Full wwPDB X-ray Structure Validation Report (i)

Aug 6, 2020 – 11:52 PM BST

PDB ID	:	6G01
Title	:	Complex of neuraminidase from H1N1 influenza virus with tamiphosphor monomethyl ester
Authors	:	Pachl, P.; Pokorna, J.
Deposited on	:	2018-03-15
Resolution	:	1.61 Å(reported)

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

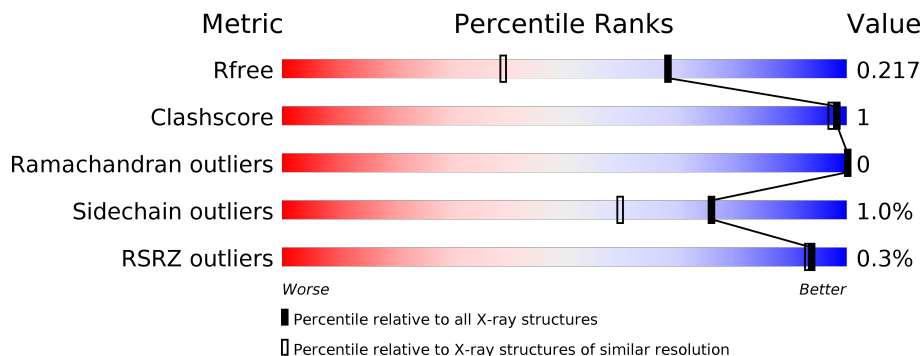|		
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.13.1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.13.1

1 Overall quality at a glance ⓘ

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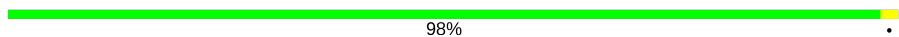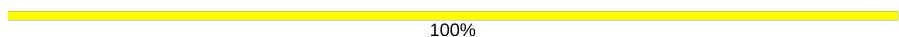*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.61 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	Percentile Ranks	Value
Rfree		0.217
Clashscore		1
Ramachandran outliers		0
Sidechain outliers		1.0%
RSRZ outliers		0.3%

Worse → *Better*

■ Percentile relative to all X-ray structures

▯ Percentile relative to X-ray structures of similar resolution

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	4693 (1.64-1.60)
Clashscore	141614	5002 (1.64-1.60)
Ramachandran outliers	138981	4888 (1.64-1.60)
Sidechain outliers	138945	4887 (1.64-1.60)
RSRZ outliers	127900	4609 (1.64-1.60)

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 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	387	98% ·
1	B	387	98% ·
2	C	2	100%
2	E	2	100%
3	D	3	33% 67%
4	F	4	100%

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	C	2	-	-	-	X
3	NAG	D	2	-	-	-	X
3	BMA	D	3	-	-	-	X
7	NAG	B	510	-	-	-	X

2 Entry composition (i)

There are 9 unique types of molecules in this entry. The entry contains 7152 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 Neuraminidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	387	Total 3030	C 1901	N 522	O 586	S 21	0	6	0
1	B	387	Total 3026	C 1900	N 523	O 582	S 21	0	6	0

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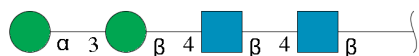

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	2	Total 28	C 16	N 2	O 10	0	0	0
2	E	2	Total 28	C 16	N 2	O 10	0	0	0

- Molecule 3 is an oligosaccharide called beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	D	3	Total 39	C 22	N 2	O 15	0	0	0

- Molecule 4 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.
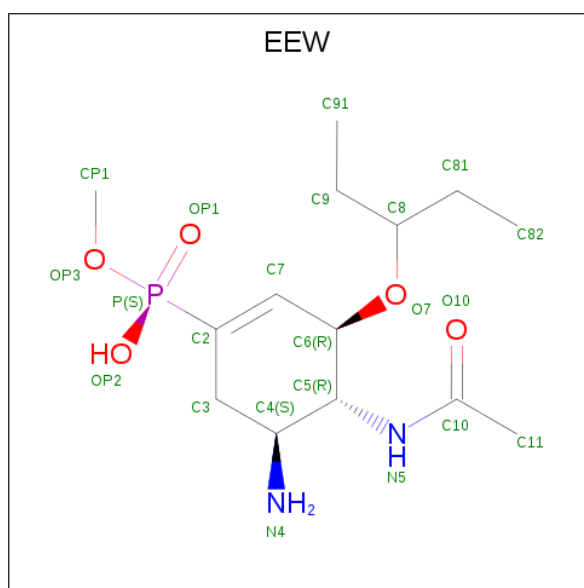
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
4	F	4	50	28	2	20	0	0	0

- Molecule 5 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Ca		
5	B	2	2	2	0	0
			Total	Ca		
5	A	2	2	2	0	0

- Molecule 6 is [(3 {R},4 {R},5 {S})-4-acetamido-5-azanyl-3-pentan-3-yloxy-cyclohexen-1-yl]-methoxy-phosphinic acid (three-letter code: EEW) (formula: $C_{14}H_{27}N_2O_5P$).
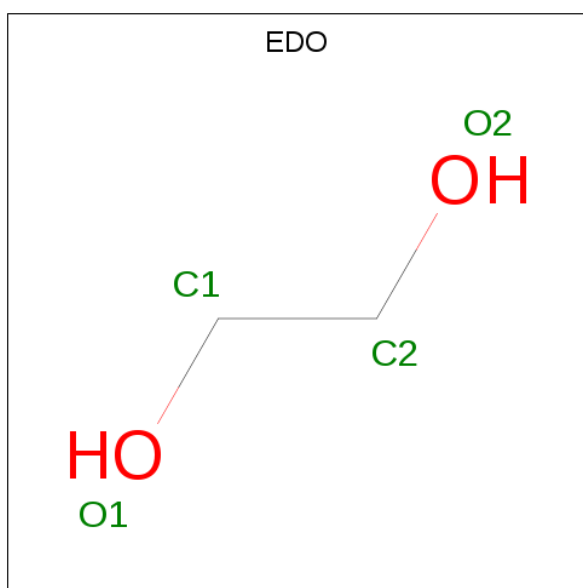

EEW

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
6	A	1	22	14	2	5	1	0	0
			Total	C	N	O	P		
6	B	1	22	14	2	5	1	0	0

- Molecule 7 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).

