Full wwPDB X-ray Structure Validation Report (i)

Oct 17, 2023 – 07:11 PM EDT

PDB ID	:	1V3E
Title	:	Structure of the hemagglutinin-neuraminidase from human parainfluenza virus type III: complex with ZANAMAVIR
Authors	:	Lawrence, M.C.; Borg, N.A.; Streltsov, V.A.; Pilling, P.A.; Epa, V.C.; Varghese, J.N.; McKimm-Breschkin, J.L.; Colman, P.M.
Deposited on	:	2003-10-30
Resolution	:	1.89 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 (i) 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 (i)) were used in the production of this report:

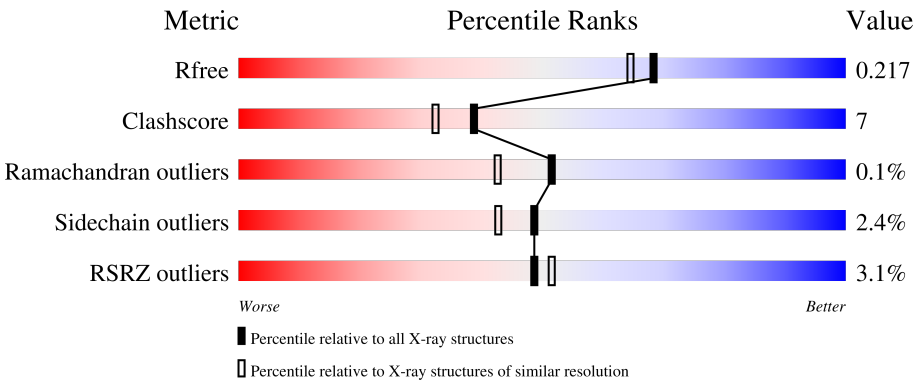
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 1.89 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	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

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	431	4% 85% 13% •
1	B	431	3% 88% 10% •
2	C	2	50% 50%
2	D	2	50% 50%

2 Entry composition (i)

There are 7 unique types of molecules in this entry. The entry contains 7921 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.

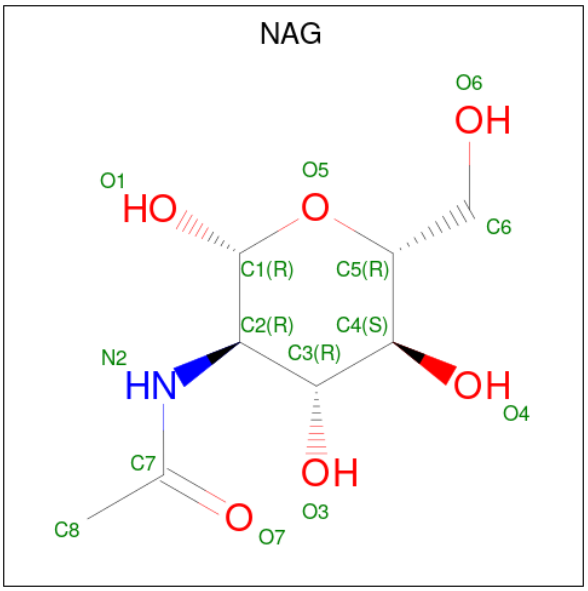
- Molecule 1 is a protein called hemagglutinin-neuraminidase glycoprotein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	431	3383	2142	581	640	20	0	1	0
1	B	431	3383	2142	581	640	20	0	1	0

- Molecule 2 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.

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

- Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).

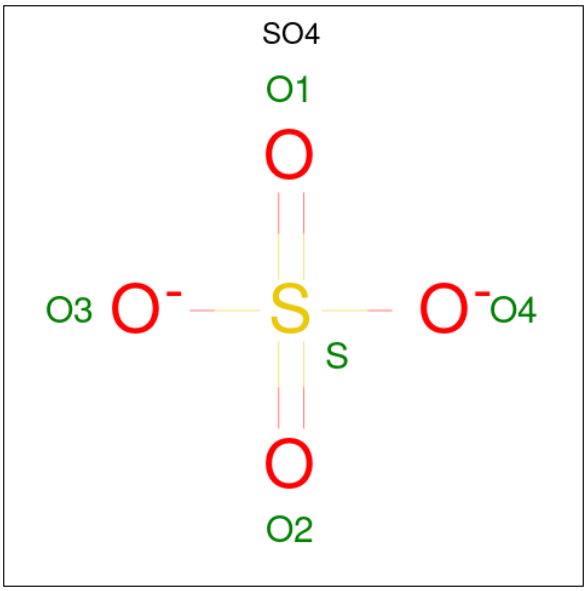


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
3	A	1	14	8	1	5	0	0
			Total	C	N	O		
3	B	1	14	8	1	5	0	0
			Total	C	N	O		
3	B	1	14	8	1	5	0	0

- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

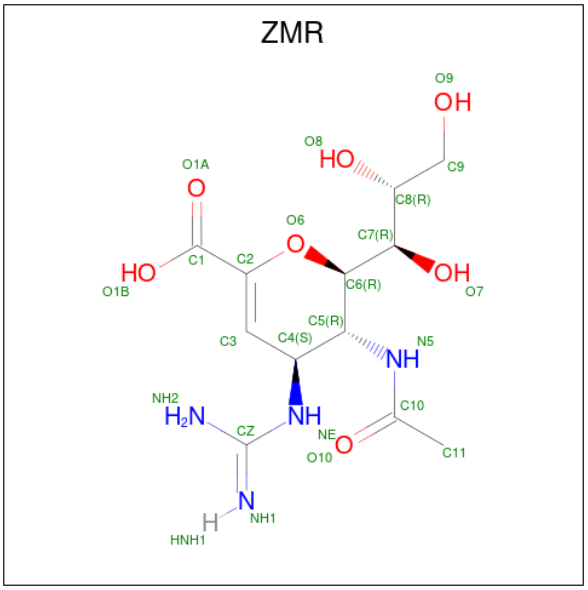
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Ca		
4	A	1	1	1	0	0
			Total	Ca		
4	B	1	1	1	0	0

- Molecule 5 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	S		
5	A	1	5	4	1	0	0
			Total	O	S		
5	B	1	5	4	1	0	0

- Molecule 6 is ZANAMIVIR (three-letter code: ZMR) (formula: $C_{12}H_{20}N_4O_7$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
6	A	1	23	12	4	7	0	0
			Total	C	N	O		
6	B	1	23	12	4	7	0	0

- Molecule 7 is water.

