



wwPDB X-ray Structure Validation Summary Report ⓘ

Aug 8, 2020 – 03:30 PM BST

PDB ID : 4NDZ
Title : Structure of Maltose Binding Protein fusion to 2-O-Sulfotransferase with bound heptasaccharide and PAP
Authors : Liu, C.; Sheng, J.; Krahn, J.M.; Perera, L.; Xu, Y.; Hsieh, P.; Liu, J.; Pedersen, L.C.
Deposited on : 2013-10-28
Resolution : 3.45 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity	:	4.02b-467
Mogul	:	1.8.5 (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

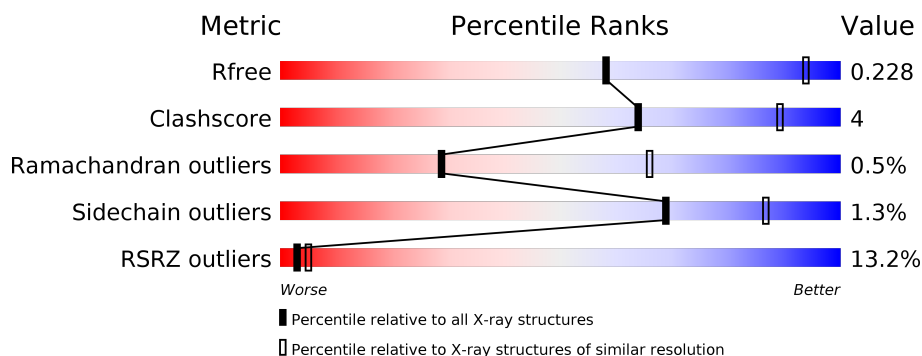
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.45 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.



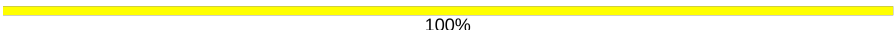



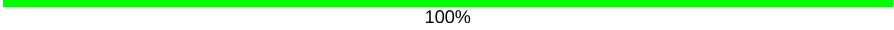
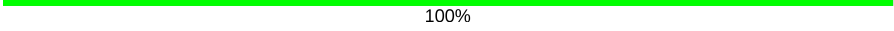

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1291 (3.52-3.40)
Clashscore	141614	1372 (3.52-3.40)
Ramachandran outliers	138981	1337 (3.52-3.40)
Sidechain outliers	138945	1338 (3.52-3.40)
RSRZ outliers	127900	1205 (3.52-3.40)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	658	<div> <div>38%</div> <div>6%</div> <div>56%</div> </div>
1	B	658	<div> <div>27%</div> <div>86%</div> <div>14%</div> <div>.</div> </div>
1	C	658	<div> <div>10%</div> <div>87%</div> <div>12%</div> </div>
1	D	658	<div> <div>6%</div> <div>86%</div> <div>12%</div> <div>..</div> </div>
1	E	658	<div> <div>5%</div> <div>87%</div> <div>13%</div> <div>.</div> </div>
1	F	658	<div> <div>23%</div> <div>86%</div> <div>13%</div> <div>.</div> </div>

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Mol	Chain	Length	Quality of chain
2	G	7	 100%
2	I	7	 86% 14%
2	L	7	 14% 86%
2	N	7	 71% 29%
2	P	7	 71% 29%
3	H	2	 100%
3	J	2	 100%
3	K	2	 50% 50%
3	M	2	 50% 50%
3	O	2	 50% 50%

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	BDP	N	5	-	-	-	X
2	BDP	N	7	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 28764 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 Maltose-binding periplasmic protein, Heparan sulfate 2-O-sulfotransferase 1 fusion.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	287	Total	C	N	O	S	0	0	1
			2352	1522	399	422	9			
1	B	654	Total	C	N	O	S	0	0	1
			5136	3314	852	955	15			
1	C	655	Total	C	N	O	S	0	0	1
			5128	3311	848	955	14			
1	D	654	Total	C	N	O	S	0	0	1
			5136	3313	853	955	15			
1	E	654	Total	C	N	O	S	0	0	1
			5110	3298	842	955	15			
1	F	654	Total	C	N	O	S	0	1	0
			5124	3307	846	956	15			

There are 42 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	359	ALA	GLU	engineered mutation	UNP P0AEX9
A	362	ALA	LYS	engineered mutation	UNP P0AEX9
A	363	ALA	ASP	engineered mutation	UNP P0AEX9
A	367	ASN	ARG	engineered mutation	UNP P0AEX9
A	368	ALA	-	linker	UNP P0AEX9
A	369	ALA	-	linker	UNP P0AEX9
A	370	ALA	-	linker	UNP P0AEX9
B	359	ALA	GLU	engineered mutation	UNP P0AEX9
B	362	ALA	LYS	engineered mutation	UNP P0AEX9
B	363	ALA	ASP	engineered mutation	UNP P0AEX9
B	367	ASN	ARG	engineered mutation	UNP P0AEX9
B	368	ALA	-	linker	UNP P0AEX9
B	369	ALA	-	linker	UNP P0AEX9
B	370	ALA	-	linker	UNP P0AEX9
C	359	ALA	GLU	engineered mutation	UNP P0AEX9
C	362	ALA	LYS	engineered mutation	UNP P0AEX9

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Chain	Residue	Modelled	Actual	Comment	Reference
C	363	ALA	ASP	engineered mutation	UNP P0AEX9
C	367	ASN	ARG	engineered mutation	UNP P0AEX9
C	368	ALA	-	linker	UNP P0AEX9
C	369	ALA	-	linker	UNP P0AEX9
C	370	ALA	-	linker	UNP P0AEX9
D	359	ALA	GLU	engineered mutation	UNP P0AEX9
D	362	ALA	LYS	engineered mutation	UNP P0AEX9
D	363	ALA	ASP	engineered mutation	UNP P0AEX9
D	367	ASN	ARG	engineered mutation	UNP P0AEX9
D	368	ALA	-	linker	UNP P0AEX9
D	369	ALA	-	linker	UNP P0AEX9
D	370	ALA	-	linker	UNP P0AEX9
E	359	ALA	GLU	engineered mutation	UNP P0AEX9
E	362	ALA	LYS	engineered mutation	UNP P0AEX9
E	363	ALA	ASP	engineered mutation	UNP P0AEX9
E	367	ASN	ARG	engineered mutation	UNP P0AEX9
E	368	ALA	-	linker	UNP P0AEX9
E	369	ALA	-	linker	UNP P0AEX9
E	370	ALA	-	linker	UNP P0AEX9
F	359	ALA	GLU	engineered mutation	UNP P0AEX9
F	362	ALA	LYS	engineered mutation	UNP P0AEX9
F	363	ALA	ASP	engineered mutation	UNP P0AEX9
F	367	ASN	ARG	engineered mutation	UNP P0AEX9
F	368	ALA	-	linker	UNP P0AEX9
F	369	ALA	-	linker	UNP P0AEX9
F	370	ALA	-	linker	UNP P0AEX9

- Molecule 2 is an oligosaccharide called beta-D-glucopyranuronic acid-(1-4)-2-acetamido-2-deoxy-alpha-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-4)-2-deoxy-2-(sulfoamino)-alpha-D-glucopyranose-(1-4)-alpha-L-idopyranuronic acid-(1-4)-2-deoxy-2-(sulfoamino)-alpha-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid.



