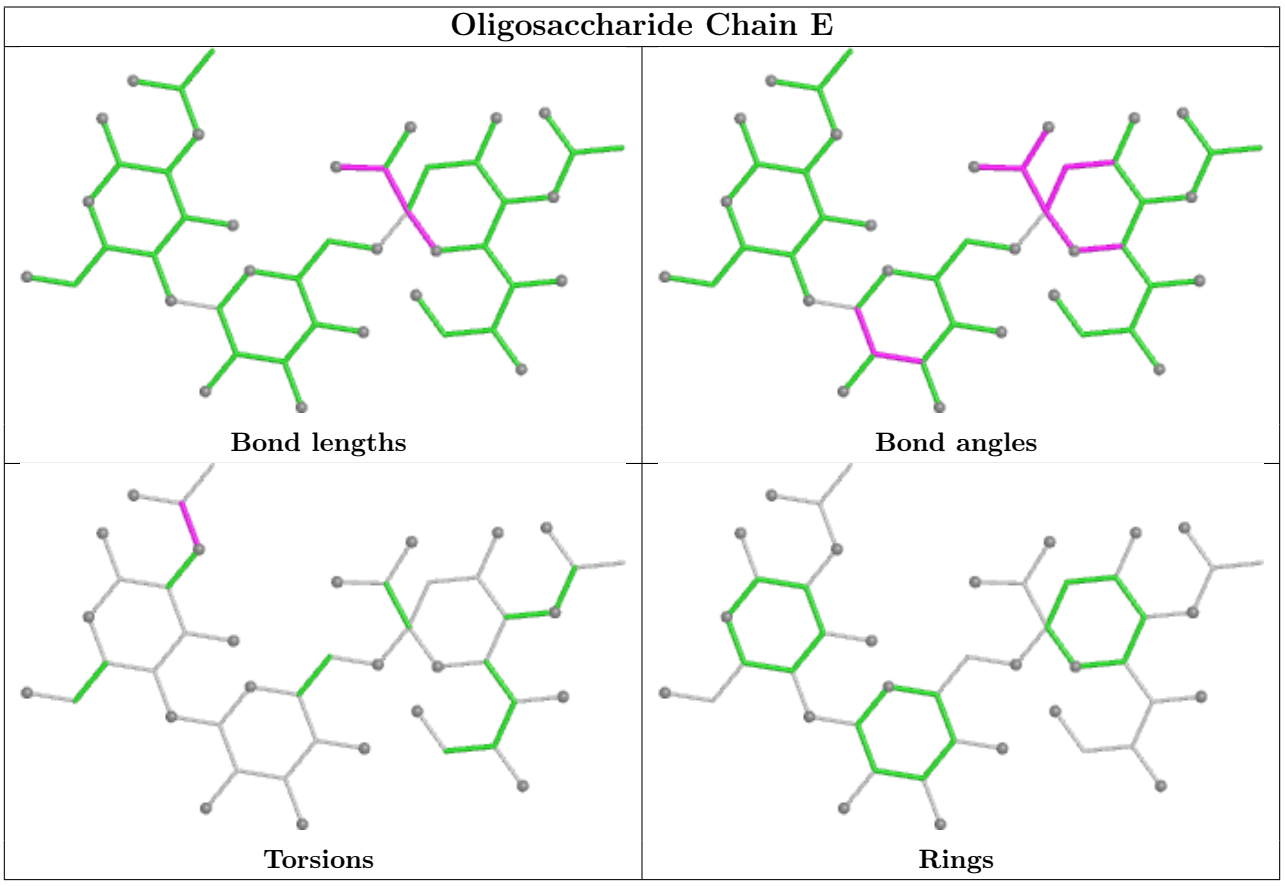
Mol	Chain	Res	Type	Atoms
3	E	1	NAG	C8-C7-N2-C2
3	E	1	NAG	O7-C7-N2-C2
3	F	3	SIA	O8-C8-C9-O9
3	F	3	SIA	C7-C8-C9-O9