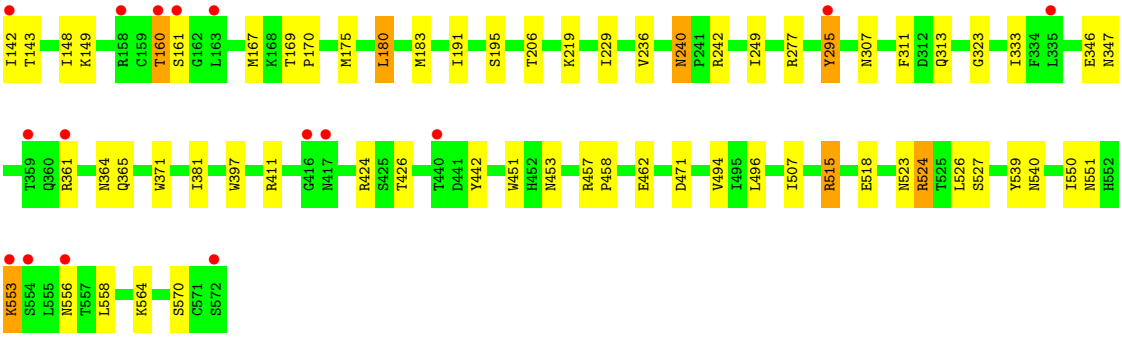
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
7	A	468	468	468	0	0
			Total	O		
7	B	531	531	531	0	0

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.

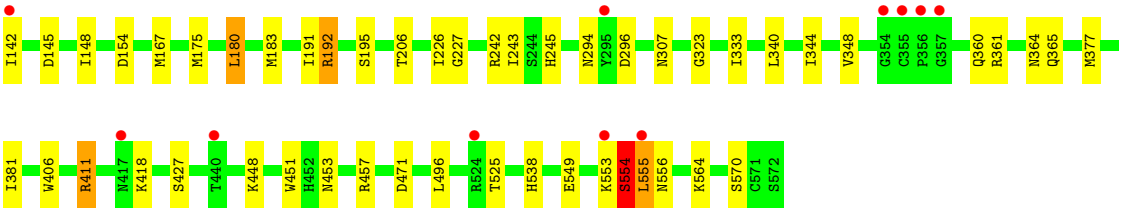
• Molecule 1: hemagglutinin-neuraminidase glycoprotein

Chain A:



• Molecule 1: hemagglutinin-neuraminidase glycoprotein

Chain B:



• Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain C:

• Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain D:

NAG1
NAG2

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 62 2 2	Depositor
Cell constants a, b, c, α, β, γ	218.45Å 218.45Å 109.90Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.97 – 1.89 29.97 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.7 (29.97-1.89) 99.8 (29.97-1.90)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	0.15	Depositor
$< I/\sigma(I) >$ [1]	2.28 (at 1.91Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.184 , 0.219 0.180 , 0.217	Depositor DCC
R_{free} test set	6391 reflections (5.31%)	wwPDB-VP
Wilson B-factor (Å^2)	27.4	Xtriage
Anisotropy	0.177	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 62.4	EDS
L-test for twinning[2]	$< \|L\| > = 0.51$, $< L^2 > = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	7921	wwPDB-VP
Average B, all atoms (Å^2)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.18% 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 (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: CA, ZMR, NAG, SO4

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	0.82	0/3469	0.94	2/4730 (0.0%)
1	B	0.85	1/3469 (0.0%)	0.94	4/4730 (0.1%)
All	All	0.83	1/6938 (0.0%)	0.94	6/9460 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	175	MET	SD-CE	-6.41	1.42	1.77

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	515	ARG	NE-CZ-NH2	-10.69	114.96	120.30
1	A	515	ARG	NE-CZ-NH1	8.45	124.53	120.30
1	B	411	ARG	NE-CZ-NH2	-5.92	117.34	120.30
1	B	496	LEU	N-CA-C	-5.50	96.14	111.00
1	B	538	HIS	N-CA-C	-5.32	96.64	111.00
1	B	175	MET	CG-SD-CE	-5.18	91.92	100.20

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	3383	0	3349	64	0
1	B	3383	0	3348	45	1
2	C	28	0	25	1	0
2	D	28	0	25	1	0
3	A	14	0	13	0	0
3	B	28	0	26	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	5	0	0	0	0
5	B	5	0	0	0	0
6	A	23	0	19	0	0
6	B	23	0	19	0	0
7	A	468	0	0	17	0
7	B	531	0	0	11	1
All	All	7921	0	6824	103	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 7.