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	G	7	Total	C	N	O	S	0	0	0
			92	44	3	43	2			
2	I	7	Total	C	N	O	S	0	0	0
			92	44	3	43	2			

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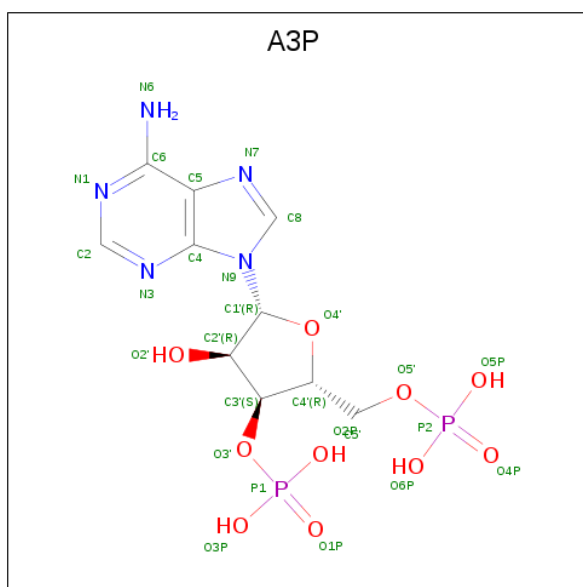
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	L	7	Total	C	N	O	S	0	0	0
			92	44	3	43	2			
2	N	7	Total	C	N	O	S	0	0	0
			92	44	3	43	2			
2	P	7	Total	C	N	O	S	0	0	0
			93	44	3	44	2			

- Molecule 3 is an oligosaccharide called alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose.



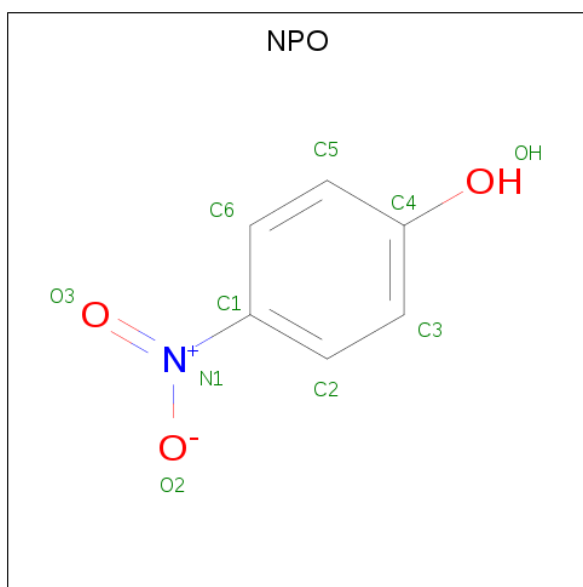
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
3	H	2	Total	C	O	0	0	0
			23	12	11			
3	J	2	Total	C	O	0	0	0
			23	12	11			
3	K	2	Total	C	O	0	0	0
			23	12	11			
3	M	2	Total	C	O	0	0	0
			23	12	11			
3	O	2	Total	C	O	0	0	0
			23	12	11			

- Molecule 4 is ADENOSINE-3'-5'-DIPHOSPHATE (three-letter code: A3P) (formula: C₁₀H₁₅N₅O₁₀P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
4	B	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
4	C	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
4	D	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
4	E	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
4	F	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

- Molecule 5 is P-NITROPHENOL (three-letter code: NPO) (formula: $C_6H_5NO_3$).

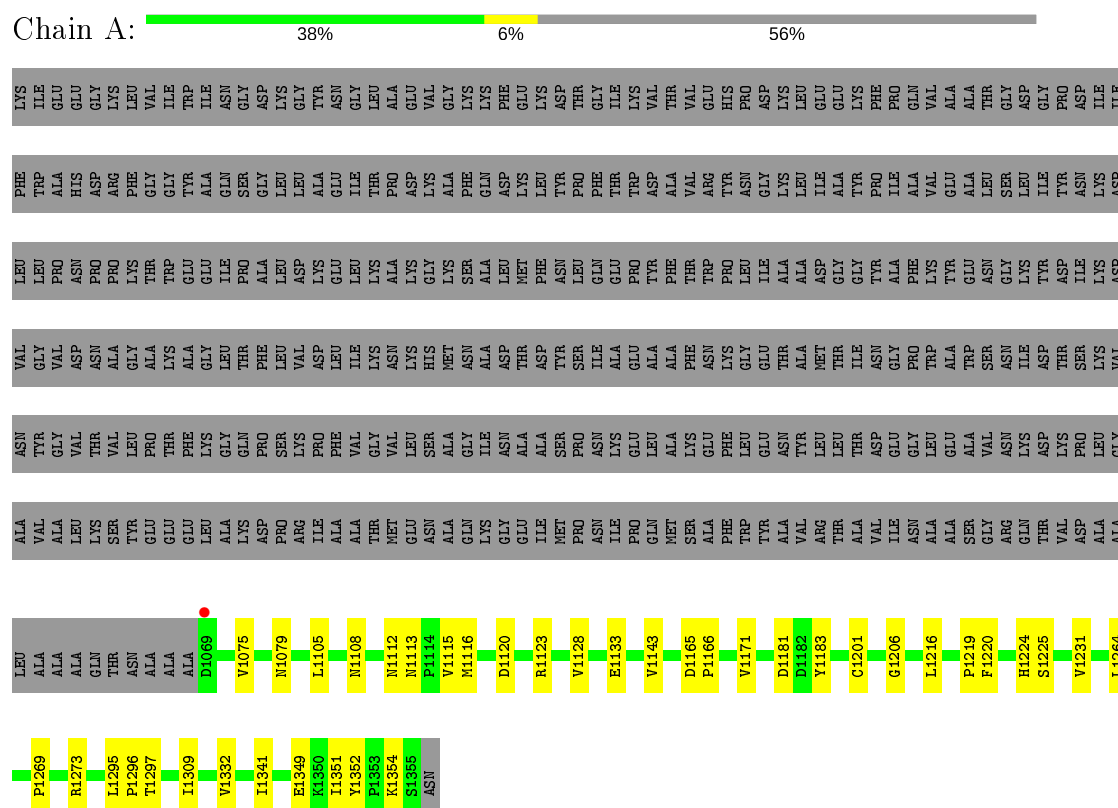


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			10	6	1	3		
5	B	1	Total	C	N	O	0	0
			10	6	1	3		
5	D	1	Total	C	N	O	0	0
			10	6	1	3		
5	E	1	Total	C	N	O	0	0
			10	6	1	3		