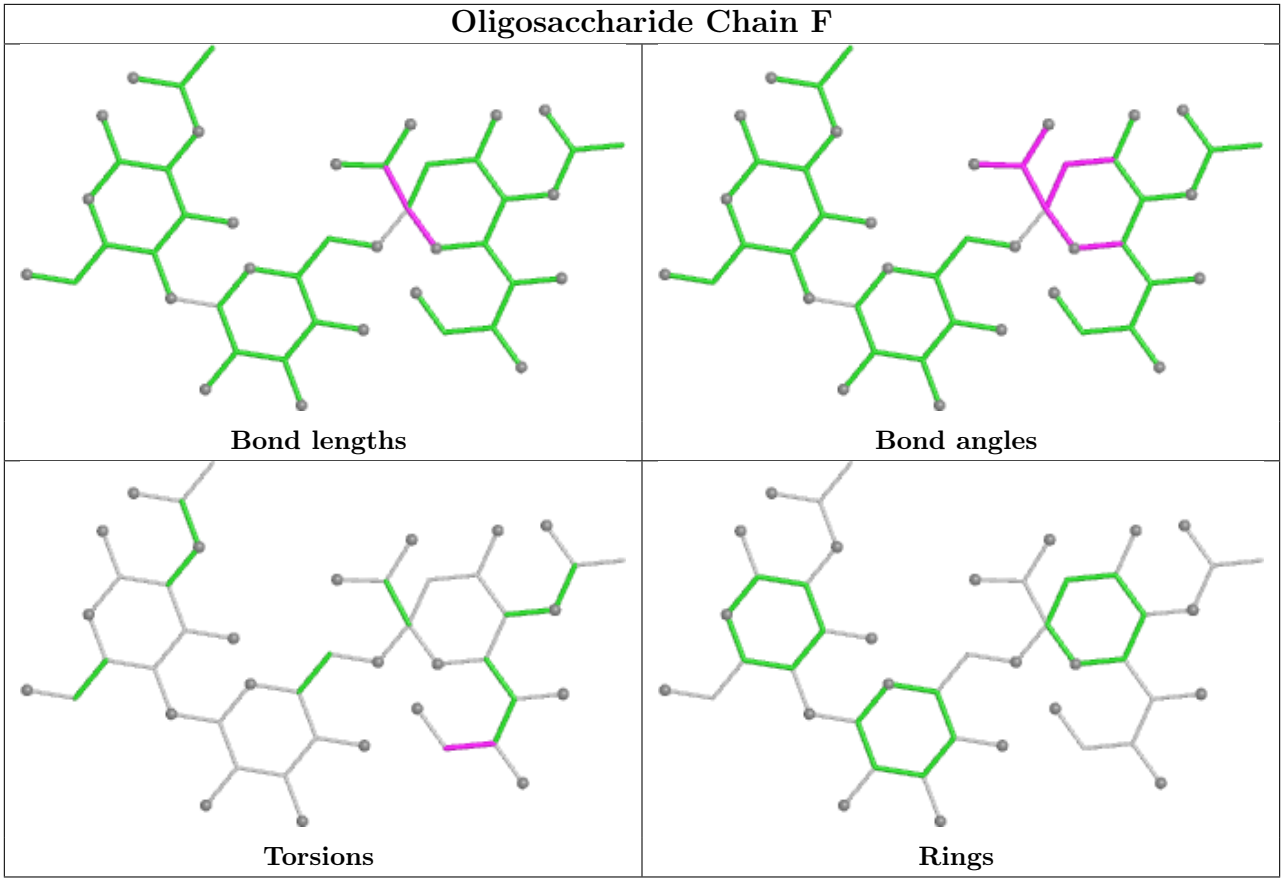
There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	F	1	NAG	1	0
3	F	3	SIA	1	0
3	E	3	SIA	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.





5.6 Ligand geometry (i)

5 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	$\#\|Z\| > 2$
4	NAG	A	401	1	14,14,15	0.22	0	17,19,21	0.39	0
4	NAG	C	402	1	14,14,15	0.22	0	17,19,21	0.44	0
4	NAG	A	403	1	14,14,15	0.24	0	17,19,21	0.56	0
4	NAG	C	401	1	14,14,15	0.25	0	17,19,21	0.41	0
4	NAG	A	402	1	14,14,15	0.32	0	17,19,21	1.34	2 (11%)

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
4	NAG	A	401	1	-	2/6/23/26	0/1/1/1
4	NAG	C	402	1	-	2/6/23/26	0/1/1/1
4	NAG	A	403	1	-	0/6/23/26	0/1/1/1
4	NAG	C	401	1	-	2/6/23/26	0/1/1/1
4	NAG	A	402	1	-	3/6/23/26	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	A	402	NAG	C2-N2-C7	4.47	129.27	122.90
4	A	402	NAG	C1-C2-N2	2.19	114.22	110.49

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	C	401	NAG	O5-C5-C6-O6
4	C	402	NAG	O5-C5-C6-O6
4	C	401	NAG	C4-C5-C6-O6
4	A	401	NAG	C8-C7-N2-C2
4	A	401	NAG	O7-C7-N2-C2
4	A	402	NAG	C8-C7-N2-C2
4	A	402	NAG	O7-C7-N2-C2
4	C	402	NAG	C4-C5-C6-O6
4	A	402	NAG	C3-C2-N2-C7

There are no ring outliers.

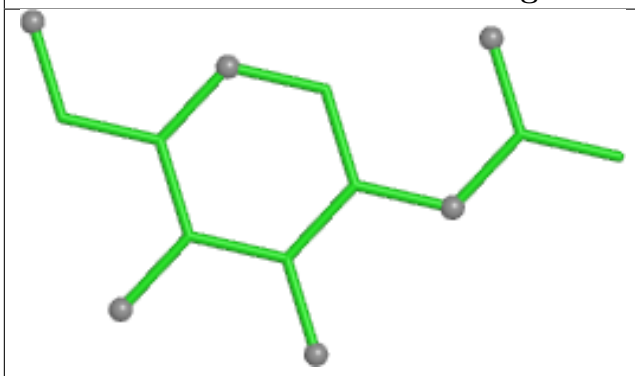
3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	401	NAG	2	0
4	A	403	NAG	1	0
4	A	402	NAG	1	0

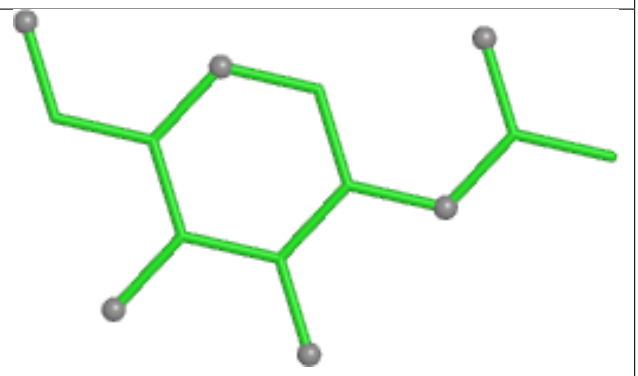
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will

also be included. For torsion angles, if less then 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