All (103) 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:195:SER:OG	7:A:4201:HOH:O	1.73	1.05
1:A:524:ARG:HD2	1:A:524:ARG:H	1.33	0.94
1:B:360:GLN:HE21	1:B:364:ASN:HD21	1.26	0.84
1:A:553:LYS:HA	7:A:4441:HOH:O	1.76	0.83
1:A:142:ILE:HG13	7:A:4452:HOH:O	1.82	0.80
1:A:524:ARG:H	1:A:524:ARG:CD	1.94	0.78
1:A:142:ILE:HG23	1:A:143:THR:HG23	1.66	0.78
1:B:360:GLN:HE21	1:B:364:ASN:ND2	1.83	0.76
1:B:365:GLN:HG2	7:B:5664:HOH:O	1.85	0.76
1:A:462:GLU:OE1	7:A:4467:HOH:O	2.02	0.76
1:B:361:ARG:O	1:B:365:GLN:HG3	1.86	0.75
1:B:525:THR:O	1:B:553:LYS:HE2	1.87	0.74
1:A:160:THR:HG23	1:A:161:SER:H	1.51	0.73
1:A:556:ASN:ND2	7:A:4441:HOH:O	2.15	0.72
1:A:551:ASN:ND2	7:A:4469:HOH:O	2.23	0.71
1:A:180:LEU:HD13	1:B:180:LEU:HD13	1.73	0.70
1:A:539:TYR:O	1:A:540:ASN:HB2	1.93	0.69
1:B:294:ASN:HB2	1:B:296:ASP:OD1	1.94	0.68
1:A:277:ARG:HD2	1:A:371:TRP:CZ2	2.29	0.67
1:A:160:THR:HG22	7:A:4449:HOH:O	1.95	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:A:4253:HOH:O	1:B:554:SER:HB3	1.93	0.66
1:B:555:LEU:O	1:B:556:ASN:HB3	1.94	0.66
1:B:418:LYS:HE3	7:B:5636:HOH:O	1.95	0.66
1:A:553:LYS:HB2	1:A:553:LYS:NZ	2.13	0.64
1:B:243:ILE:HD12	1:B:245[B]:HIS:NE2	2.12	0.64
1:A:277:ARG:HD3	7:A:4459:HOH:O	1.97	0.62
1:A:160:THR:HG23	1:A:161:SER:N	2.14	0.62
1:A:142:ILE:N	7:A:4295:HOH:O	2.34	0.60
1:A:307:ASN:HB2	7:A:4398:HOH:O	2.00	0.60
1:A:553:LYS:HG2	7:A:4441:HOH:O	2.03	0.58
1:B:554:SER:HB2	1:B:555:LEU:HD22	1.84	0.58
1:B:226:ILE:HD12	1:B:243:ILE:HD11	1.85	0.57
1:A:524:ARG:HD2	1:A:524:ARG:N	2.12	0.57
1:A:183:MET:HE3	1:B:183:MET:HE3	1.87	0.56
1:B:553:LYS:HA	7:B:5727:HOH:O	2.06	0.56
1:A:539:TYR:O	1:A:540:ASN:CB	2.53	0.56
1:B:142:ILE:N	7:B:5707:HOH:O	2.40	0.55
1:A:515:ARG:HD2	1:A:518:GLU:OE1	2.07	0.54
1:A:277:ARG:HD2	1:A:371:TRP:CE2	2.43	0.54
1:A:551:ASN:HA	1:A:558:LEU:HD23	1.90	0.54
1:A:527:SER:HB2	1:A:551:ASN:HB3	1.90	0.53
1:A:149:LYS:HD2	7:A:4366:HOH:O	2.07	0.53
1:A:183:MET:CE	1:B:183:MET:HE3	2.39	0.53
1:A:183:MET:CE	1:B:183:MET:CE	2.88	0.52
1:B:555:LEU:HD22	1:B:555:LEU:H	1.75	0.52
1:A:183:MET:HE3	1:B:183:MET:CE	2.41	0.50
1:B:377:MET:HG3	1:B:406:TRP:CZ2	2.47	0.50
1:B:555:LEU:O	1:B:556:ASN:CB	2.59	0.49
1:A:195:SER:HB3	1:A:206:THR:HG22	1.94	0.49
1:A:361:ARG:O	1:A:365:GLN:HG3	2.12	0.49
1:B:427:SER:HB3	7:B:5320:HOH:O	2.13	0.48
1:A:175:MET:SD	1:A:236:VAL:HG21	2.54	0.48
1:A:515:ARG:CD	1:A:518:GLU:OE1	2.62	0.47
1:A:323:GLY:HA3	1:A:411:ARG:HB3	1.95	0.47
1:B:360:GLN:NE2	1:B:364:ASN:HD21	2.04	0.47
1:B:333:ILE:CG2	1:B:381:ILE:HG23	2.45	0.47
1:B:451:TRP:CH2	1:B:453:ASN:HB2	2.50	0.47
1:A:169:THR:HA	1:A:170:PRO:C	2.35	0.47
1:A:397:TRP:CD2	1:A:442:TYR:HB3	2.50	0.47
1:A:523:ASN:HB2	1:A:524:ARG:NH1	2.30	0.47
7:A:4270:HOH:O	2:C:1:NAG:H81	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:340:LEU:HD13	1:B:344:ILE:HD12	1.97	0.47
2:D:1:NAG:H62	2:D:2:NAG:C8	2.44	0.46
1:B:525:THR:O	1:B:553:LYS:CE	2.61	0.46
1:A:160:THR:CG2	1:A:161:SER:H	2.26	0.46
1:B:167:MET:HG2	7:B:5652:HOH:O	2.15	0.46
1:B:471:ASP:HB2	7:B:5509:HOH:O	2.15	0.46
1:A:277:ARG:HD2	1:A:371:TRP:CH2	2.51	0.46
1:A:553:LYS:HB2	1:A:553:LYS:HZ3	1.79	0.46
1:A:523:ASN:HB2	1:A:524:ARG:CZ	2.46	0.46
1:A:553:LYS:CA	7:A:4441:HOH:O	2.52	0.45
1:A:240:ASN:HD22	1:A:240:ASN:C	2.20	0.44
1:A:180:LEU:HD13	1:B:180:LEU:CD1	2.44	0.44
1:B:183:MET:HE3	1:B:183:MET:HB3	1.88	0.44
1:A:451:TRP:CH2	1:A:453:ASN:HB2	2.54	0.43
1:A:180:LEU:HD22	1:B:180:LEU:HD22	2.01	0.43
1:A:424:ARG:HG2	1:A:426:THR:OG1	2.18	0.43
1:B:340:LEU:CD1	1:B:344:ILE:HD12	2.49	0.43
1:A:167:MET:HG2	1:A:570:SER:HB3	2.00	0.43
1:A:311:PHE:HB3	1:A:313:GLN:O	2.19	0.43
1:A:526:LEU:HD11	1:A:550:ILE:HG23	2.00	0.42
1:A:149:LYS:HB3	1:A:149:LYS:HE2	1.77	0.42
1:A:346:GLU:HG2	7:A:4410:HOH:O	2.19	0.42
1:A:160:THR:CG2	1:A:161:SER:N	2.82	0.42
1:B:307:ASN:HB3	7:B:5713:HOH:O	2.20	0.42
1:B:348:VAL:HG22	1:B:364:ASN:HD22	1.85	0.41
1:B:145:ASP:HB2	1:B:148:ILE:HD12	2.02	0.41
1:B:192:ARG:NH2	7:B:5620:HOH:O	2.40	0.41
1:B:192:ARG:HB2	1:B:549:GLU:OE2	2.19	0.41
1:A:295:TYR:HD1	1:A:295:TYR:HA	1.81	0.41
1:A:307:ASN:ND2	7:A:4425:HOH:O	2.54	0.41
1:B:323:GLY:HA3	1:B:411:ARG:HB3	2.01	0.41
1:A:333:ILE:CG2	1:A:381:ILE:HG23	2.50	0.41
1:B:195:SER:HB3	1:B:206:THR:HG22	2.03	0.41
1:A:180:LEU:CD1	1:B:180:LEU:HD13	2.48	0.41
1:A:494:VAL:HG22	1:A:507:ILE:HG12	2.03	0.41
1:B:227:GLY:HA2	1:B:242:ARG:HG2	2.03	0.41
1:A:458:PRO:HD3	1:A:496:LEU:O	2.21	0.41
1:B:448:LYS:HE2	7:B:5631:HOH:O	2.20	0.41
1:A:219:LYS:HB2	1:A:249:ILE:HG12	2.04	0.40
1:A:347:ASN:ND2	1:A:364:ASN:HD21	2.18	0.40
1:B:570:SER:HB2	7:B:5579:HOH:O	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:148:ILE:HD13	1:A:229:ILE:HG22	2.03	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)
1:B:296:ASP:OD2	1:B:296:ASP:OD2[10_665]	1.51	0.69
7:B:5722:HOH:O	7:B:5722:HOH:O[10_665]	1.86	0.34

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	430/431 (100%)	407 (95%)	23 (5%)	0	100	100
1	B	430/431 (100%)	409 (95%)	20 (5%)	1 (0%)	47	38
All	All	860/862 (100%)	816 (95%)	43 (5%)	1 (0%)	51	42

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	554	SER

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	391/390 (100%)	380 (97%)	11 (3%)	43	36
1	B	391/390 (100%)	383 (98%)	8 (2%)	55	51
All	All	782/780 (100%)	763 (98%)	19 (2%)	49	43