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

- Molecule 8 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	A	1	Total 4	C 2	O 2	0	0
8	A	1	Total 4	C 2	O 2	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	A	1	Total 4	C 2	O 2	0	0
8	A	1	Total 4	C 2	O 2	0	0
8	B	1	Total 8	C 4	O 4	0	1
8	B	1	Total 4	C 2	O 2	0	0
8	B	1	Total 4	C 2	O 2	0	0

- Molecule 9 is water.

| Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	412	Total 422	O 422	0	10
9	B	412	Total 421	O 421	0	11

3 Residue-property plots ⓘ

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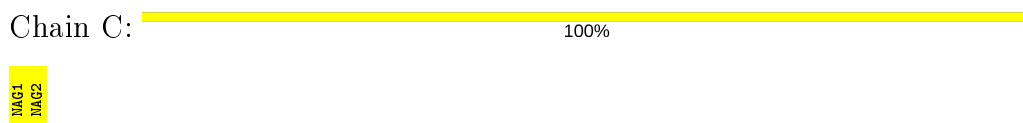
• Molecule 1: Neuraminidase
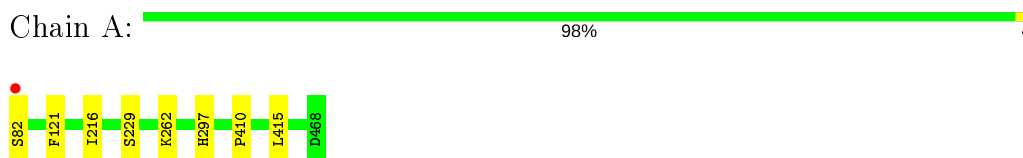
• Molecule 1: Neuraminidase
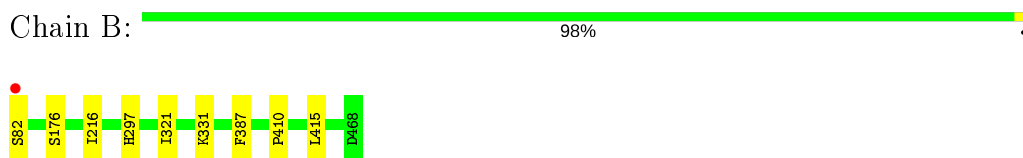
• Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

• Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose
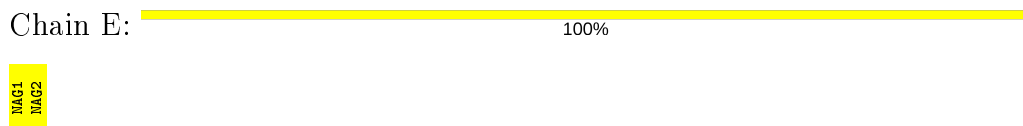
• Molecule 3: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

• Molecule 4: alpha-D-mannopyranose-(1-3)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain F: _____ 100%

NAG1
NAG2
BMA3
MAN4

4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α, β, γ	118.58Å 137.41Å 118.45Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.89 – 1.61 44.89 – 1.61	Depositor EDS
% Data completeness (in resolution range)	98.2 (44.89-1.61) 98.2 (44.89-1.61)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$<I/\sigma(I)>$ [1]	1.81 (at 1.62Å)	Xtriage
Refinement program	REFMAC 5.8.0189	Depositor
R, R_{free}	0.178 , 0.205 0.190 , 0.217	Depositor DCC
R_{free} test set	1821 reflections (1.50%)	wwPDB-VP
Wilson B-factor (Å^2)	17.3	Xtriage
Anisotropy	0.091	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 48.9	EDS
L-test for twinning[2]	$<|L|> = 0.47$, $<L^2> = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	7152	wwPDB-VP
Average B, all atoms (Å^2)	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 24.66 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 3.6778e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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[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: BMA, NAG, EEW, EDO, CA, 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	0.73	0/3118	0.78	0/4238
1	B	0.71	0/3124	0.78	0/4245
All	All	0.72	0/6242	0.78	0/8483

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 (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	3030	0	2860	3	0
1	B	3026	0	2865	3	0
2	C	28	0	25	0	0
2	E	28	0	25	0	0
3	D	39	0	34	0	0
4	F	50	0	43	0	0
5	A	2	0	0	0	0
5	B	2	0	0	0	0
6	A	22	0	0	0	0
6	B	22	0	0	0	0
7	A	14	0	13	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	B	14	0	13	0	0
8	A	16	0	24	1	0
8	B	16	0	24	0	0
9	A	422	0	0	1	0
9	B	421	0	0	2	0
All	All	7152	0	5926	7	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 1.

All (7) 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:262:LYS:NZ	9:A:601:HOH:O	2.24	0.52
8:A:513:EDO:H12	9:B:674:HOH:O	2.15	0.46
1:B:410:PRO:HB3	1:B:415:LEU:O	2.16	0.44
1:B:331:LYS:HE3	9:B:911:HOH:O	2.20	0.41
1:B:321[B]:ILE:CD1	1:B:387:PHE:HB3	2.50	0.41
1:A:121:PHE:CG	1:A:229:SER:HA	2.55	0.41
1:A:410:PRO:HB3	1:A:415:LEU:O	2.21	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	391/387 (101%)	376 (96%)	15 (4%)	0	100	100
1	B	391/387 (101%)	377 (96%)	14 (4%)	0	100	100
All	All	782/774 (101%)	753 (96%)	29 (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	340/334 (102%)	337 (99%)	3 (1%)	78	64
1	B	340/334 (102%)	335 (98%)	5 (2%)	65	43
All	All	680/668 (102%)	672 (99%)	8 (1%)	76	52

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	82	SER
1	A	216	ILE
1	A	297	HIS
1	B	82	SER
1	B	176	SER
1	B	216	ILE
1	B	297[A]	HIS
1	B	297[B]	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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)

11 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	0.36	0	17,19,21	1.85	6 (35%)
2	NAG	C	2	2	14,14,15	0.43	0	17,19,21	1.15	1 (5%)
3	NAG	D	1	1,3	14,14,15	0.46	0	17,19,21	1.22	3 (17%)
3	NAG	D	2	3	14,14,15	0.56	0	17,19,21	1.16	1 (5%)
3	BMA	D	3	3	11,11,12	0.54	0	15,15,17	0.85	0
2	NAG	E	1	1,2	14,14,15	0.54	0	17,19,21	1.54	2 (11%)
2	NAG	E	2	2	14,14,15	0.55	0	17,19,21	2.01	2 (11%)
4	NAG	F	1	1,4	14,14,15	0.52	0	17,19,21	1.20	2 (11%)
4	NAG	F	2	4	14,14,15	0.49	0	17,19,21	1.61	3 (17%)
4	BMA	F	3	4	11,11,12	0.69	0	15,15,17	0.74	1 (6%)
4	MAN	F	4	4	11,11,12	0.60	0	15,15,17	1.33	2 (13%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	C	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	C	2	2	-	1/6/23/26	0/1/1/1
3	NAG	D	1	1,3	-	2/6/23/26	0/1/1/1
3	NAG	D	2	3	-	0/6/23/26	0/1/1/1
3	BMA	D	3	3	-	0/2/19/22	0/1/1/1
2	NAG	E	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	E	2	2	-	0/6/23/26	0/1/1/1
4	NAG	F	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	F	2	4	-	2/6/23/26	0/1/1/1
4	BMA	F	3	4	-	0/2/19/22	0/1/1/1
4	MAN	F	4	4	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	E	2	NAG	C1-O5-C5	6.92	121.56	112.19
2	E	1	NAG	C1-C2-N2	-4.27	103.20	110.49
2	C	1	NAG	C2-N2-C7	4.04	128.66	122.90
2	C	1	NAG	C8-C7-N2	3.64	122.25	116.10
2	C	2	NAG	C1-O5-C5	3.61	117.08	112.19
4	F	2	NAG	C3-C4-C5	3.30	116.12	110.24
2	E	2	NAG	C3-C4-C5	3.23	116.00	110.24
4	F	2	NAG	C8-C7-N2	3.07	121.31	116.10
2	E	1	NAG	O5-C5-C6	2.82	111.63	107.20
3	D	2	NAG	O5-C5-C6	2.80	111.59	107.20
4	F	2	NAG	O5-C1-C2	-2.66	107.08	111.29
2	C	1	NAG	O5-C5-C6	2.63	111.33	107.20
4	F	4	MAN	O3-C3-C4	2.55	116.25	110.35
4	F	4	MAN	C1-O5-C5	2.43	115.49	112.19
3	D	1	NAG	O5-C1-C2	-2.37	107.55	111.29
4	F	1	NAG	C1-C2-N2	2.36	114.51	110.49
2	C	1	NAG	C4-C3-C2	2.17	114.19	111.02
4	F	3	BMA	O3-C3-C2	-2.12	105.92	109.99
3	D	1	NAG	C1-O5-C5	2.05	114.96	112.19
2	C	1	NAG	C1-C2-N2	-2.04	107.01	110.49
2	C	1	NAG	O7-C7-C8	-2.03	118.28	122.06
3	D	1	NAG	C2-N2-C7	2.03	125.79	122.90
4	F	1	NAG	O4-C4-C3	-2.03	105.67	110.35

There are no chirality outliers.