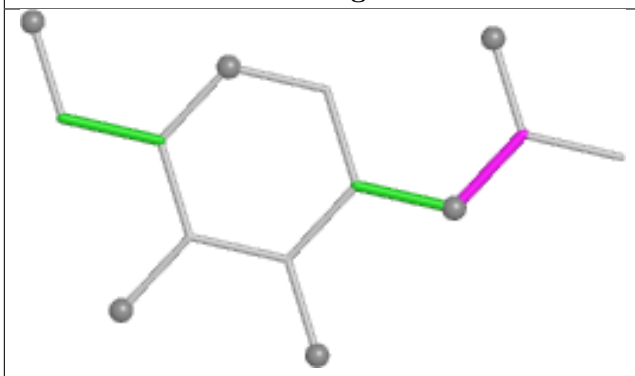
Ligand NAG A 401



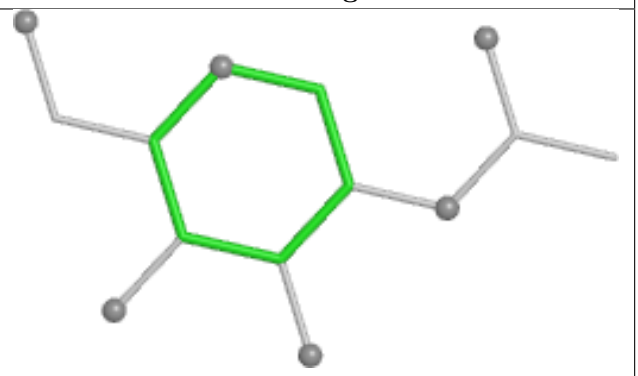
Bond lengths



Bond angles

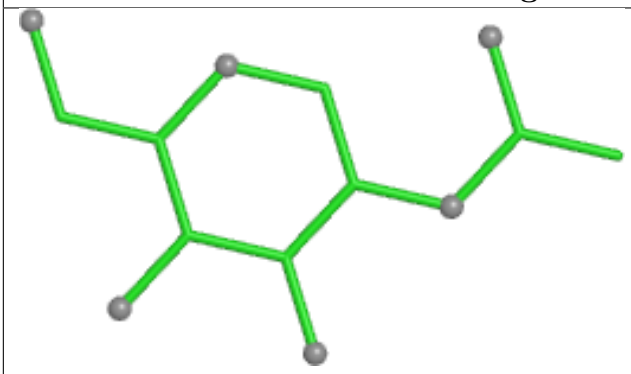


Torsions

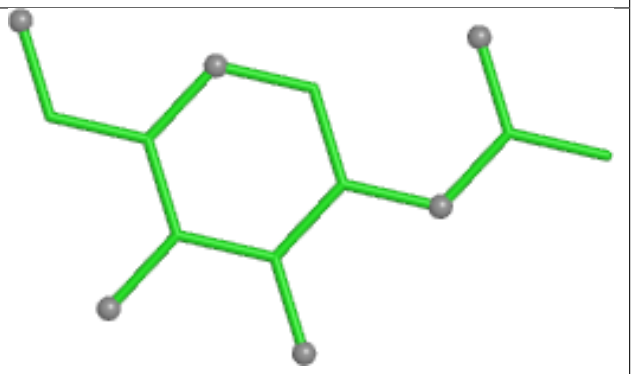


Rings

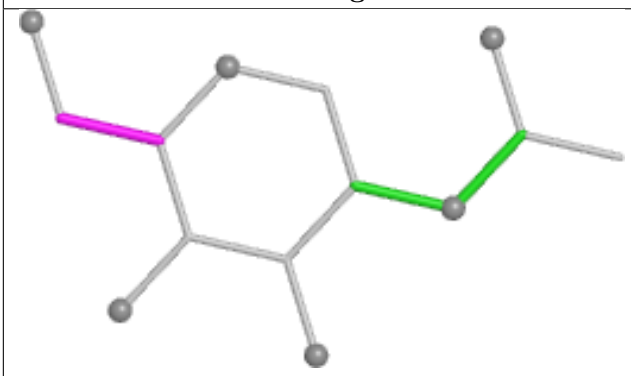
Ligand NAG C 402



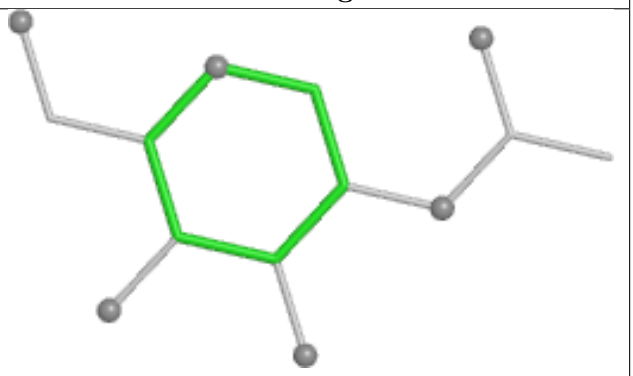
Bond lengths



Bond angles

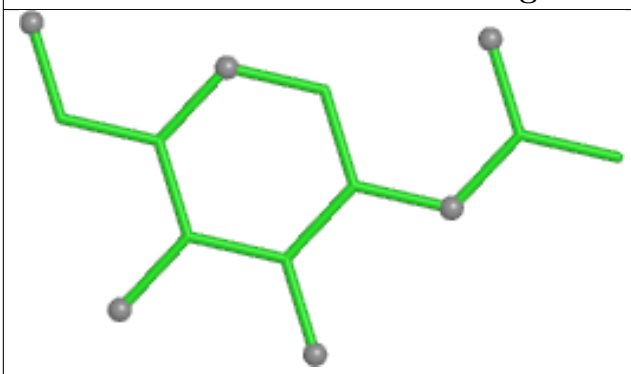


Torsions

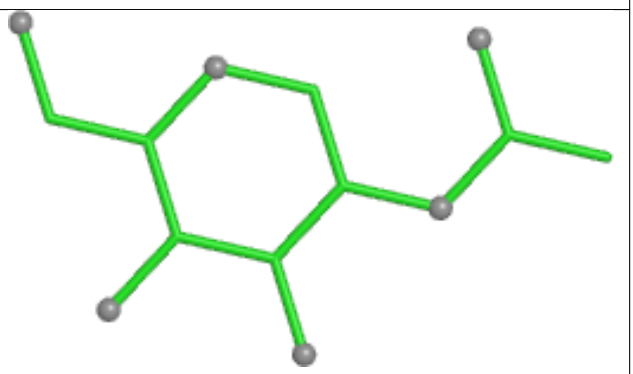


Rings

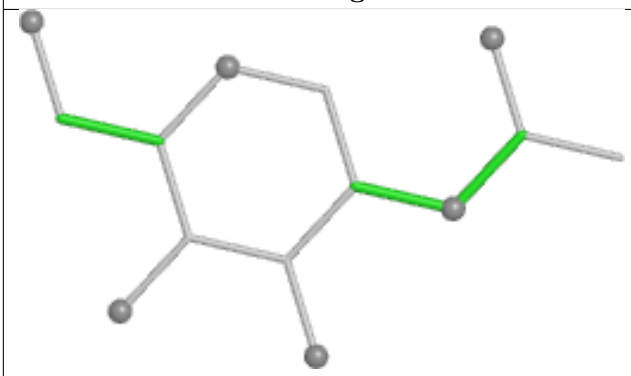
Ligand NAG A 403



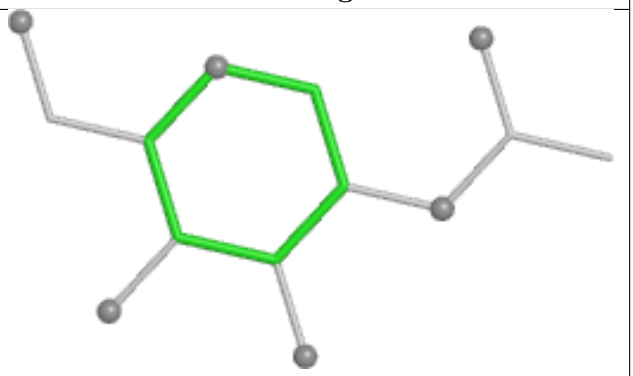