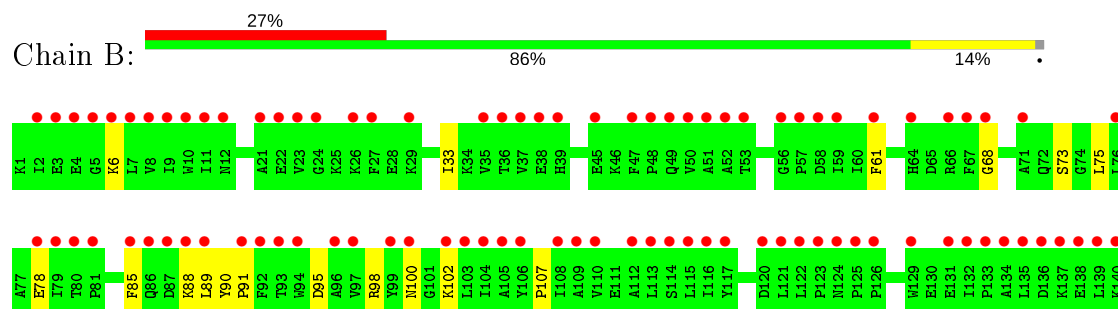
3 Residue-property plots

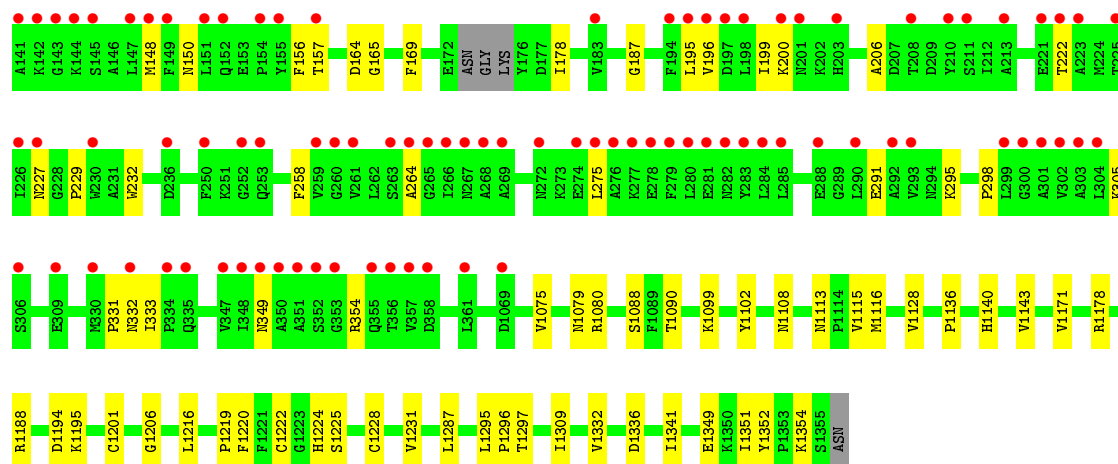
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Maltose-binding periplasmic protein, Heparan sulfate 2-O-sulfotransferase 1 fusion

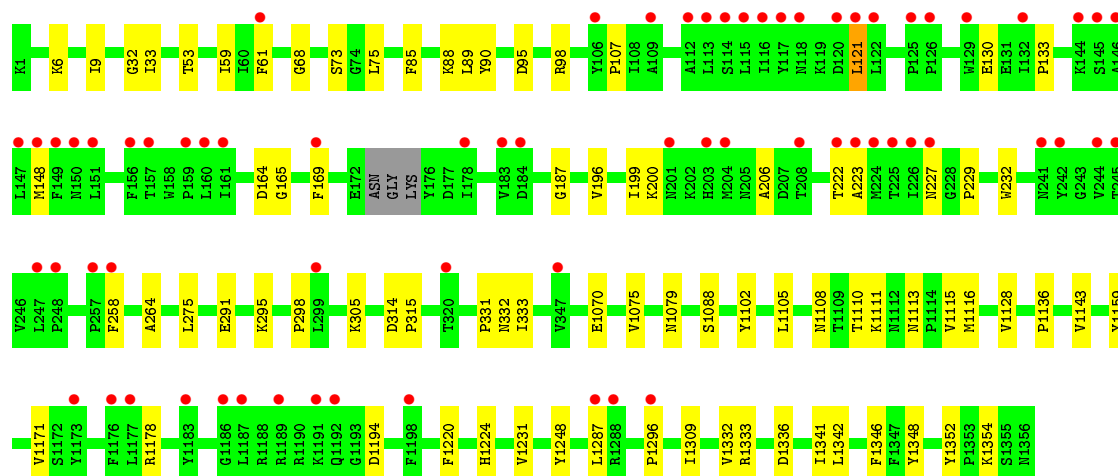
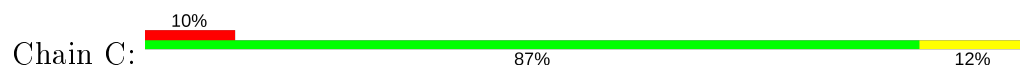


- Molecule 1: Maltose-binding periplasmic protein, Heparan sulfate 2-O-sulfotransferase 1 fusion

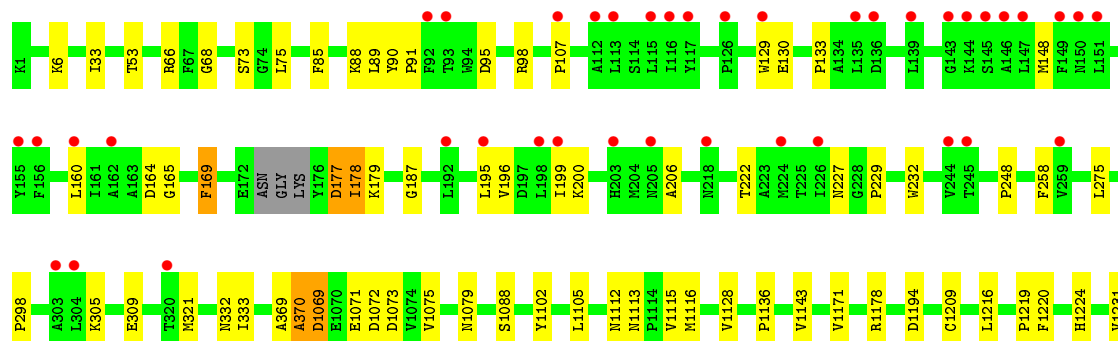
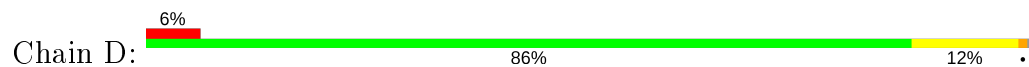




- Molecule 1: Maltose-binding periplasmic protein, Heparan sulfate 2-O-sulfotransferase 1 fusion

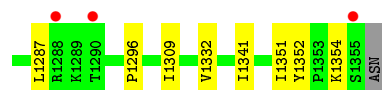
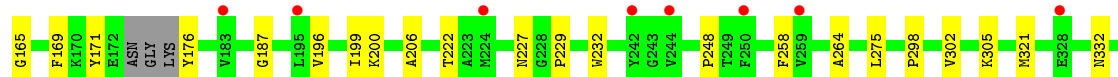
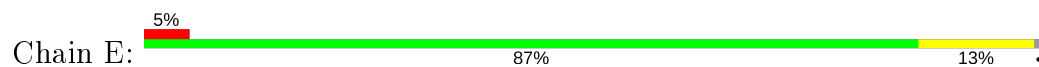


- Molecule 1: Maltose-binding periplasmic protein, Heparan sulfate 2-O-sulfotransferase 1 fusion