All (7) torsion outliers are listed below:

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

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 oligosaccharide.

Oligosaccharide Chain C

Bond lengths

Bond angles

Torsions

Rings

Oligosaccharide Chain E

Bond lengths

Bond angles

Torsions

Rings

Oligosaccharide Chain D

Bond lengths

Bond angles

Torsions

Rings

Oligosaccharide Chain F

Bond lengths

Bond angles

Torsions

Rings
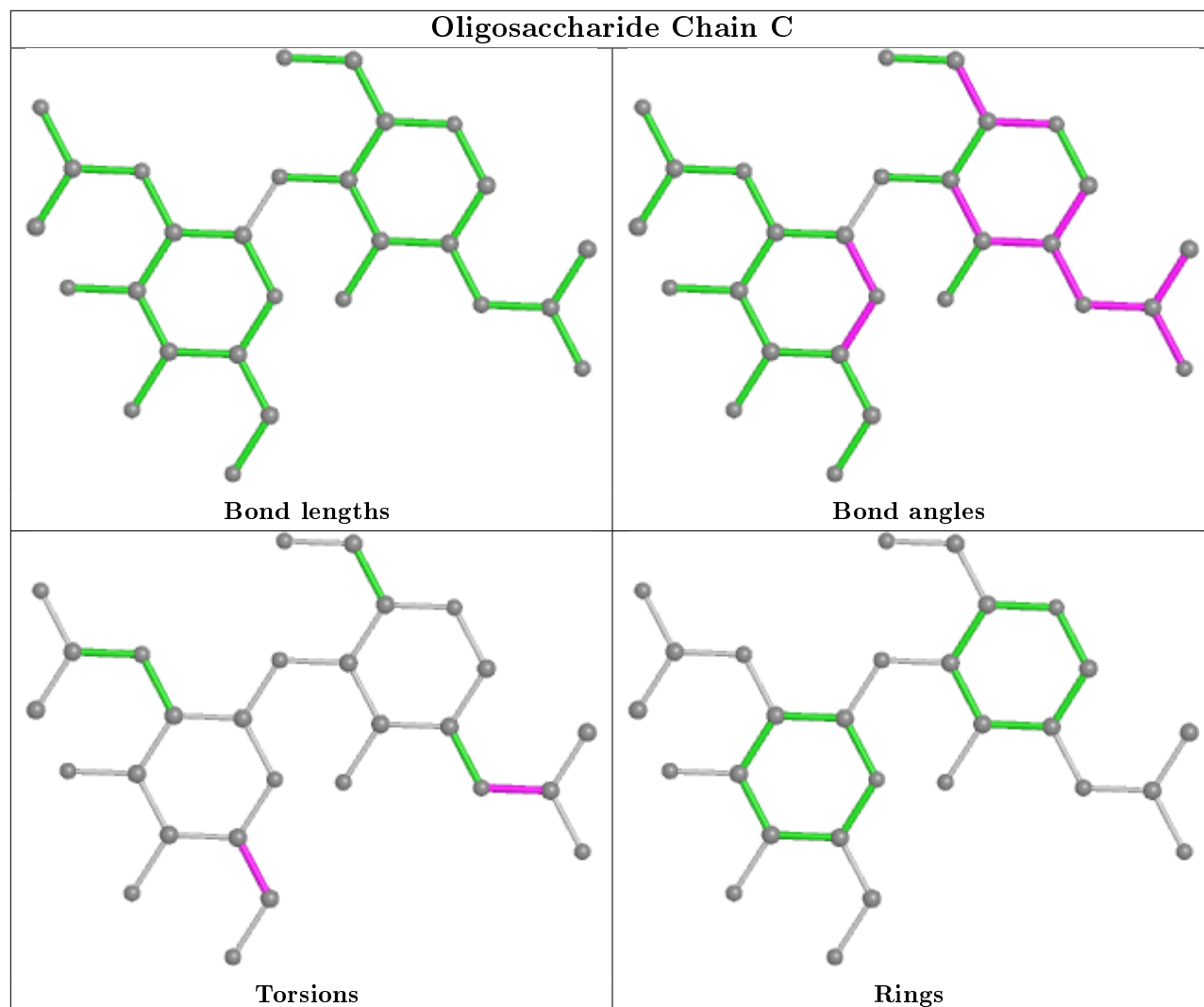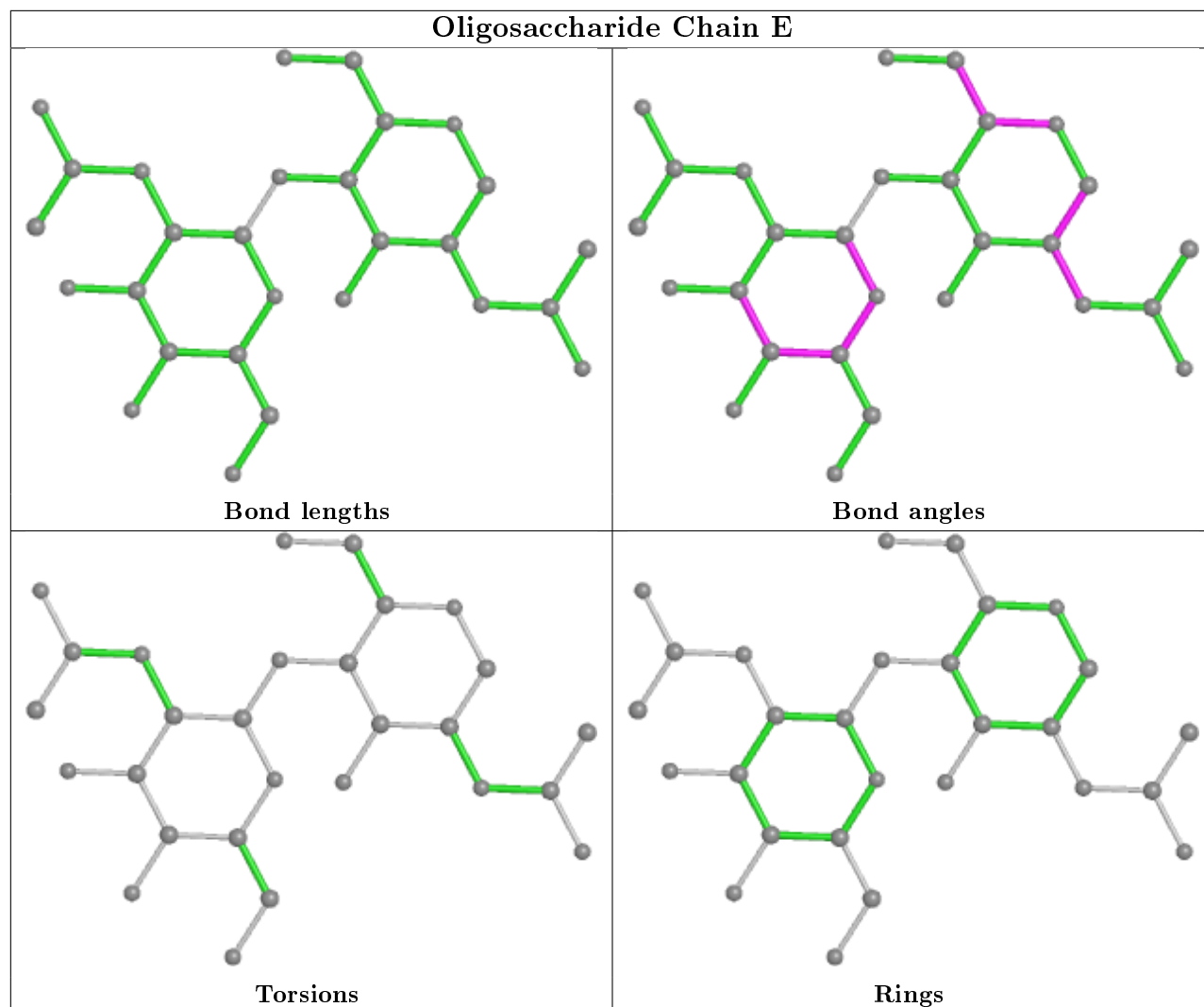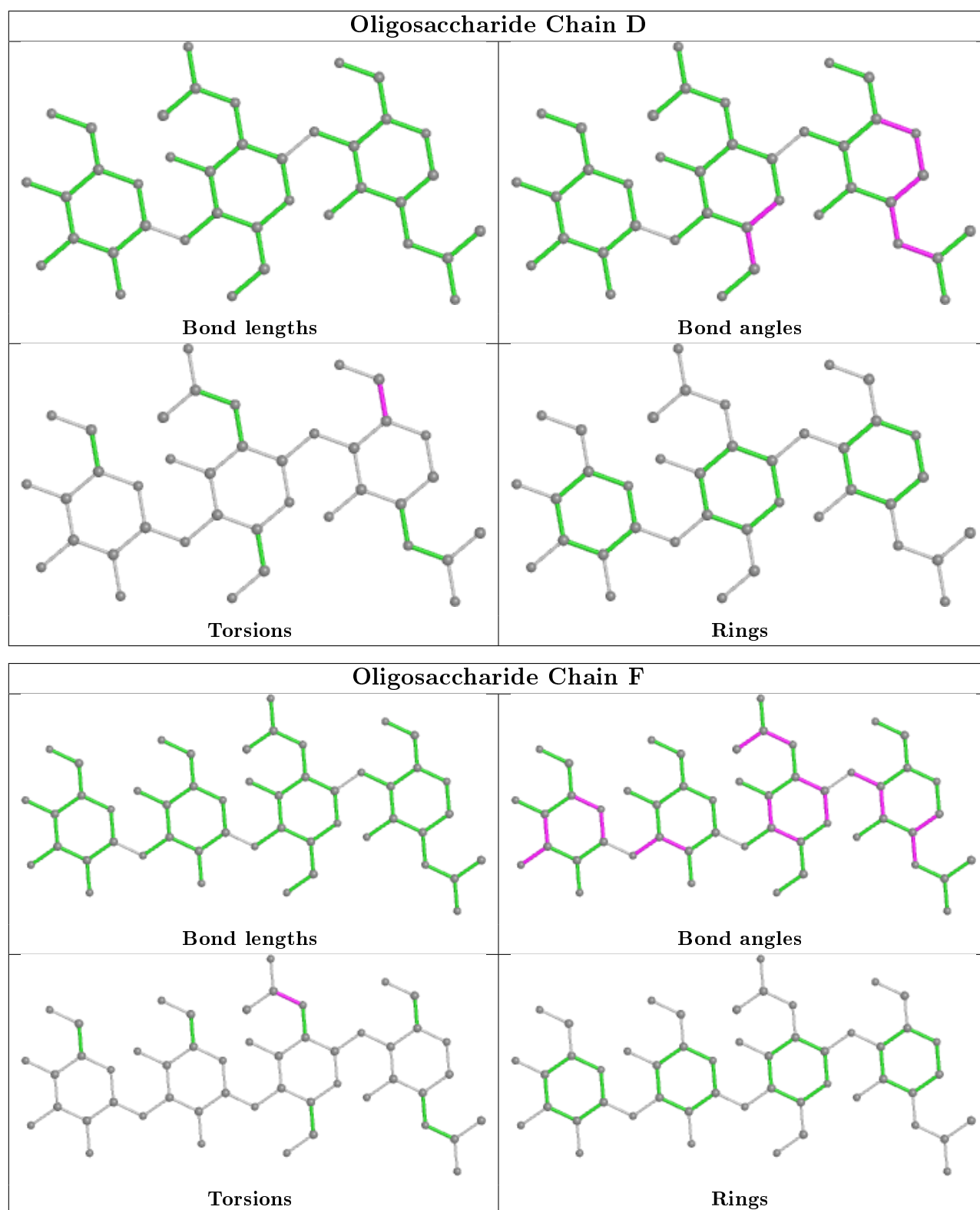
5.6 Ligand geometry (i)