Bond lengths



Bond angles

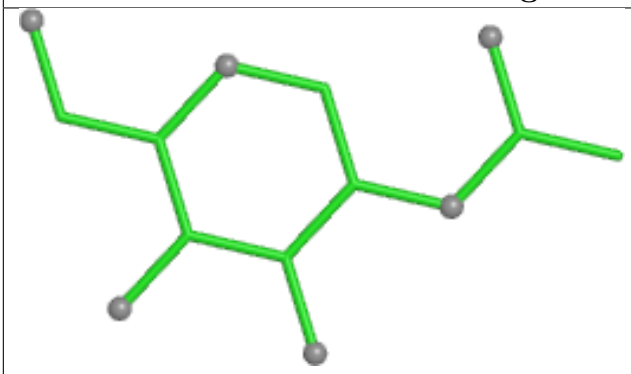


Torsions

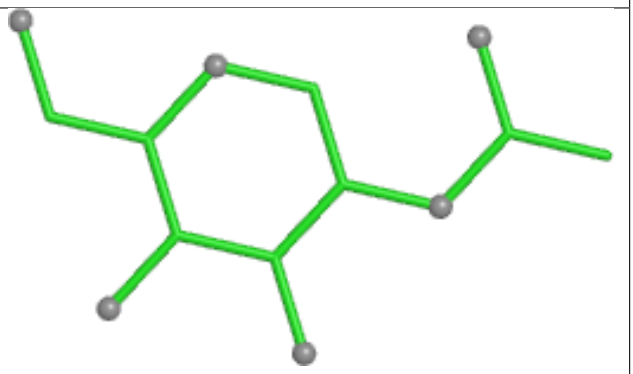


Rings

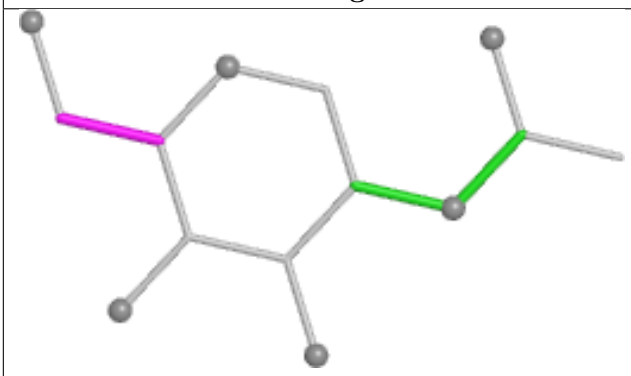
Ligand NAG C 401



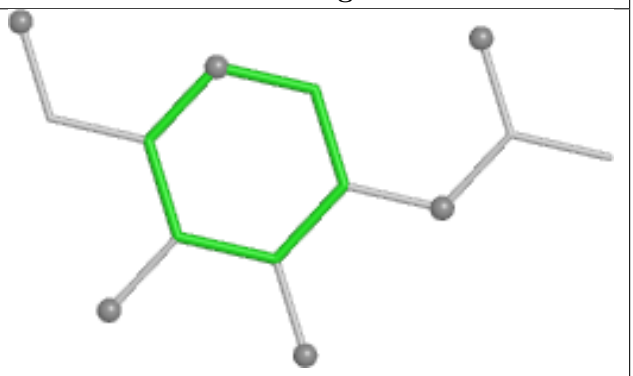
Bond lengths



Bond angles

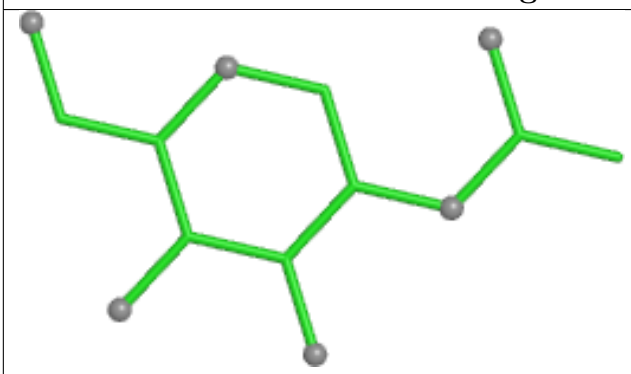


Torsions

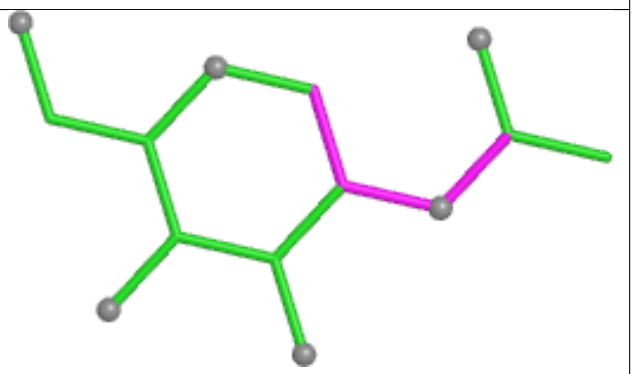


Rings

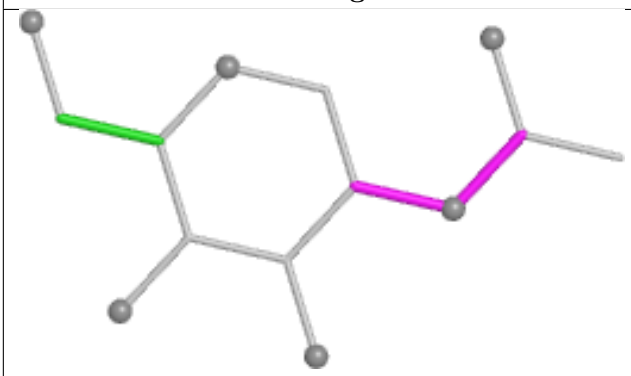
Ligand NAG A 402



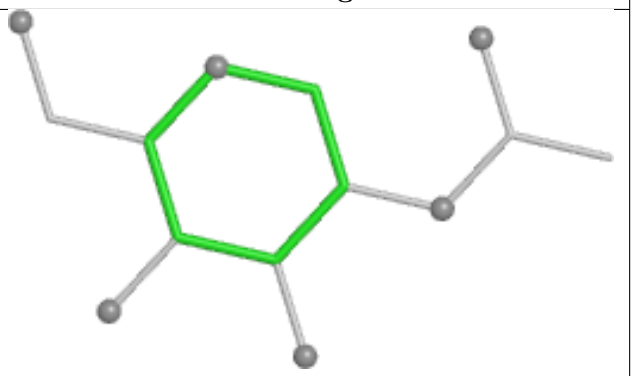
Bond lengths



Bond angles



Torsions



Rings