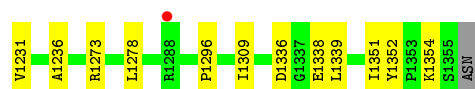
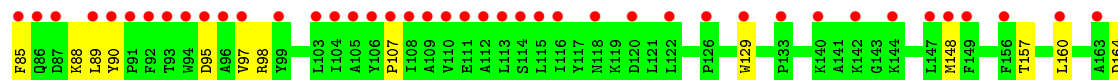
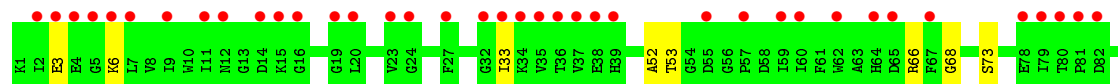
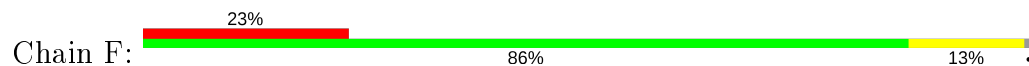




- Molecule 1: Maltose-binding periplasmic protein, Heparan sulfate 2-O-sulfotransferase 1 fusion



- Molecule 1: Maltose-binding periplasmic protein, Heparan sulfate 2-O-sulfotransferase 1 fusion





- Molecule 2: beta-D-glucopyranuronic acid-(1-4)-2-acetamido-2-deoxy-alpha-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-4)-2-deoxy-2-(sulfoamino)-alpha-D-glucopyranose-(1-4)-alpha-L-idopyranuronic acid-(1-4)-2-deoxy-2-(sulfoamino)-alpha-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid

Chain G:  100%



BDP1	GNS2	IDR3	GNS4	BDP5	IDG6	BDP7
------	------	------	------	------	------	------

- Molecule 2: beta-D-glucopyranuronic acid-(1-4)-2-acetamido-2-deoxy-alpha-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-4)-2-deoxy-2-(sulfoamino)-alpha-D-glucopyranose-(1-4)-alpha-L-idopyranuronic acid-(1-4)-2-deoxy-2-(sulfoamino)-alpha-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid

Chain I:  86%  14%

BDP1	GNS2	IDR3	GNS4	BDP5	IDG6	BDP7
------	------	------	------	------	------	------

- Molecule 2: beta-D-glucopyranuronic acid-(1-4)-2-acetamido-2-deoxy-alpha-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-4)-2-deoxy-2-(sulfoamino)-alpha-D-glucopyranose-(1-4)-alpha-L-idopyranuronic acid-(1-4)-2-deoxy-2-(sulfoamino)-alpha-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid

Chain L:  14%  86%

BDP1	GNS2	IDR3	GNS4	BDP5	IDG6	BDP7
------	------	------	------	------	------	------

- Molecule 2: beta-D-glucopyranuronic acid-(1-4)-2-acetamido-2-deoxy-alpha-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-4)-2-deoxy-2-(sulfoamino)-alpha-D-glucopyranose-(1-4)-alpha-L-idopyranuronic acid-(1-4)-2-deoxy-2-(sulfoamino)-alpha-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid

Chain N:  71%  29%

BDP1	GNS2	IDR3	GNS4	BDP5	IDG6	BDP7
------	------	------	------	------	------	------

- Molecule 2: beta-D-glucopyranuronic acid-(1-4)-2-acetamido-2-deoxy-alpha-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-4)-2-deoxy-2-(sulfoamino)-alpha-D-glucopyranose-(1-4)-alpha-L-idopyranuronic acid-(1-4)-2-deoxy-2-(sulfoamino)-alpha-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid

Chain P:  71%  29%

BDP1	GNS2	IDR3	GNS4	BDP5	IDG6	BDP7
------	------	------	------	------	------	------

- Molecule 3: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain H:  100%

GLC1
GLC2

- Molecule 3: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain J:  100%

GLC1
GLC2

- Molecule 3: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain K:  50% 50%

GLC1
GLC2

- Molecule 3: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain M:  50% 50%

GLC1
GLC2

- Molecule 3: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain O:  50% 50%

GLC1
GLC2

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	155.72Å 170.69Å 183.96Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.66 – 3.45 49.80 – 3.45	Depositor EDS
% Data completeness (in resolution range)	98.5 (49.66-3.45) 91.1 (49.80-3.45)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.08 (at 3.48Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.2_1309)	Depositor
R, R_{free}	0.176 , 0.228 0.180 , 0.228	Depositor DCC
R_{free} test set	1962 reflections (3.05%)	wwPDB-VP
Wilson B-factor (Å ²)	88.5	Xtriage
Anisotropy	0.167	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 85.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	28764	wwPDB-VP
Average B, all atoms (Å ²)	138.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.94% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 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: GNS, NPO, A3P, GLC, IDR, NDG, BDP

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	0.38	0/2416	0.47	0/3270
1	B	0.30	0/5267	0.42	0/7153
1	C	0.29	0/5259	0.43	0/7144
1	D	0.30	0/5267	0.43	0/7152
1	E	0.30	0/5241	0.44	0/7123
1	F	0.29	0/5256	0.42	0/7142
All	All	0.30	0/28706	0.43	0/38984

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	2352	0	2265	21	0
1	B	5136	0	4988	49	0
1	C	5128	0	4973	41	0
1	D	5136	0	4988	45	0
1	E	5110	0	4933	47	0
1	F	5124	0	4950	48	0
2	G	92	0	54	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	I	92	0	54	2	0
2	L	92	0	54	0	0
2	N	92	0	54	2	0
2	P	93	0	56	2	0
3	H	23	0	21	0	0
3	J	23	0	21	0	0
3	K	23	0	21	1	0
3	M	23	0	21	1	0
3	O	23	0	21	1	0
4	A	27	0	11	0	0
4	B	27	0	11	0	0
4	C	27	0	11	0	0
4	D	27	0	11	0	0
4	E	27	0	11	0	0
4	F	27	0	11	0	0
5	A	10	0	4	0	0
5	B	10	0	4	0	0
5	D	10	0	4	0	0
5	E	10	0	4	0	0
All	All	28764	0	27556	239	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 4.

The worst 5 of 239 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:D:369:ALA:O	1:D:1069:ASP:N	2.24	0.71
1:E:1131:TRP:NE1	1:F:366:THR:HG21	2.08	0.68
1:E:1131:TRP:CD1	1:F:366:THR:HG21	2.30	0.66
1:D:1071:GLU:O	1:D:1073:ASP:N	2.30	0.65
1:A:1120:ASP:OD1	1:A:1123:ARG:NH1	2.31	0.63

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	285/658 (43%)	272 (95%)	13 (5%)	0	100	100
1	B	650/658 (99%)	616 (95%)	31 (5%)	3 (0%)	29	66
1	C	651/658 (99%)	619 (95%)	28 (4%)	4 (1%)	25	62
1	D	650/658 (99%)	614 (94%)	30 (5%)	6 (1%)	17	54
1	E	650/658 (99%)	619 (95%)	28 (4%)	3 (0%)	29	66
1	F	651/658 (99%)	619 (95%)	29 (4%)	3 (0%)	29	66
All	All	3537/3948 (90%)	3359 (95%)	159 (4%)	19 (0%)	29	66

5 of 19 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	370	ALA
1	D	1072	ASP
1	E	369	ALA
1	B	1336	ASP
1	B	169	PHE

5.3.2 Protein sidechains

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	245/548 (45%)	240 (98%)	5 (2%)	55	79
1	B	524/548 (96%)	518 (99%)	6 (1%)	73	88
1	C	522/548 (95%)	518 (99%)	4 (1%)	81	92