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	160	THR
1	A	180	LEU
1	A	191	ILE
1	A	240	ASN
1	A	242	ARG
1	A	295	TYR
1	A	457	ARG
1	A	471	ASP
1	A	524	ARG
1	A	553	LYS
1	A	564	LYS
1	B	154	ASP
1	B	180	LEU
1	B	191	ILE
1	B	192	ARG
1	B	457	ARG
1	B	554	SER
1	B	555	LEU
1	B	564	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	240	ASN
1	A	262	ASN
1	A	347	ASN
1	A	365	GLN
1	A	504	ASN
1	A	556	ASN
1	B	240	ASN
1	B	250	ASN
1	B	262	ASN
1	B	307	ASN
1	B	347	ASN

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Mol	Chain	Res	Type
1	B	364	ASN
1	B	504	ASN
1	B	556	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)

4 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	$\#\|Z\| > 2$
2	NAG	C	1	1,2	14,14,15	0.63	0	17,19,21	1.37	2 (11%)
2	NAG	C	2	2	14,14,15	0.58	0	17,19,21	0.86	0
2	NAG	D	1	1,2	14,14,15	0.82	0	17,19,21	1.14	2 (11%)
2	NAG	D	2	2	14,14,15	0.76	0	17,19,21	0.72	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	C	1	1,2	-	4/6/23/26	0/1/1/1
2	NAG	C	2	2	-	0/6/23/26	0/1/1/1
2	NAG	D	1	1,2	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	D	2	2	-	3/6/23/26	0/1/1/1

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	C	1	NAG	C2-N2-C7	-3.37	118.11	122.90
2	C	1	NAG	C4-C3-C2	-3.28	106.20	111.02
2	D	1	NAG	C2-N2-C7	-2.72	119.04	122.90
2	D	1	NAG	O5-C1-C2	-2.23	107.77	111.29

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	1	NAG	C8-C7-N2-C2
2	C	1	NAG	O7-C7-N2-C2
2	C	1	NAG	O5-C5-C6-O6
2	C	1	NAG	C4-C5-C6-O6
2	D	2	NAG	C8-C7-N2-C2
2	D	2	NAG	O7-C7-N2-C2
2	D	2	NAG	C3-C2-N2-C7

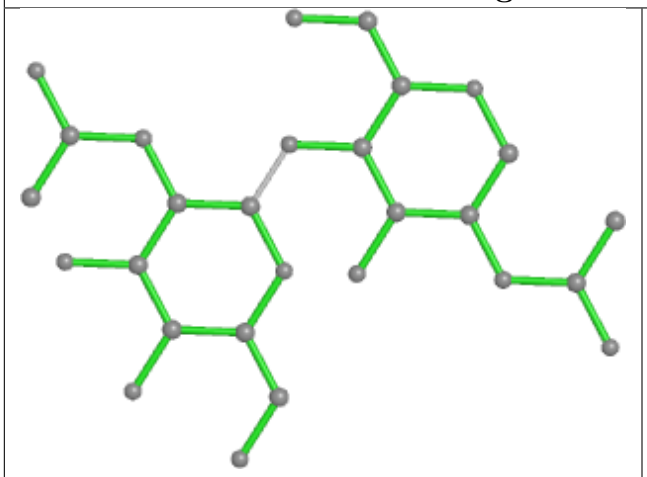
There are no ring outliers.

3 monomers are involved in 2 short contacts:

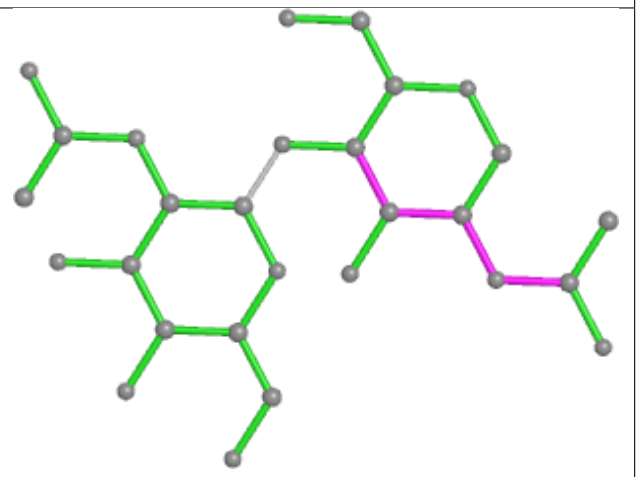
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	1	NAG	1	0
2	D	1	NAG	1	0
2	D	2	NAG	1	0