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, 95^{th} 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	316/316 (100%)	-1.44	0	100	100	16, 28, 42, 56	0
1	C	316/316 (100%)	-1.46	0	100	100	15, 28, 41, 50	0
2	B	162/176 (92%)	-1.29	0	100	100	15, 35, 61, 65	0
2	D	162/176 (92%)	-1.49	0	100	100	14, 25, 43, 51	0
All	All	956/984 (97%)	-1.43	0	100	100	14, 28, 48, 65	0

There are no RSRZ outliers to report.

6.2 Non-standard residues in protein, DNA, RNA chains

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates

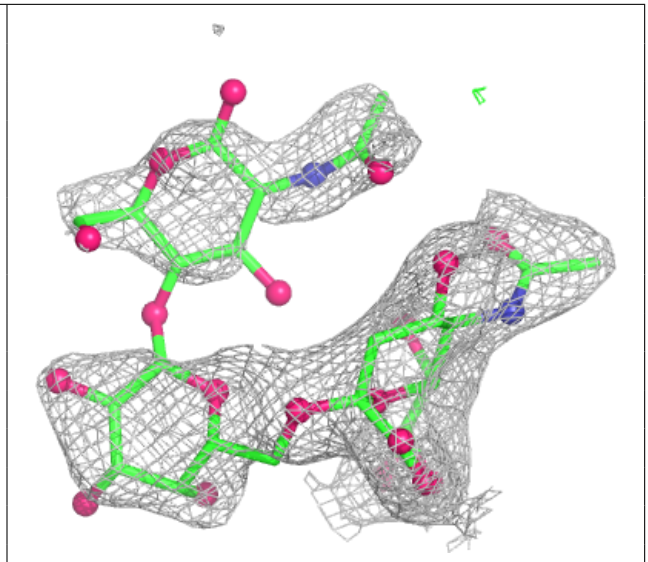
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