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	524/548 (96%)	515 (98%)	9 (2%)	60	82
1	E	519/548 (95%)	514 (99%)	5 (1%)	76	89
1	F	521/548 (95%)	512 (98%)	9 (2%)	60	82
All	All	2855/3288 (87%)	2817 (99%)	38 (1%)	69	86

5 of 38 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	227	ASN
1	D	1209	CYS
1	F	1126	LYS
1	D	321	MET
1	D	1220	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 8 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	1127	ASN
1	F	1127	ASN
1	E	1127	ASN
1	C	1106	HIS
1	D	1127	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 ⓘ

45 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	BDP	G	1	2,5	9,12,13	2.94	5 (55%)	12,17,19	1.53	2 (16%)
2	GNS	G	2	2	14,15,16	2.56	7 (50%)	17,22,24	2.28	3 (17%)
2	IDR	G	3	2	9,12,13	2.62	4 (44%)	12,17,19	0.87	0
2	GNS	G	4	2	14,15,16	2.38	9 (64%)	17,22,24	2.40	5 (29%)
2	BDP	G	5	2	9,12,13	2.63	4 (44%)	12,17,19	1.44	1 (8%)
2	NDG	G	6	2	14,14,15	0.61	1 (7%)	17,19,21	0.50	0
2	BDP	G	7	2	9,12,13	2.76	5 (55%)	12,17,19	1.16	2 (16%)
3	GLC	H	1	3	12,12,12	0.53	0	17,17,17	0.63	0
3	GLC	H	2	3	11,11,12	0.63	0	15,15,17	0.71	0
2	BDP	I	1	2,5	9,12,13	2.94	5 (55%)	12,17,19	1.27	2 (16%)
2	GNS	I	2	2	14,15,16	2.51	7 (50%)	17,22,24	2.22	3 (17%)
2	IDR	I	3	2	9,12,13	2.59	4 (44%)	12,17,19	0.92	0
2	GNS	I	4	2	14,15,16	2.55	8 (57%)	17,22,24	2.42	4 (23%)
2	BDP	I	5	2	9,12,13	2.62	4 (44%)	12,17,19	1.21	1 (8%)
2	NDG	I	6	2	14,14,15	0.56	0	17,19,21	0.63	0
2	BDP	I	7	2	9,12,13	2.83	4 (44%)	12,17,19	1.07	1 (8%)
3	GLC	J	1	3	12,12,12	0.52	0	17,17,17	0.59	0
3	GLC	J	2	3	11,11,12	0.60	0	15,15,17	0.78	0
3	GLC	K	1	3	12,12,12	0.52	0	17,17,17	0.61	0
3	GLC	K	2	3	11,11,12	0.66	0	15,15,17	0.77	0
2	BDP	L	1	2,5	9,12,13	2.62	4 (44%)	12,17,19	1.38	1 (8%)
2	GNS	L	2	2	14,15,16	2.57	8 (57%)	17,22,24	2.36	6 (35%)
2	IDR	L	3	2	9,12,13	2.56	4 (44%)	12,17,19	0.56	0
2	GNS	L	4	2	14,15,16	2.50	7 (50%)	17,22,24	2.51	3 (17%)
2	BDP	L	5	2	9,12,13	2.53	3 (33%)	12,17,19	1.26	1 (8%)
2	NDG	L	6	2	14,14,15	0.50	0	17,19,21	0.66	0
2	BDP	L	7	2	9,12,13	2.69	5 (55%)	12,17,19	1.22	2 (16%)
3	GLC	M	1	3	12,12,12	0.53	0	17,17,17	0.63	0
3	GLC	M	2	3	11,11,12	0.65	0	15,15,17	0.84	0
2	BDP	N	1	2,5	9,12,13	2.87	4 (44%)	12,17,19	1.28	2 (16%)
2	GNS	N	2	2	14,15,16	2.52	7 (50%)	17,22,24	2.40	4 (23%)
2	IDR	N	3	2	9,12,13	2.42	4 (44%)	12,17,19	0.79	0
2	GNS	N	4	2	14,15,16	2.40	8 (57%)	17,22,24	2.41	3 (17%)
2	BDP	N	5	2	9,12,13	2.82	3 (33%)	12,17,19	0.90	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NDG	N	6	2	14,14,15	0.60	1 (7%)	17,19,21	0.49	0
2	BDP	N	7	2	9,12,13	2.77	5 (55%)	12,17,19	1.14	1 (8%)
3	GLC	O	1	3	12,12,12	0.51	0	17,17,17	0.64	0
3	GLC	O	2	3	11,11,12	0.65	0	15,15,17	0.64	0
2	BDP	P	1	2	10,13,13	2.01	2 (20%)	15,19,19	1.37	1 (6%)
2	GNS	P	2	2	14,15,16	2.52	8 (57%)	17,22,24	2.50	4 (23%)
2	IDR	P	3	2	9,12,13	2.44	4 (44%)	12,17,19	0.77	0
2	GNS	P	4	2	14,15,16	2.43	8 (57%)	17,22,24	2.39	4 (23%)
2	BDP	P	5	2	9,12,13	2.81	3 (33%)	12,17,19	1.12	1 (8%)
2	NDG	P	6	2	14,14,15	0.58	1 (7%)	17,19,21	0.49	0
2	BDP	P	7	2	9,12,13	2.66	5 (55%)	12,17,19	1.19	2 (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
2	BDP	G	1	2,5	-	0/0/21/24	0/1/1/1
2	GNS	G	2	2	-	0/7/24/27	0/1/1/1
2	IDR	G	3	2	-	0/0/21/24	0/1/1/1
2	GNS	G	4	2	-	1/7/24/27	0/1/1/1
2	BDP	G	5	2	-	0/0/21/24	0/1/1/1
2	NDG	G	6	2	-	0/6/23/26	0/1/1/1
2	BDP	G	7	2	-	0/0/21/24	0/1/1/1
3	GLC	H	1	3	-	0/2/22/22	0/1/1/1
3	GLC	H	2	3	-	0/2/19/22	0/1/1/1
2	BDP	I	1	2,5	-	0/0/21/24	0/1/1/1
2	GNS	I	2	2	-	0/7/24/27	0/1/1/1
2	IDR	I	3	2	-	0/0/21/24	0/1/1/1
2	GNS	I	4	2	-	0/7/24/27	0/1/1/1
2	BDP	I	5	2	-	0/0/21/24	0/1/1/1
2	NDG	I	6	2	-	4/6/23/26	0/1/1/1
2	BDP	I	7	2	-	0/0/21/24	0/1/1/1
3	GLC	J	1	3	-	0/2/22/22	0/1/1/1
3	GLC	J	2	3	-	0/2/19/22	0/1/1/1
3	GLC	K	1	3	-	0/2/22/22	0/1/1/1
3	GLC	K	2	3	-	0/2/19/22	0/1/1/1
2	BDP	L	1	2,5	-	0/0/21/24	0/1/1/1
2	GNS	L	2	2	-	0/7/24/27	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	IDR	L	3	2	-	0/0/21/24	0/1/1/1
2	GNS	L	4	2	-	1/7/24/27	0/1/1/1
2	BDP	L	5	2	-	0/0/21/24	0/1/1/1
2	NDG	L	6	2	-	2/6/23/26	0/1/1/1
2	BDP	L	7	2	-	0/0/21/24	0/1/1/1
3	GLC	M	1	3	-	0/2/22/22	0/1/1/1
3	GLC	M	2	3	-	0/2/19/22	0/1/1/1
2	BDP	N	1	2,5	-	0/0/21/24	0/1/1/1
2	GNS	N	2	2	-	0/7/24/27	0/1/1/1
2	IDR	N	3	2	-	0/0/21/24	0/1/1/1
2	GNS	N	4	2	-	0/7/24/27	0/1/1/1
2	BDP	N	5	2	-	0/0/21/24	0/1/1/1
2	NDG	N	6	2	-	0/6/23/26	0/1/1/1
2	BDP	N	7	2	-	0/0/21/24	0/1/1/1
3	GLC	O	1	3	-	0/2/22/22	0/1/1/1
3	GLC	O	2	3	-	0/2/19/22	0/1/1/1
2	BDP	P	1	2	-	0/0/24/24	0/1/1/1
2	GNS	P	2	2	-	0/7/24/27	0/1/1/1
2	IDR	P	3	2	-	0/0/21/24	0/1/1/1
2	GNS	P	4	2	-	0/7/24/27	0/1/1/1
2	BDP	P	5	2	-	0/0/21/24	0/1/1/1
2	NDG	P	6	2	-	0/6/23/26	0/1/1/1
2	BDP	P	7	2	-	0/0/21/24	0/1/1/1