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

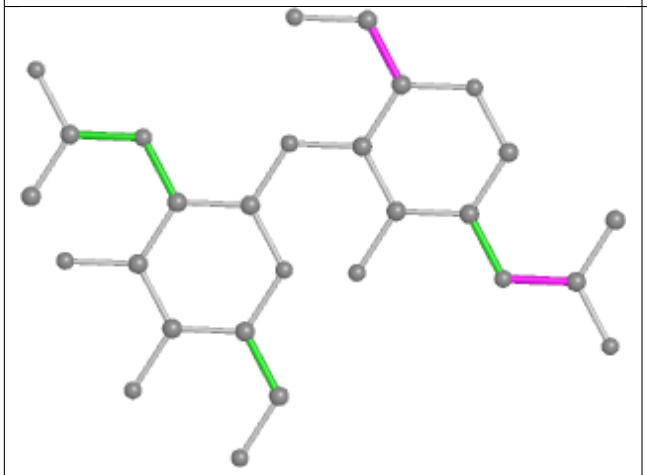
Oligosaccharide Chain C



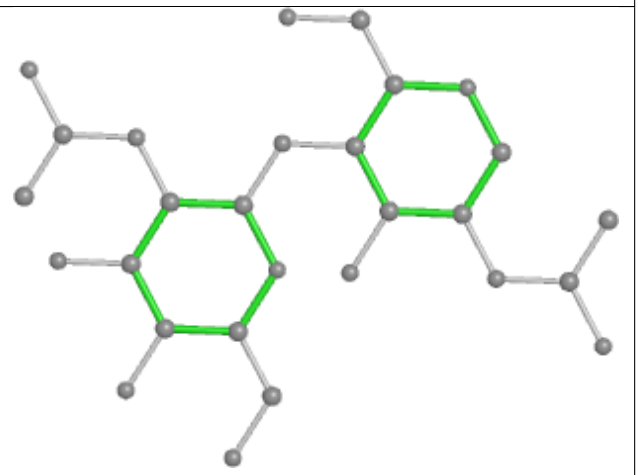
Bond lengths



Bond angles

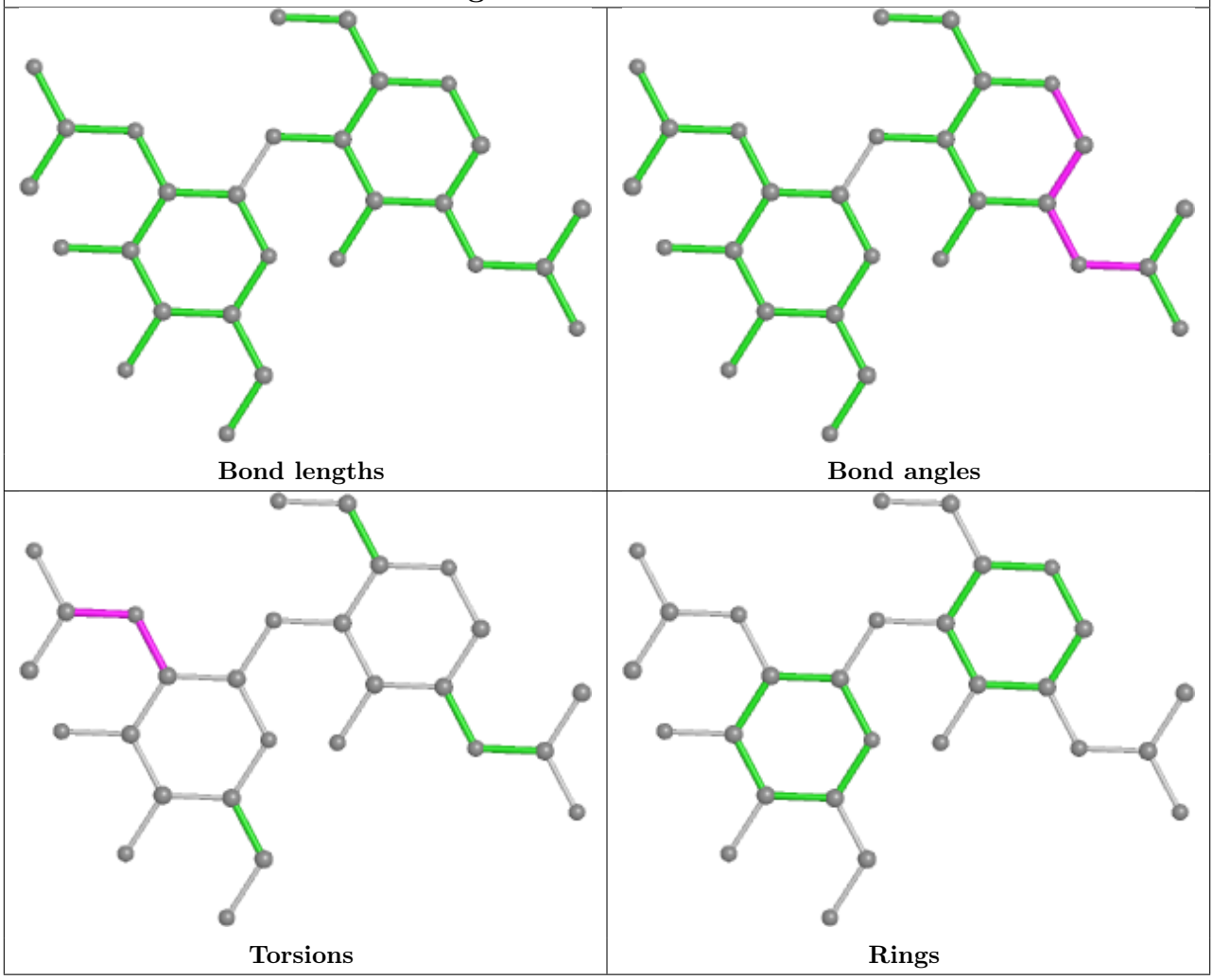


Torsions



Rings

Oligosaccharide Chain D



5.6 Ligand geometry (i)

Of 9 ligands modelled in this entry, 2 are monoatomic - leaving 7 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	$\#\|Z\| > 2$
6	ZMR	A	1200	-	22,23,23	2.10	7 (31%)	24,32,32	1.67	5 (20%)
3	NAG	B	5231	1	14,14,15	0.74	0	17,19,21	0.73	1 (5%)
3	NAG	A	3081	1	14,14,15	0.62	0	17,19,21	0.70	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
5	SO4	A	3002	-	4,4,4	0.27	0	6,6,6	0.16	0
5	SO4	B	3001	-	4,4,4	0.26	0	6,6,6	0.09	0
3	NAG	B	3081	1	14,14,15	0.65	0	17,19,21	0.72	0
6	ZMR	B	2200	-	22,23,23	2.07	7 (31%)	24,32,32	1.82	5 (20%)

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
6	ZMR	A	1200	-	-	0/22/38/38	0/1/1/1
3	NAG	B	5231	1	-	2/6/23/26	0/1/1/1
3	NAG	A	3081	1	-	2/6/23/26	0/1/1/1
3	NAG	B	3081	1	-	0/6/23/26	0/1/1/1
6	ZMR	B	2200	-	-	0/22/38/38	0/1/1/1

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	2200	ZMR	C7-C6	5.49	1.59	1.53
6	A	1200	ZMR	C7-C6	5.00	1.59	1.53
6	A	1200	ZMR	C3-C2	4.93	1.40	1.33
6	B	2200	ZMR	C3-C2	4.87	1.40	1.33
6	A	1200	ZMR	CZ-NE	3.40	1.39	1.33
6	B	2200	ZMR	C2-C1	-2.88	1.41	1.48
6	A	1200	ZMR	C11-C10	2.70	1.56	1.50
6	A	1200	ZMR	O1B-C1	-2.47	1.23	1.30
6	B	2200	ZMR	C9-C8	2.26	1.58	1.52
6	A	1200	ZMR	C2-C1	-2.24	1.43	1.48
6	B	2200	ZMR	O1B-C1	-2.20	1.24	1.30
6	B	2200	ZMR	CZ-NE	2.20	1.37	1.33
6	A	1200	ZMR	C9-C8	2.10	1.57	1.52
6	B	2200	ZMR	O7-C7	2.02	1.47	1.43

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
6	B	2200	ZMR	O6-C2-C3	-5.53	118.16	124.61
6	A	1200	ZMR	O6-C2-C3	-5.02	118.75	124.61
6	B	2200	ZMR	C8-C7-C6	-3.49	106.42	113.03

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
6	A	1200	ZMR	O1B-C1-C2	3.34	122.52	114.20
6	B	2200	ZMR	O1B-C1-C2	3.00	121.69	114.20
6	B	2200	ZMR	C6-C5-N5	-2.59	106.60	110.91
6	A	1200	ZMR	O6-C2-C1	2.52	117.10	112.06
6	A	1200	ZMR	C8-C7-C6	-2.45	108.39	113.03
6	B	2200	ZMR	O6-C2-C1	2.43	116.93	112.06
6	A	1200	ZMR	C6-C5-N5	-2.09	107.45	110.91
3	B	5231	NAG	C2-N2-C7	-2.06	119.97	122.90

There are no chirality outliers.