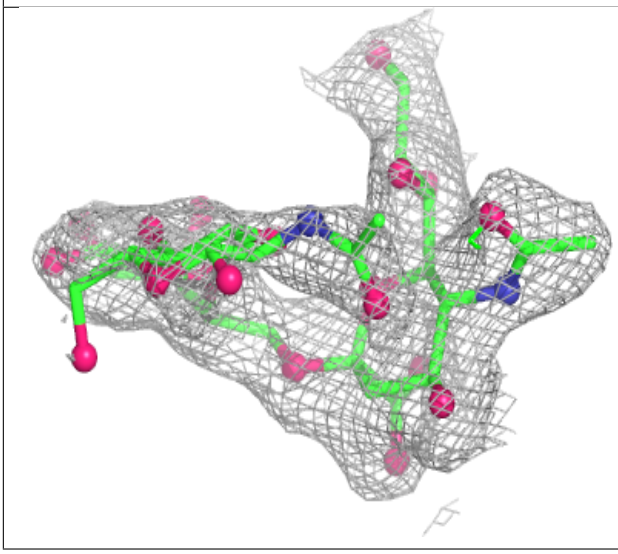
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	NAG	E	1	15/15	0.95	0.07	57,63,66,66	0
3	GAL	F	2	11/12	0.96	0.06	43,58,66,70	0
3	NAG	F	1	15/15	0.97	0.06	55,61,64,65	0
3	GAL	E	2	11/12	0.98	0.04	42,57,66,69	0
3	SIA	E	3	20/21	0.99	0.03	28,40,46,48	0
3	SIA	F	3	20/21	0.99	0.03	27,39,45,47	0

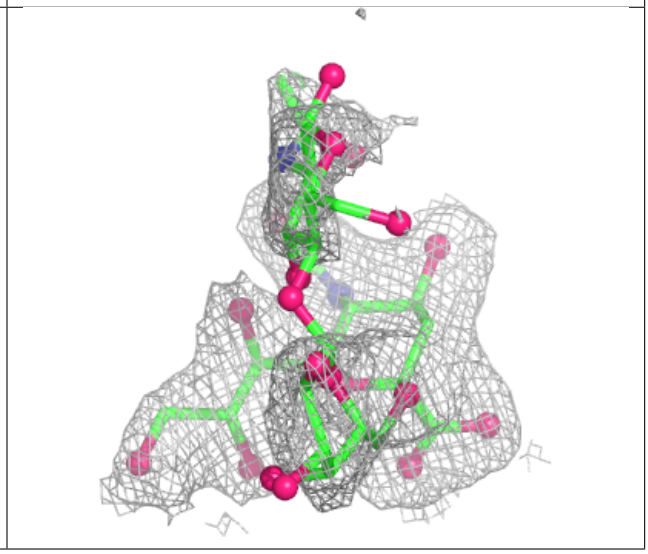
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Chain E:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

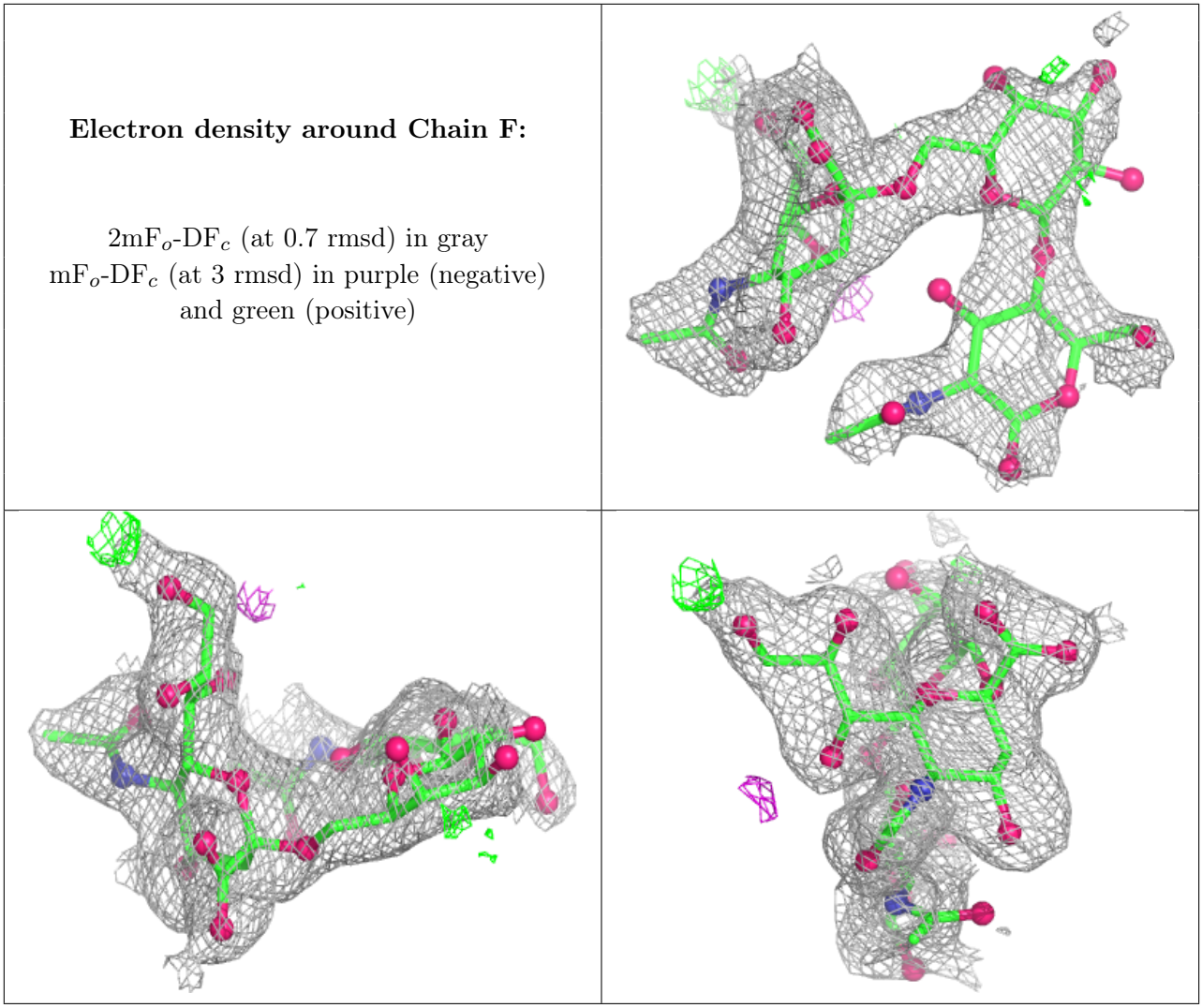






Electron density around Chain F:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)



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. The B-factors column lists the minimum, median, 95^{th} 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.