The worst 5 of 161 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	G	1	BDP	O5-C5	-6.33	1.36	1.43
2	I	1	BDP	O5-C5	-6.03	1.37	1.43
2	N	1	BDP	O5-C5	-6.02	1.37	1.43
2	P	5	BDP	O5-C5	-5.98	1.37	1.43
2	I	7	BDP	O5-C5	-5.94	1.37	1.43

The worst 5 of 59 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	P	4	GNS	O3S-S1-O2S	-7.14	103.30	120.16
2	L	4	GNS	O3S-S1-O2S	-7.11	103.37	120.16
2	N	4	GNS	O3S-S1-O2S	-7.07	103.45	120.16
2	L	2	GNS	O3S-S1-O2S	-6.79	104.11	120.16
2	I	4	GNS	O3S-S1-O2S	-6.77	104.15	120.16

There are no chirality outliers.

5 of 8 torsion outliers are listed below:

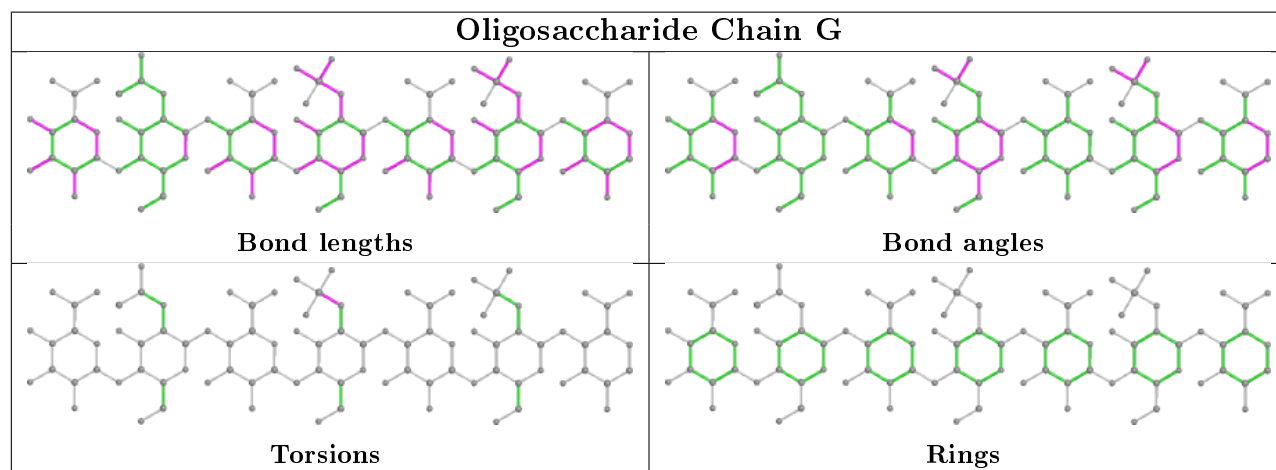
Mol	Chain	Res	Type	Atoms
2	I	6	NDG	C1-C2-N2-C7
2	I	6	NDG	C4-C5-C6-O6
2	L	6	NDG	C4-C5-C6-O6
2	I	6	NDG	O5-C5-C6-O6
2	L	6	NDG	O5-C5-C6-O6

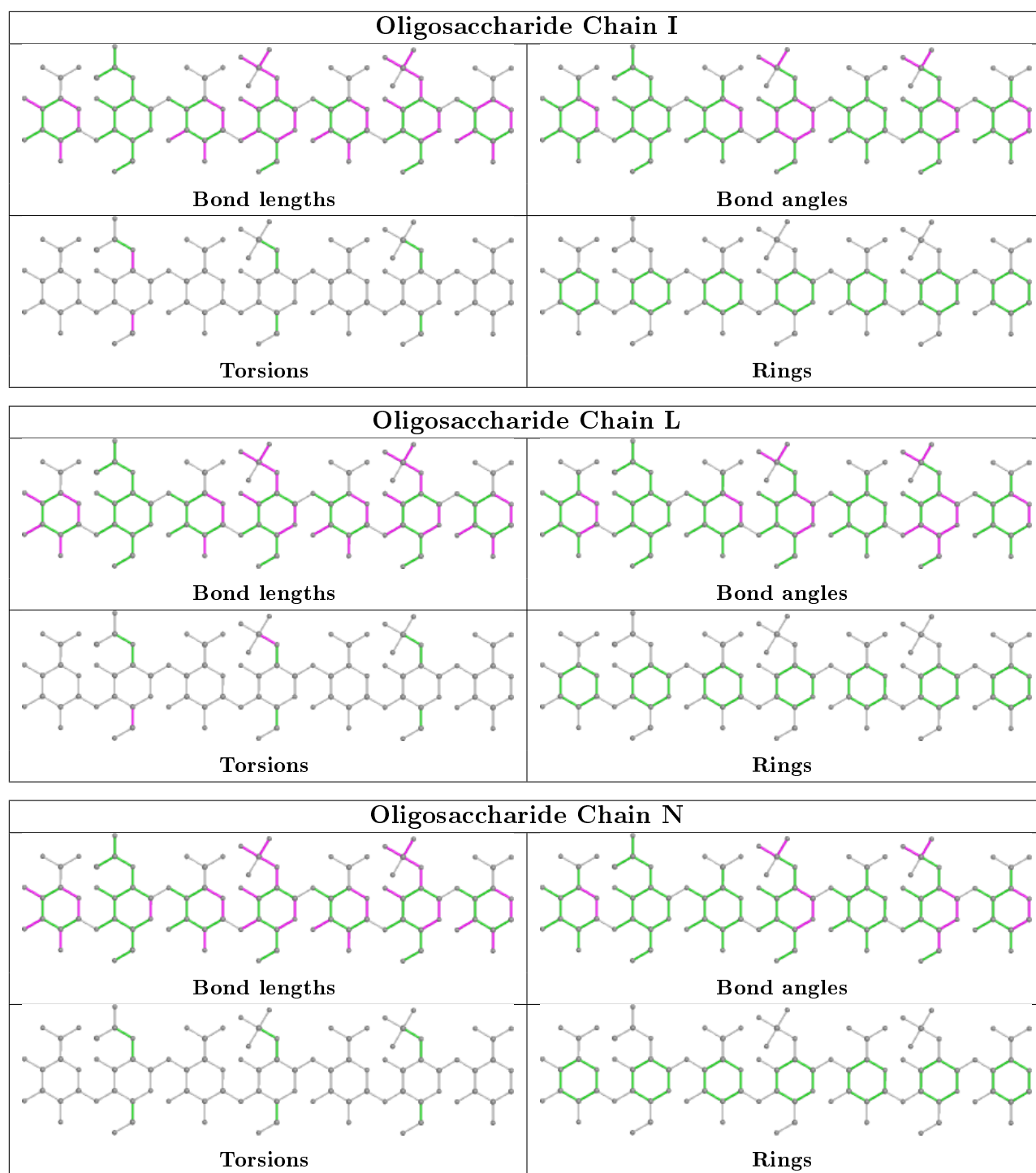
There are no ring outliers.