All (4) torsion outliers are listed below:

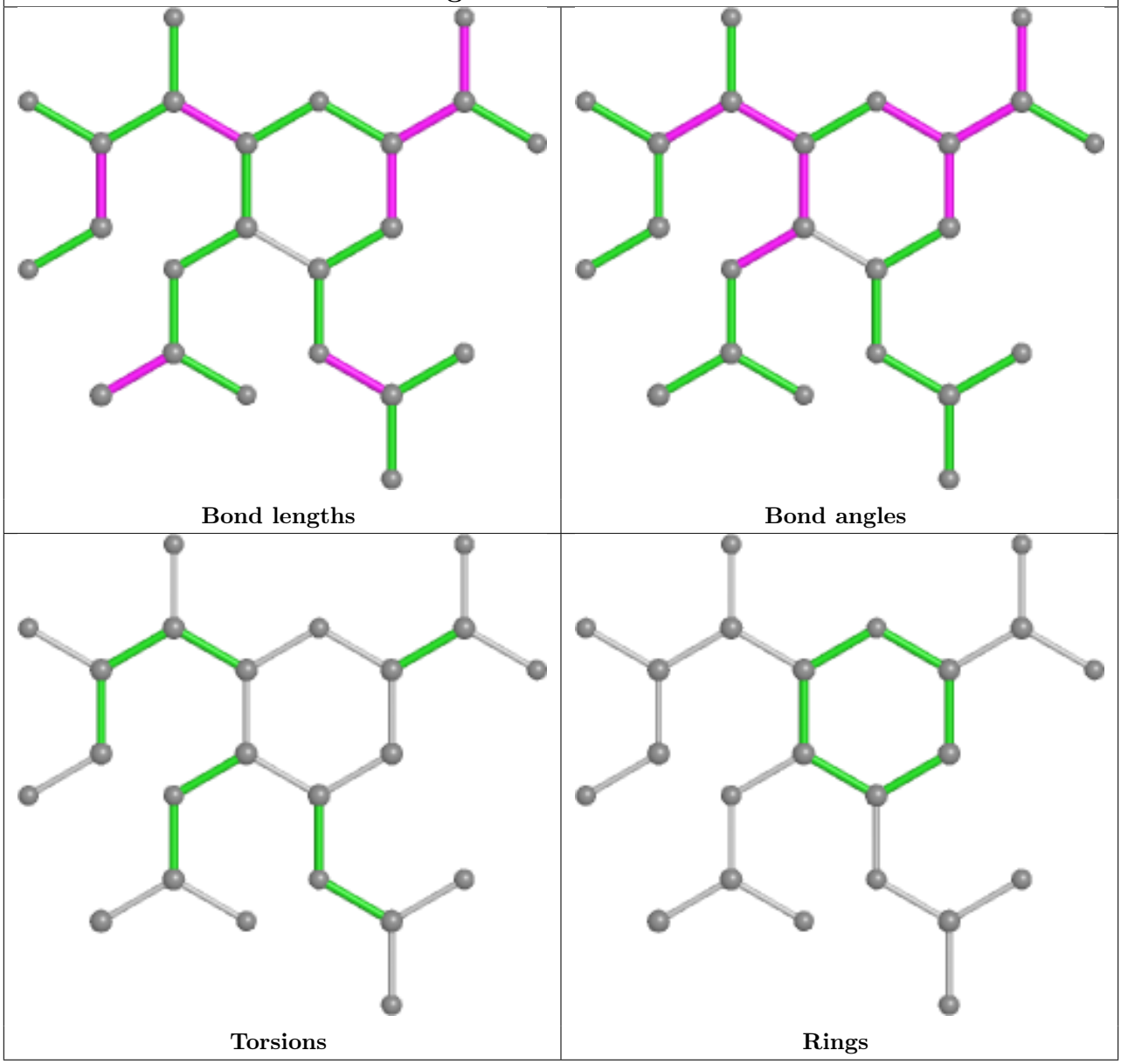
Mol	Chain	Res	Type	Atoms
3	B	5231	NAG	C8-C7-N2-C2
3	B	5231	NAG	O7-C7-N2-C2
3	A	3081	NAG	C8-C7-N2-C2
3	A	3081	NAG	O7-C7-N2-C2

There are no ring outliers.

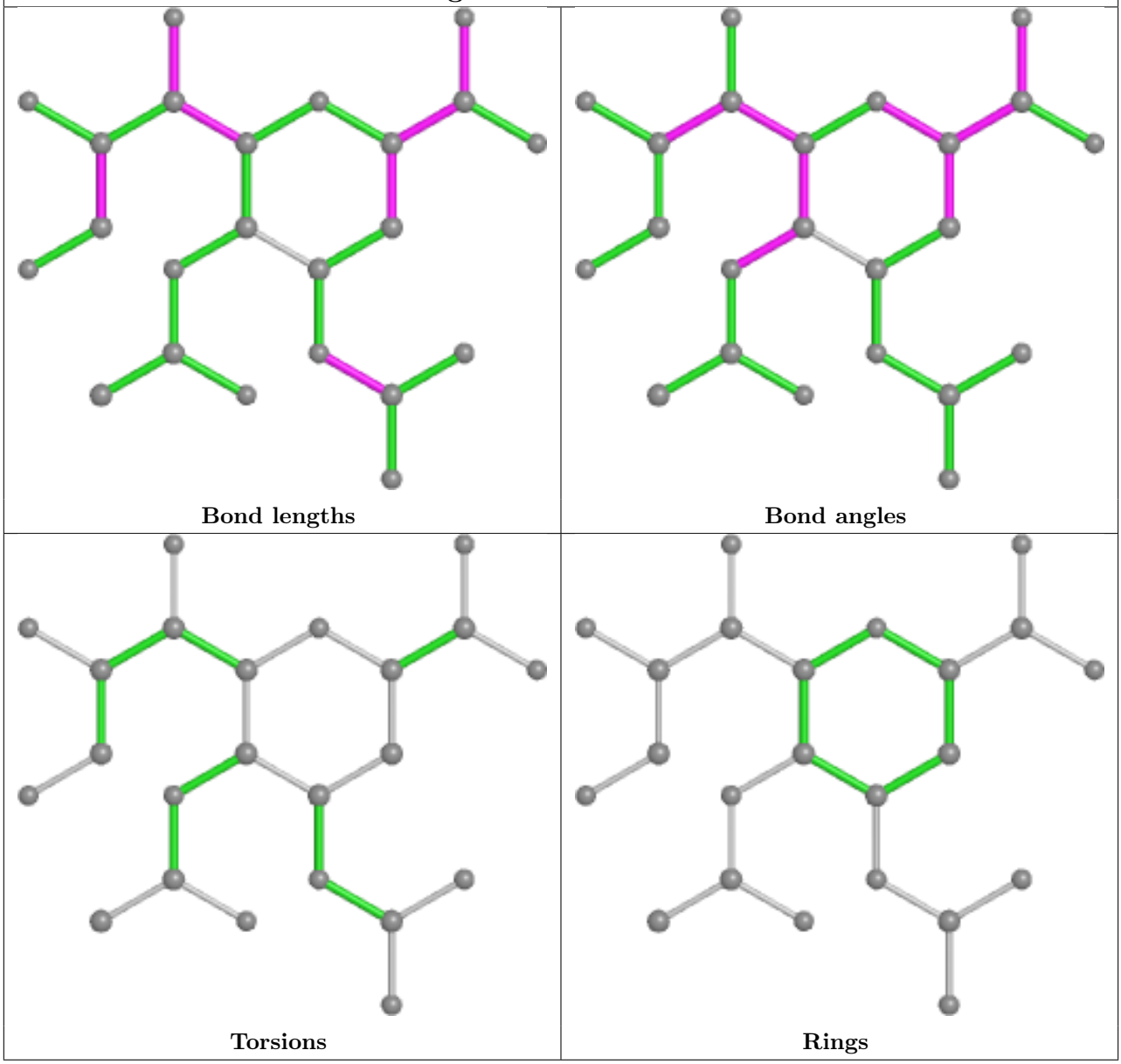
No monomer is involved in short contacts.