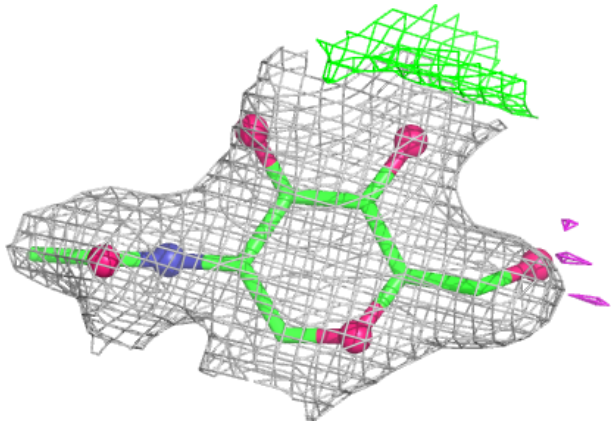
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
4	NAG	C	402	14/15	0.98	0.04	30,36,44,44	0
4	NAG	A	402	14/15	0.99	0.04	32,44,52,52	0
4	NAG	A	403	14/15	0.99	0.04	33,45,51,52	0
4	NAG	C	401	14/15	0.99	0.03	32,38,43,44	0
4	NAG	A	401	14/15	0.99	0.03	27,38,41,42	0

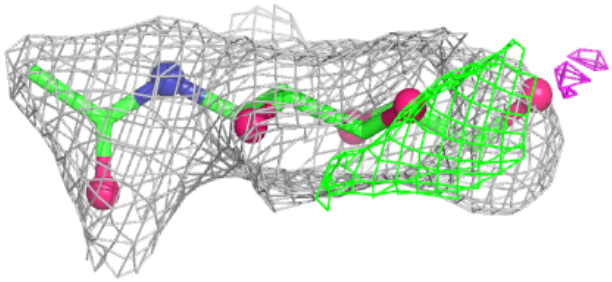
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different

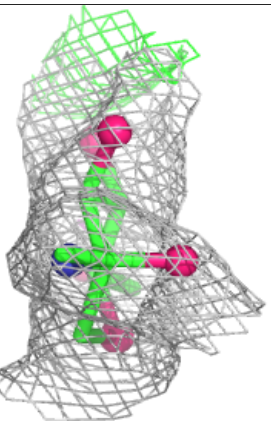
orientation to approximate a three-dimensional view.

Electron density around NAG C 402:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

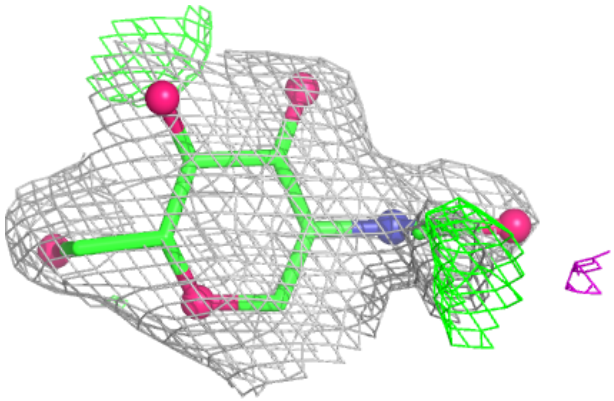


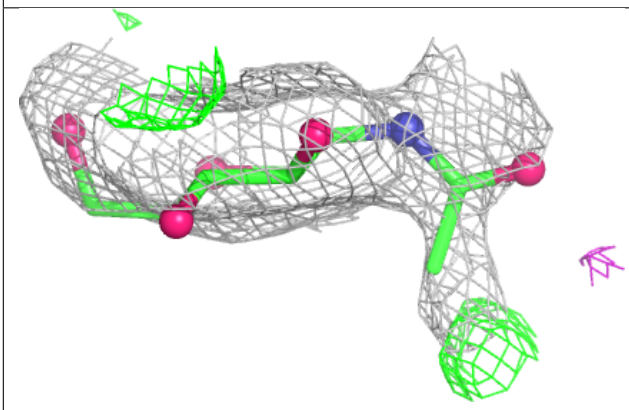


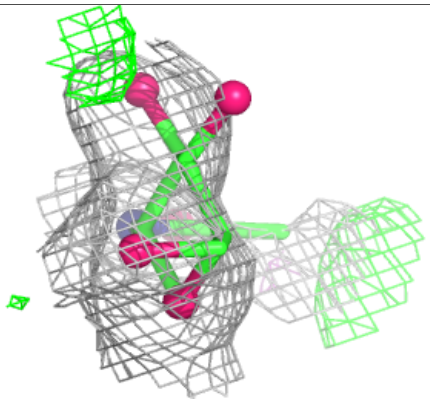


Electron density around NAG A 402:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