Of 16 ligands modelled in this entry, 4 are monoatomic - leaving 12 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul

statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	#$|Z| > 2$	Counts	RMSZ	#$|Z| > 2$
8	EDO	B	511[B]	-	3,3,3	0.51	0	2,2,2	0.01	0
8	EDO	A	510	-	3,3,3	0.37	0	2,2,2	0.33	0
8	EDO	B	512	-	3,3,3	0.43	0	2,2,2	0.29	0
7	NAG	B	510	1	14,14,15	0.33	0	17,19,21	1.22	2 (11%)
8	EDO	A	511	-	3,3,3	0.72	0	2,2,2	0.07	0
8	EDO	B	511[A]	-	3,3,3	0.64	0	2,2,2	0.04	0
6	EEW	B	503	-	18,22,22	1.85	4 (22%)	18,31,31	1.02	2 (11%)
8	EDO	A	512	-	3,3,3	0.47	0	2,2,2	0.70	0
7	NAG	A	506	1	14,14,15	0.69	0	17,19,21	1.43	1 (5%)
8	EDO	A	513	-	3,3,3	0.57	0	2,2,2	0.14	0
8	EDO	B	513	-	3,3,3	0.63	0	2,2,2	0.22	0
6	EEW	A	503	-	18,22,22	1.75	4 (22%)	18,31,31	1.03	1 (5%)