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 ZMR A 1200



Ligand ZMR B 2200



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	431/431 (100%)	-0.05	16 (3%)	41	44	18, 29, 53, 82	0
1	B	431/431 (100%)	-0.13	11 (2%)	56	58	18, 28, 51, 73	0
All	All	862/862 (100%)	-0.09	27 (3%)	49	51	18, 28, 53, 82	0

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	553	LYS	5.2
1	A	160	THR	4.8
1	B	142	ILE	4.3
1	A	163	LEU	3.9
1	A	572	SER	3.8
1	B	295	TYR	3.7
1	A	295	TYR	3.6
1	A	161	SER	3.2
1	A	416	GLY	3.1
1	A	142	ILE	3.0
1	A	440	THR	2.9
1	A	158	ARG	2.9
1	B	355	CYS	2.8
1	A	417	ASN	2.8
1	B	357	GLY	2.7
1	A	556	ASN	2.6
1	A	361	ARG	2.6
1	B	356	PRO	2.6
1	A	359	THR	2.5
1	B	555	LEU	2.5
1	B	524	ARG	2.5
1	B	417	ASN	2.4
1	B	354	GLY	2.3
1	A	554	SER	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	553	LYS	2.1
1	B	440	THR	2.1
1	A	335	LEU	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)

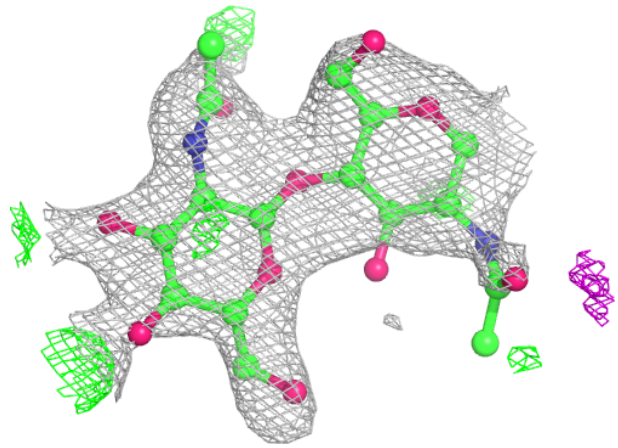
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

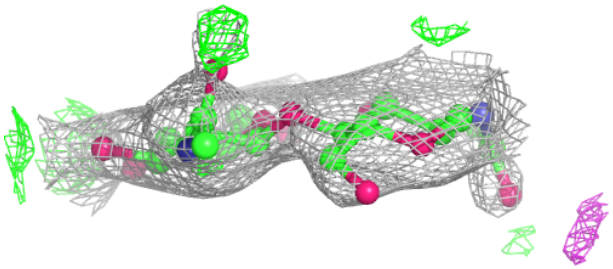
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	NAG	D	2	14/15	0.82	0.22	65,74,79,80	0
2	NAG	C	2	14/15	0.85	0.20	75,81,85,87	0
2	NAG	C	1	14/15	0.86	0.21	68,78,85,85	0
2	NAG	D	1	14/15	0.91	0.15	67,69,74,75	0

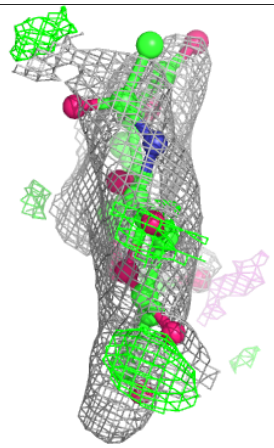
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Chain C:

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

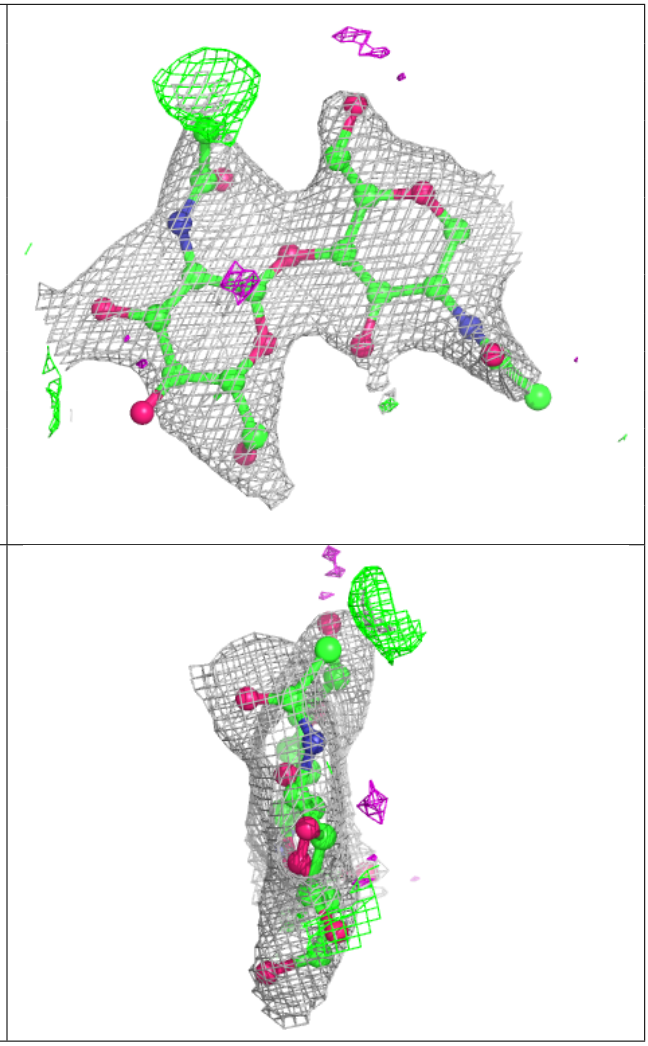






Electron density around Chain D:

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



6.4 Ligands (i)

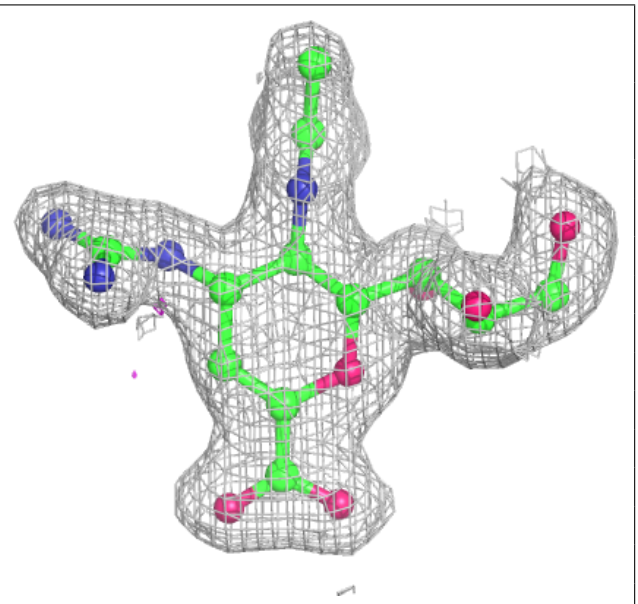
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

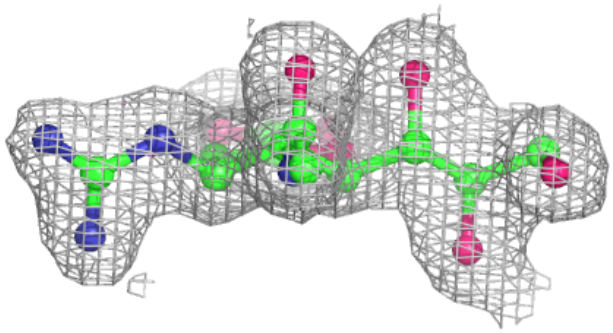
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	NAG	A	3081	14/15	0.81	0.38	59,67,73,78	0
3	NAG	B	5231	14/15	0.89	0.23	47,60,70,70	0
3	NAG	B	3081	14/15	0.90	0.23	42,47,55,59	0
5	SO4	A	3002	5/5	0.91	0.19	91,93,97,98	0
5	SO4	B	3001	5/5	0.95	0.21	82,82,88,88	0
4	CA	A	4001	1/1	0.96	0.04	32,32,32,32	0
6	ZMR	A	1200	23/23	0.97	0.07	20,25,30,33	0
4	CA	B	2001	1/1	0.98	0.05	26,26,26,26	0
6	ZMR	B	2200	23/23	0.98	0.07	18,23,31,32	0

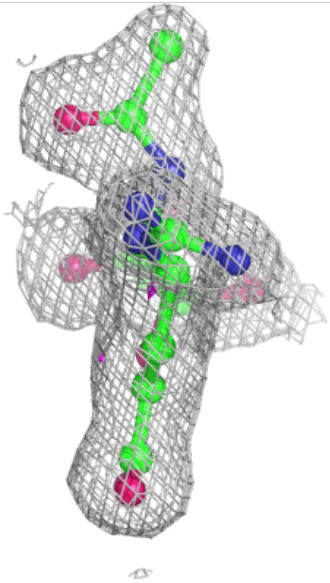
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around ZMR A 1200:

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

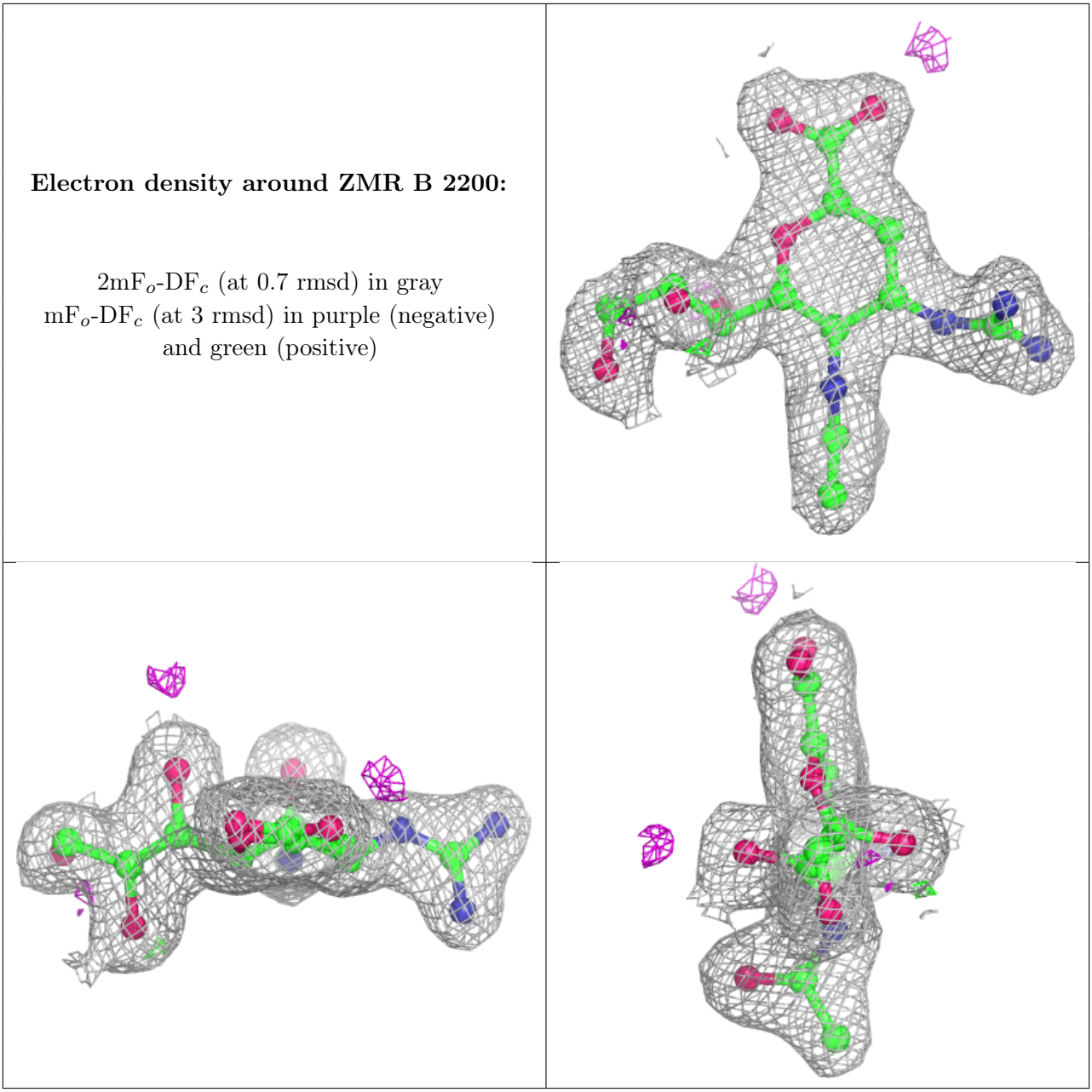






Electron density around ZMR B 2200:

$2mF_o$-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.