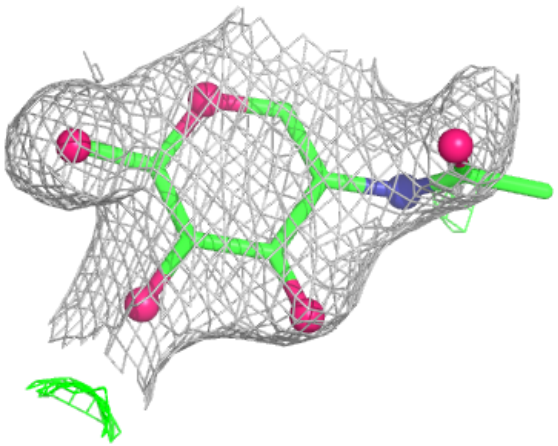


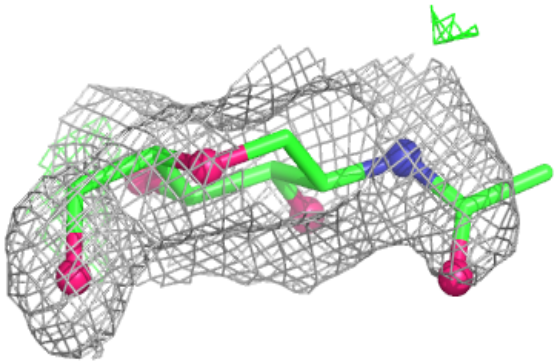


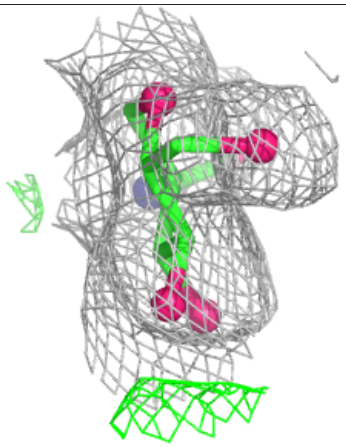


Electron density around NAG A 403:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

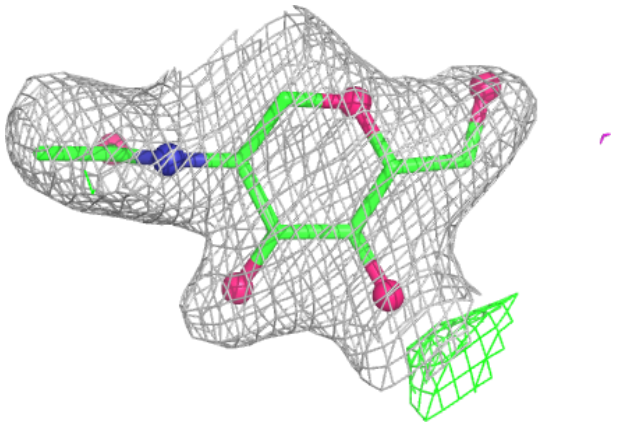


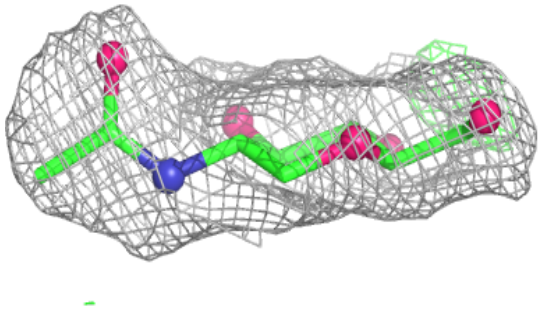


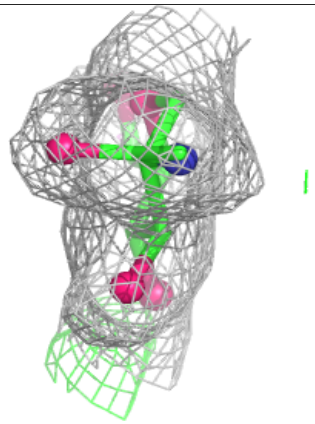


Electron density around NAG C 401:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

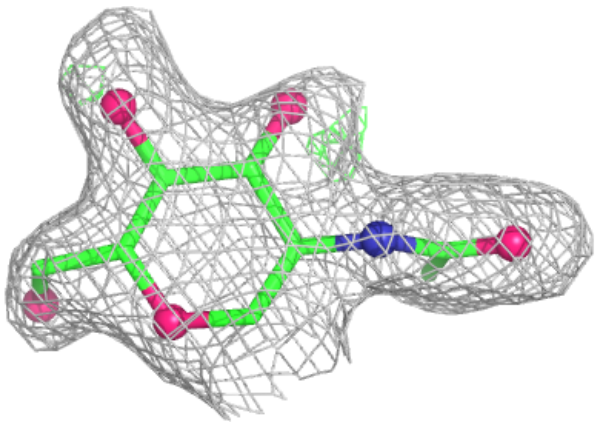


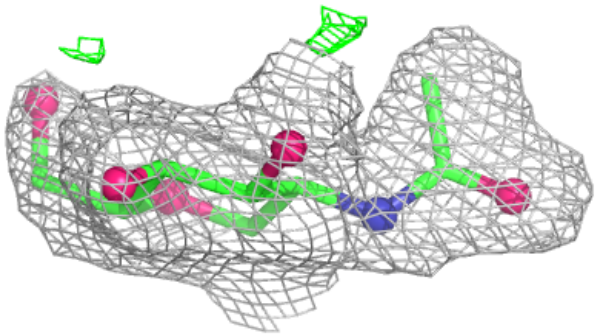


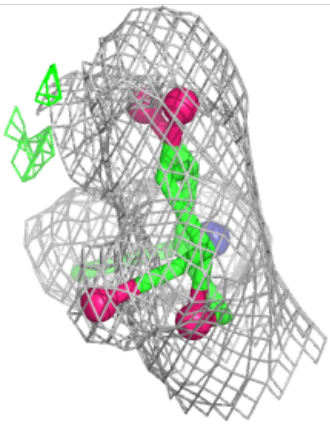


Electron density around NAG A 401:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)







6.5 Other polymers (i)

There are no such residues in this entry.