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
8	EDO	B	511[B]	-	-	1/1/1/1	-
8	EDO	A	510	-	-	1/1/1/1	-
8	EDO	B	512	-	-	1/1/1/1	-
7	NAG	B	510	1	-	2/6/23/26	0/1/1/1
8	EDO	A	511	-	-	0/1/1/1	-
8	EDO	B	511[A]	-	-	1/1/1/1	-
6	EEW	B	503	-	-	0/15/37/37	0/1/1/1
8	EDO	A	512	-	-	0/1/1/1	-
7	NAG	A	506	1	-	0/6/23/26	0/1/1/1
8	EDO	A	513	-	-	1/1/1/1	-
8	EDO	B	513	-	-	1/1/1/1	-
6	EEW	A	503	-	-	0/15/37/37	0/1/1/1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	503	EEW	P-OP3	4.90	1.65	1.57
6	A	503	EEW	C6-C7	-4.08	1.43	1.50
6	A	503	EEW	C4-C5	3.74	1.57	1.52
6	B	503	EEW	P-C2	-3.43	1.72	1.78
6	B	503	EEW	C6-C7	-3.39	1.44	1.50
6	A	503	EEW	P-OP3	3.08	1.62	1.57
6	A	503	EEW	O7-C8	2.17	1.47	1.44
6	B	503	EEW	C3-C4	2.00	1.57	1.54

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
7	A	506	NAG	C1-O5-C5	3.86	117.42	112.19
7	B	510	NAG	C1-O5-C5	3.12	116.42	112.19
6	A	503	EEW	C4-C3-C2	2.57	112.80	109.75
6	B	503	EEW	C4-C3-C2	2.53	112.74	109.75
7	B	510	NAG	C2-N2-C7	2.20	126.03	122.90
6	B	503	EEW	C6-C7-C2	2.09	125.70	122.46

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	B	510	NAG	C4-C5-C6-O6
7	B	510	NAG	O5-C5-C6-O6
8	A	510	EDO	O1-C1-C2-O2
8	A	513	EDO	O1-C1-C2-O2
8	B	511[A]	EDO	O1-C1-C2-O2
8	B	512	EDO	O1-C1-C2-O2
8	B	511[B]	EDO	O1-C1-C2-O2
8	B	513	EDO	O1-C1-C2-O2

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	A	513	EDO	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 EEW B 503

Bond lengths

Bond angles

Torsions

Rings

Ligand EEW A 503

Bond lengths

Bond angles

Torsions

Rings
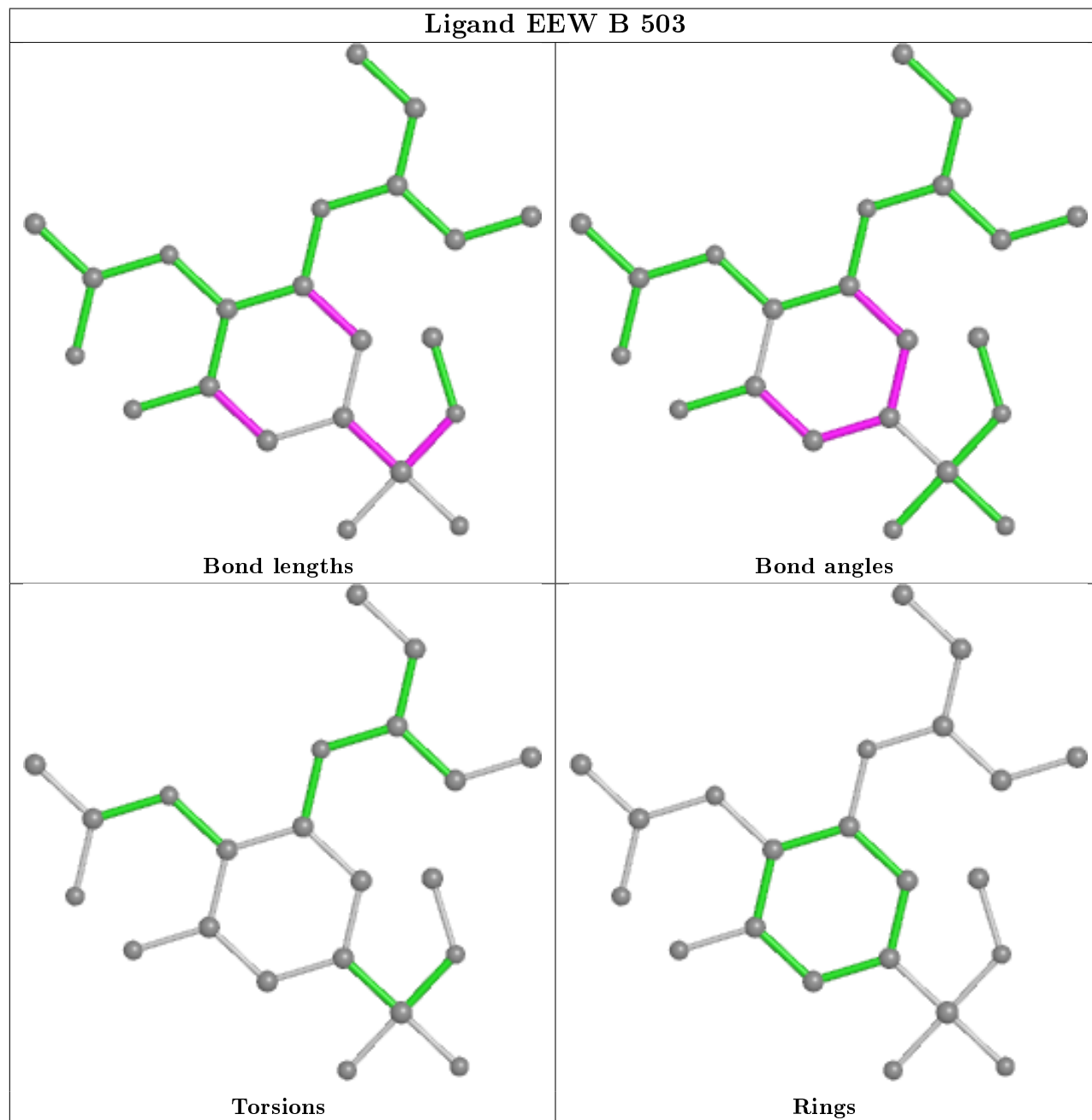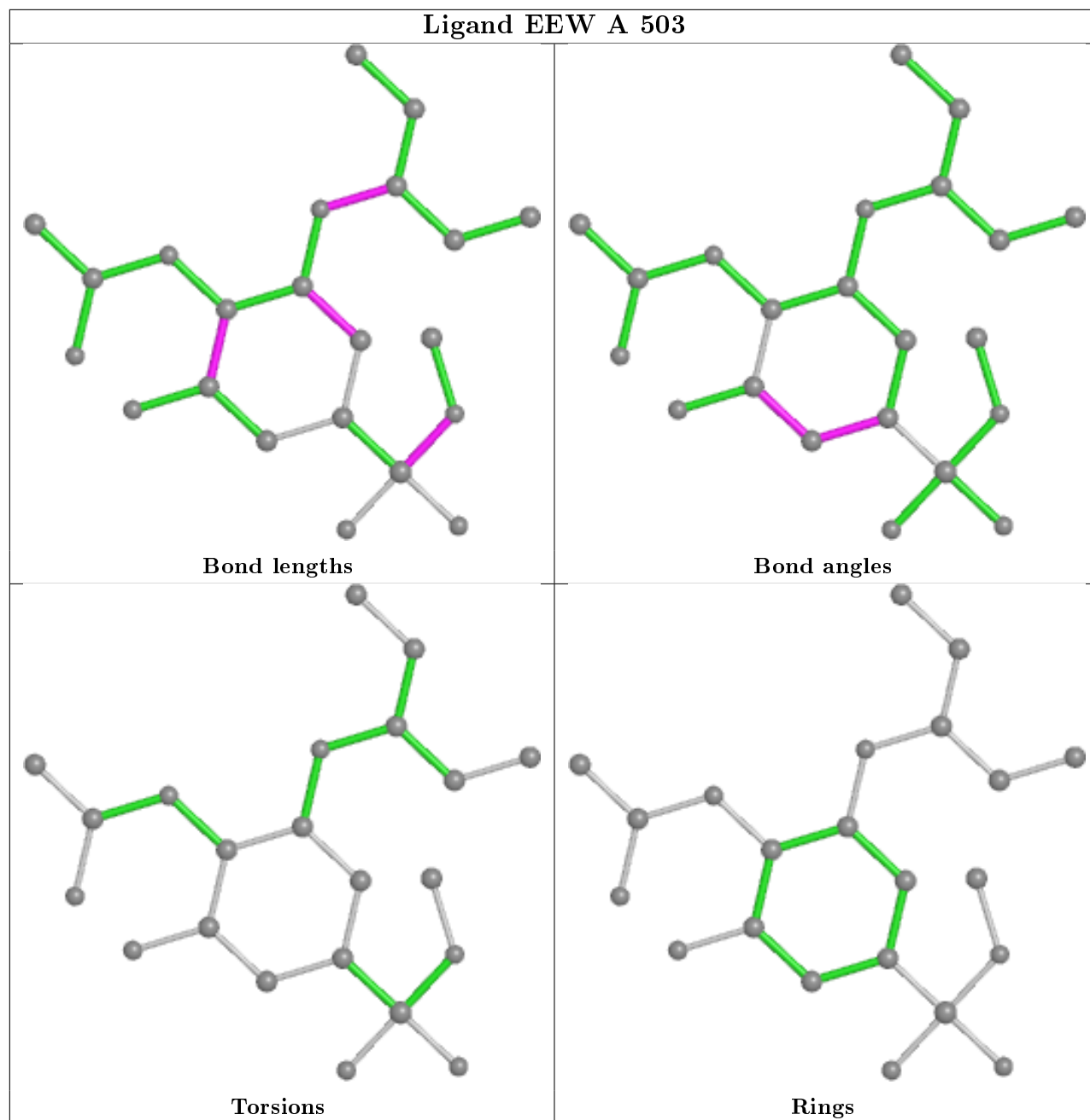
5.7 Other polymers (i)

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 (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(Å²)	Q<0.9
1	A	387/387 (100%)	-0.15	1 (0%)	94	93	12, 16, 23, 36	0
1	B	387/387 (100%)	-0.14	1 (0%)	94	93	12, 16, 23, 36	0
All	All	774/774 (100%)	-0.14	2 (0%)	94	93	12, 16, 23, 36	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	82	SER	2.3
1	A	82	SER	2.3

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(Å²)	Q<0.9
3	NAG	D	2	14/15	0.34	0.49	65,69,75,81	0
3	BMA	D	3	11/12	0.41	0.56	74,86,91,96	0
2	NAG	C	2	14/15	0.51	0.42	41,49,56,59	0
2	NAG	E	2	14/15	0.64	0.39	41,50,58,59	0
2	NAG	C	1	14/15	0.67	0.38	37,42,50,56	0
2	NAG	E	1	14/15	0.67	0.36	37,45,47,47	0
3	NAG	D	1	14/15	0.68	0.33	38,42,51,56	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($Å^2$)	Q<0.9
4	NAG	F	2	14/15	0.75	0.28	29,33,41,50	0
4	NAG	F	1	14/15	0.80	0.19	22,27,30,32	0
4	BMA	F	3	11/12	0.80	0.24	25,26,28,33	0
4	MAN	F	4	11/12	0.88	0.14	25,26,26,26	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)
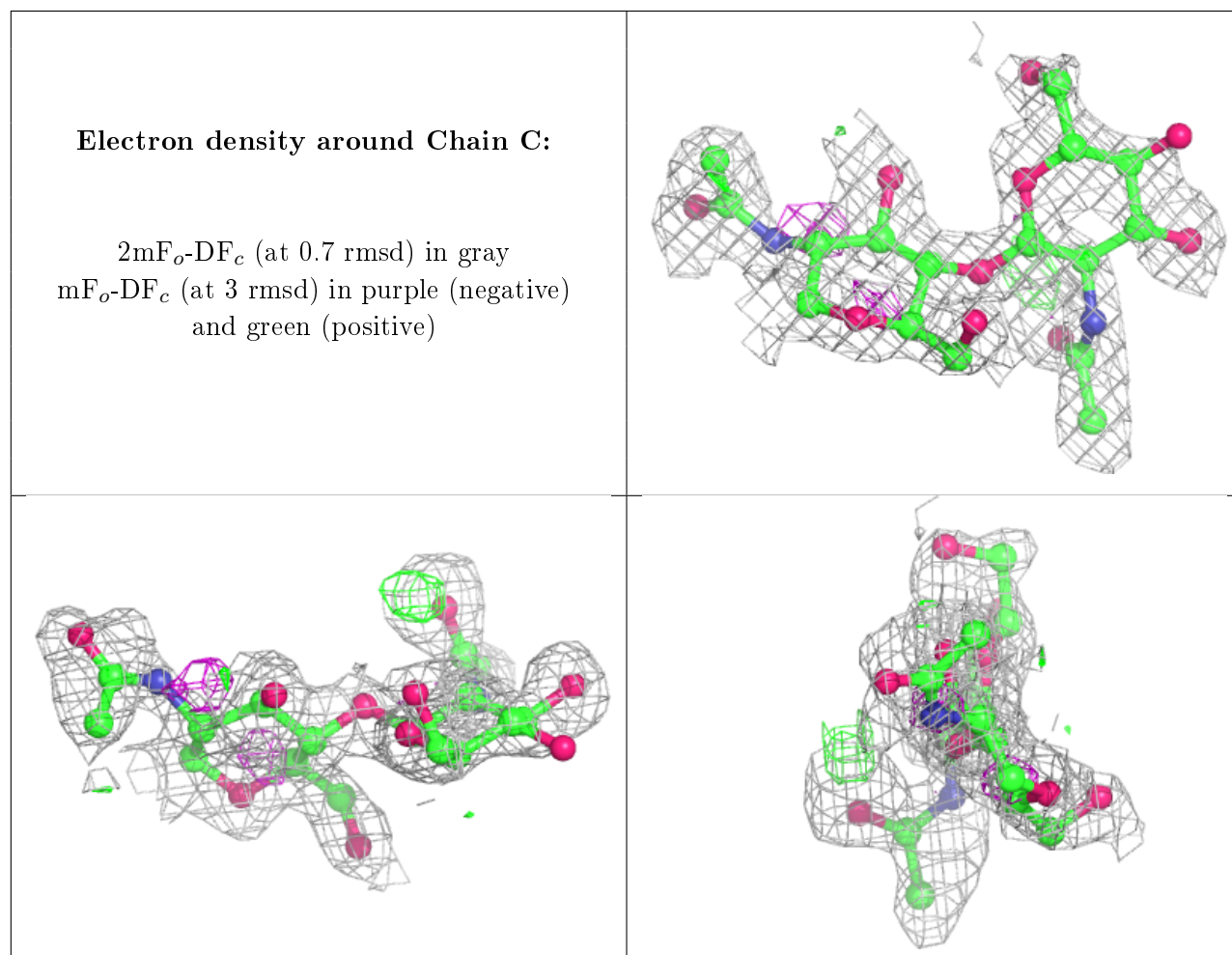
Electron density around Chain E:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)
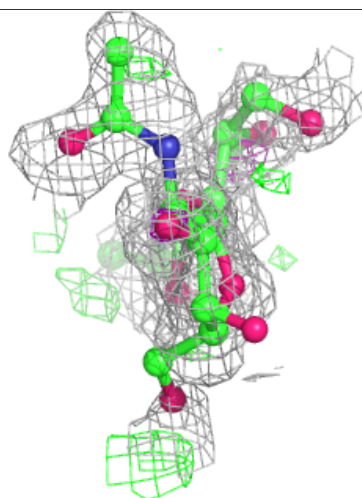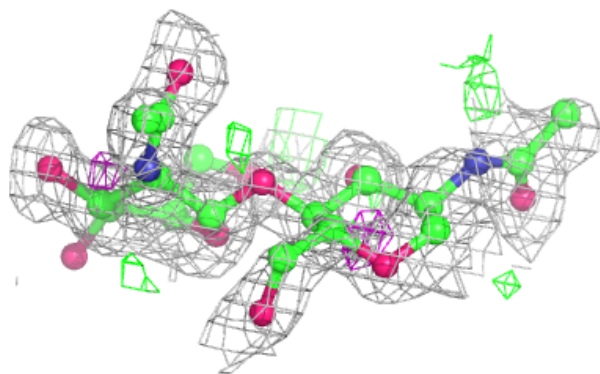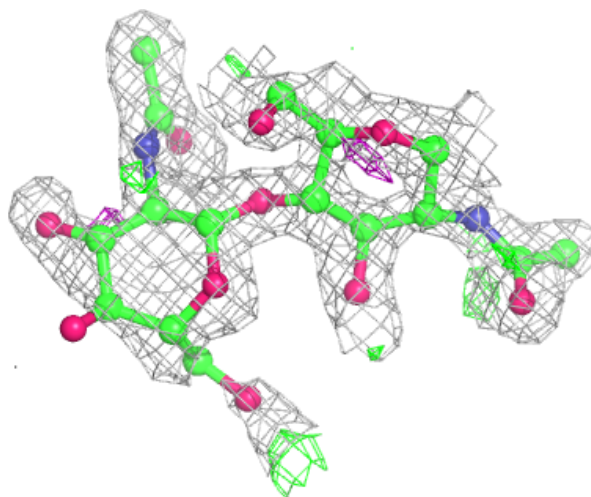
Electron density around Chain D:

2mF$_o$-DF$_c$ (at 0.7 rmsd) in gray
mF$_o$-DF$_c$ (at 3 rmsd) in purple (negative)
and green (positive)
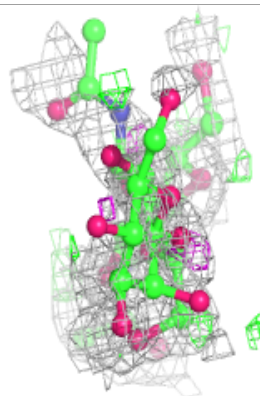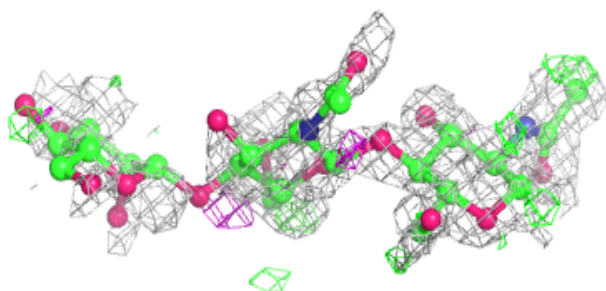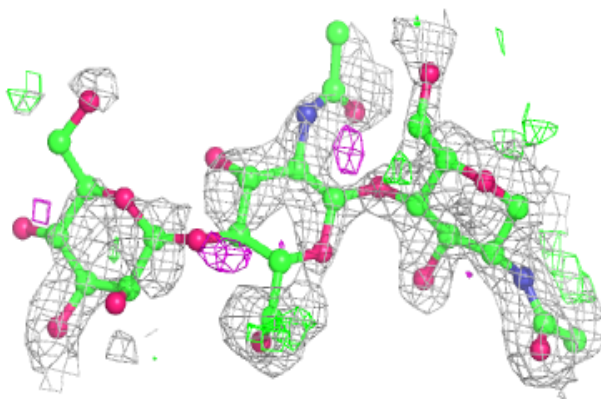
Electron density around Chain F:

2mF$_o$-DF$_c$ (at 0.7 rmsd) in gray
mF$_o$-DF$_c$ (at 3 rmsd) in purple (negative)
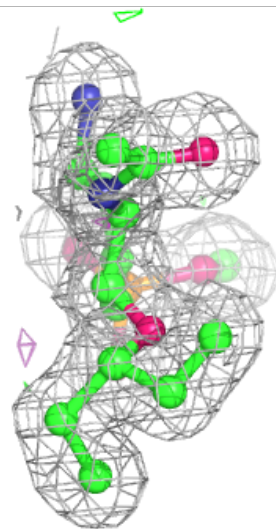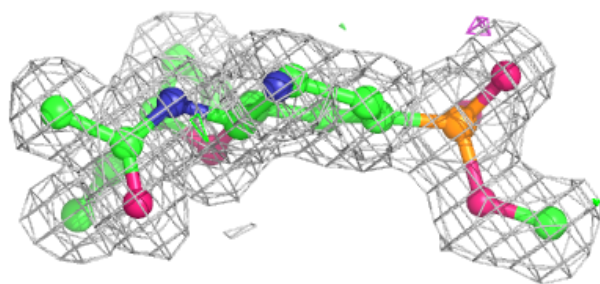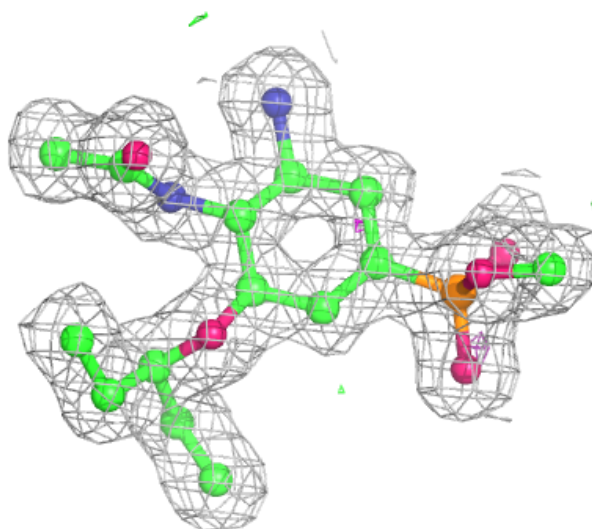and green (positive)
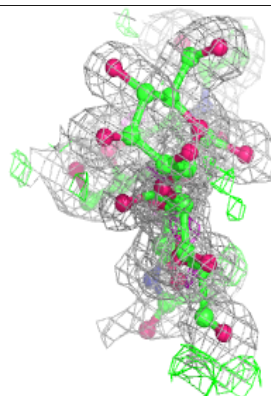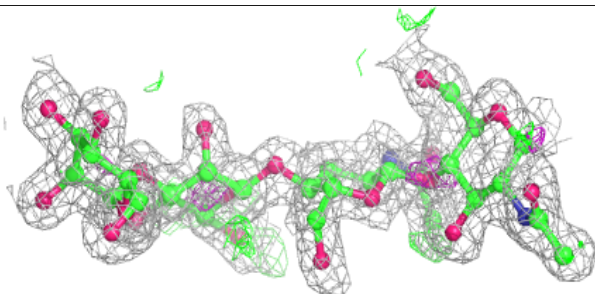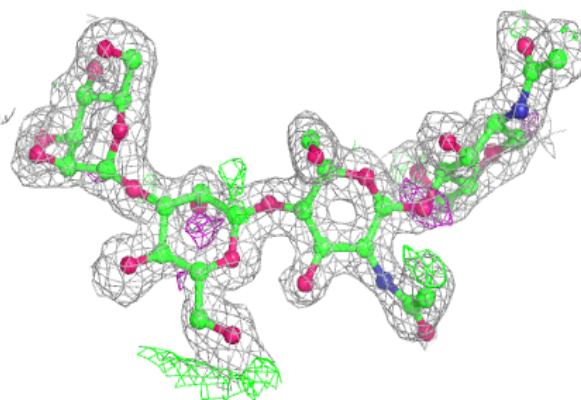
6.4 Ligands (i)

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.
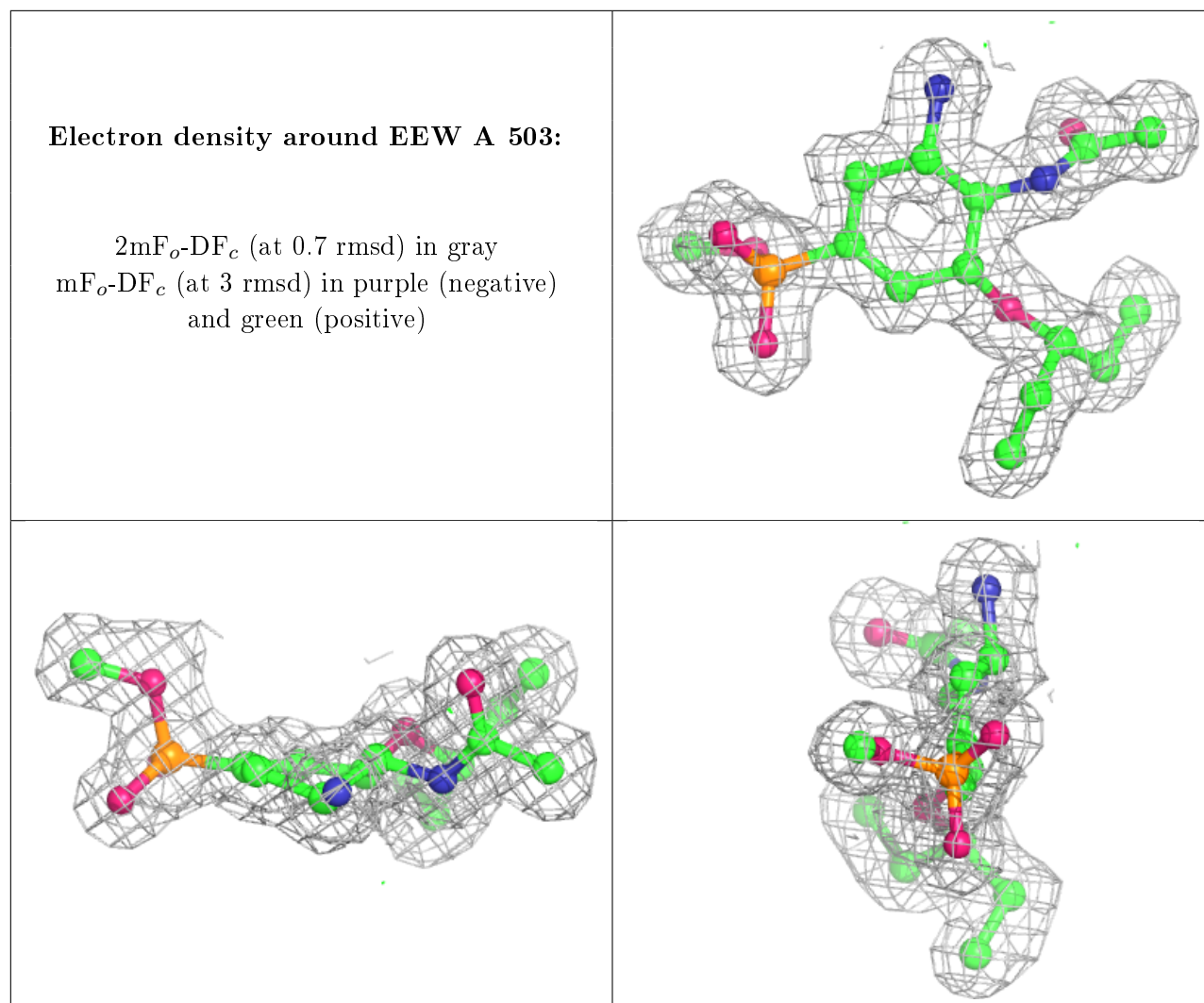
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
7	NAG	B	510	14/15	0.68	0.40	40,46,53,56	0
8	EDO	A	511	4/4	0.68	0.15	31,36,37,41	0
8	EDO	A	513	4/4	0.69	0.18	37,37,38,39	0
7	NAG	A	506	14/15	0.75	0.23	35,42,47,48	0
8	EDO	B	511[B]	4/4	0.83	0.19	17,19,19,22	4
8	EDO	B	511[A]	4/4	0.83	0.19	28,29,30,31	4
8	EDO	A	512	4/4	0.87	0.12	31,33,35,40	0
8	EDO	B	513	4/4	0.89	0.11	28,34,36,45	0
8	EDO	A	510	4/4	0.91	0.15	38,38,39,43	0
8	EDO	B	512	4/4	0.94	0.25	44,45,46,47	0
6	EEW	B	503	22/22	0.96	0.07	13,14,16,17	0
6	EEW	A	503	22/22	0.97	0.07	12,14,17,19	0
5	CA	B	502	1/1	0.98	0.05	20,20,20,20	0
5	CA	A	502	1/1	0.99	0.05	19,19,19,19	0
5	CA	B	501	1/1	0.99	0.04	15,15,15,15	0
5	CA	A	501	1/1	1.00	0.05	14,14,14,14	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 EEW B 503:

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

Electron density around EEW A 503:

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

6.5 Other polymers (i)

There are no such residues in this entry.