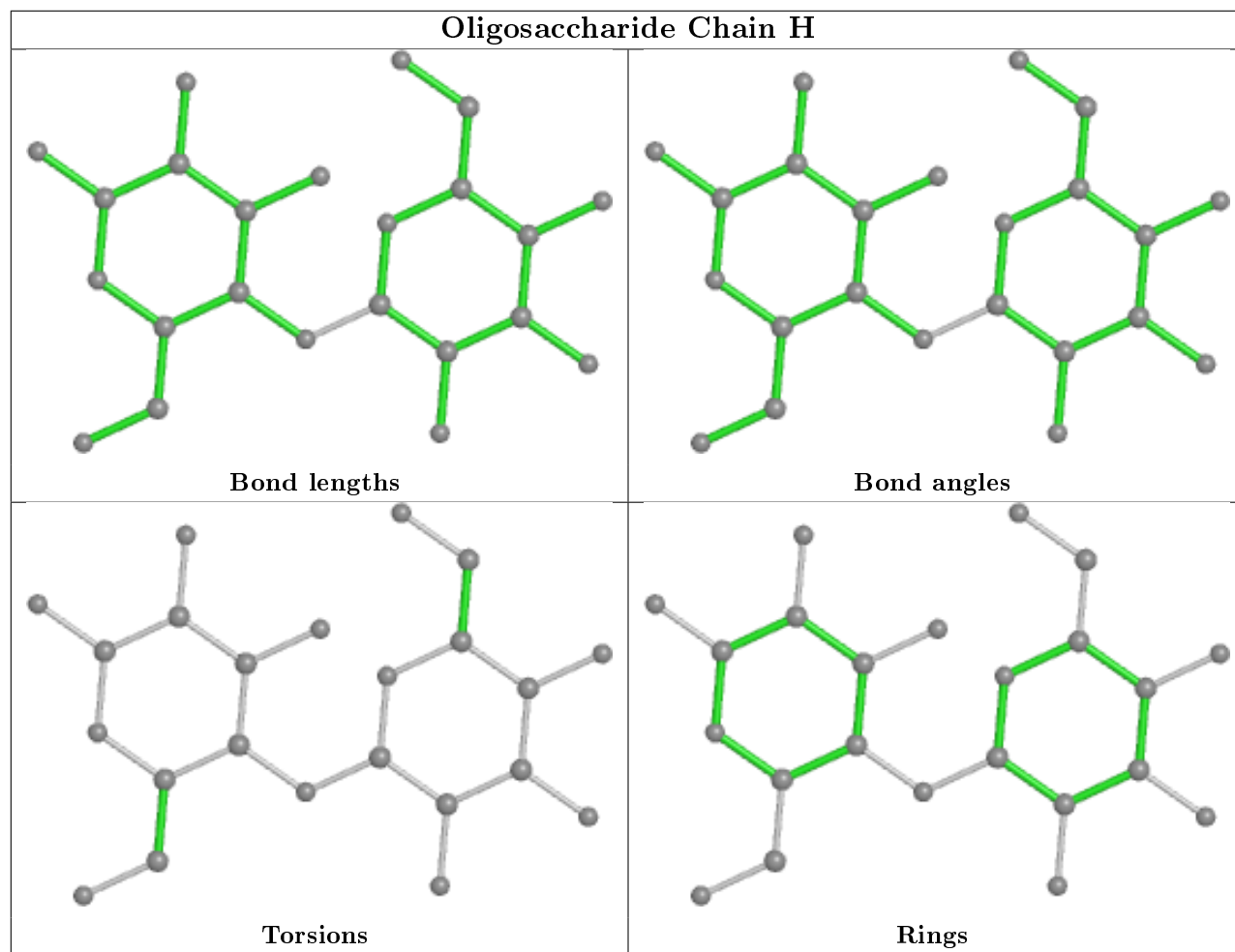
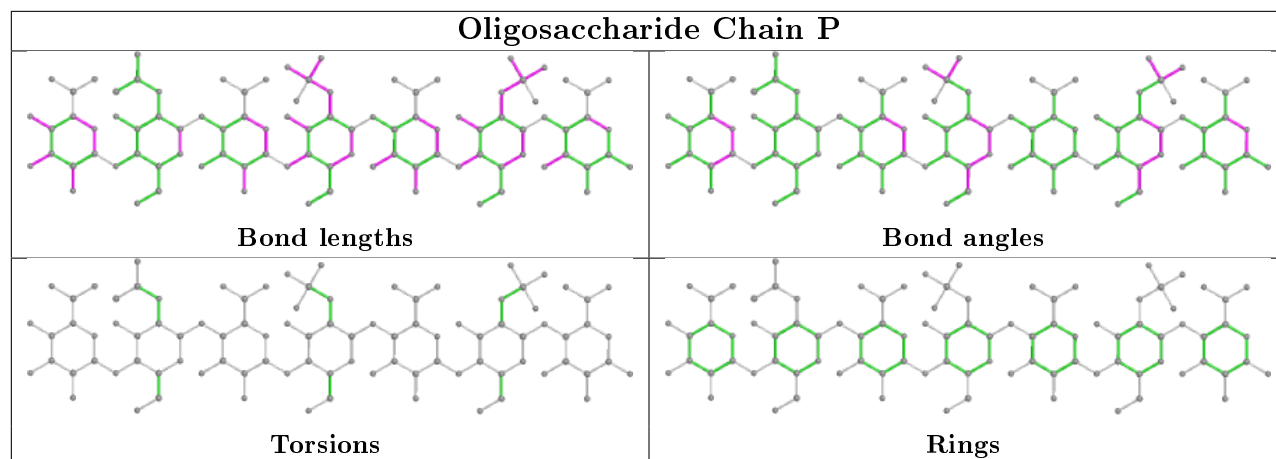
9 monomers are involved in 9 short contacts:

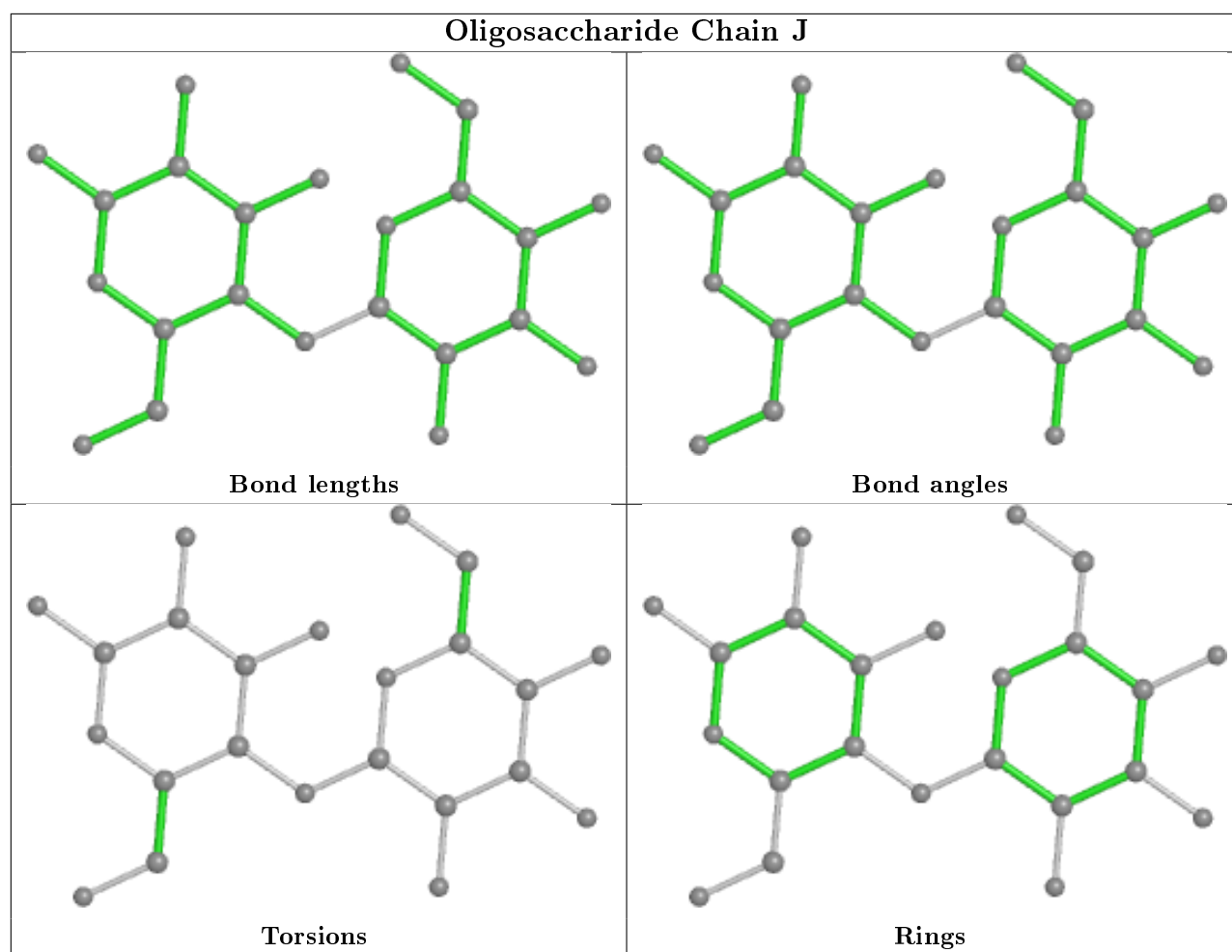
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	M	2	GLC	1	0
3	K	2	GLC	1	0
3	O	2	GLC	1	0
2	N	1	BDP	1	0
2	N	4	GNS	1	0
2	P	4	GNS	1	0
2	I	6	NDG	1	0
2	P	1	BDP	1	0
2	I	4	GNS	1	0

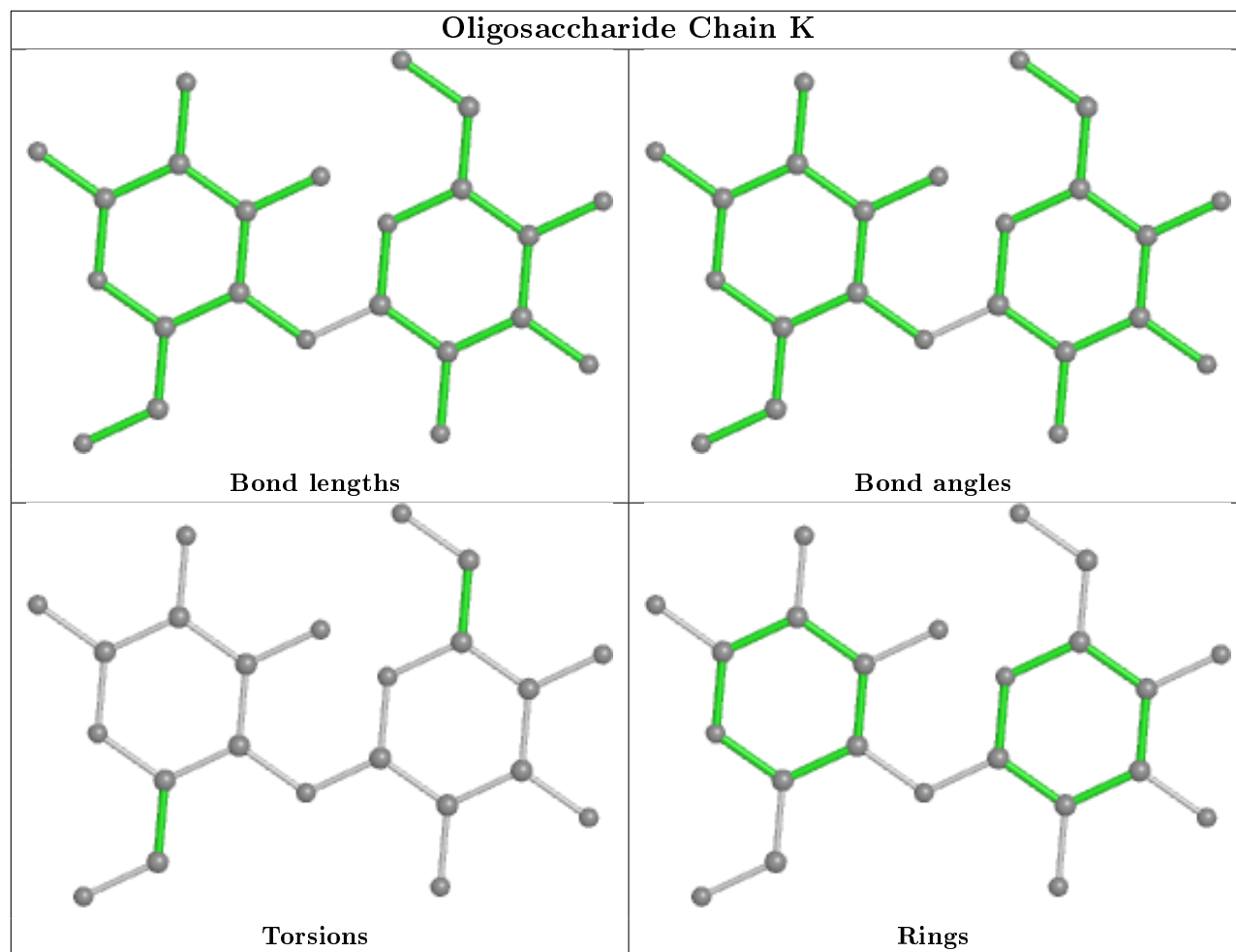
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

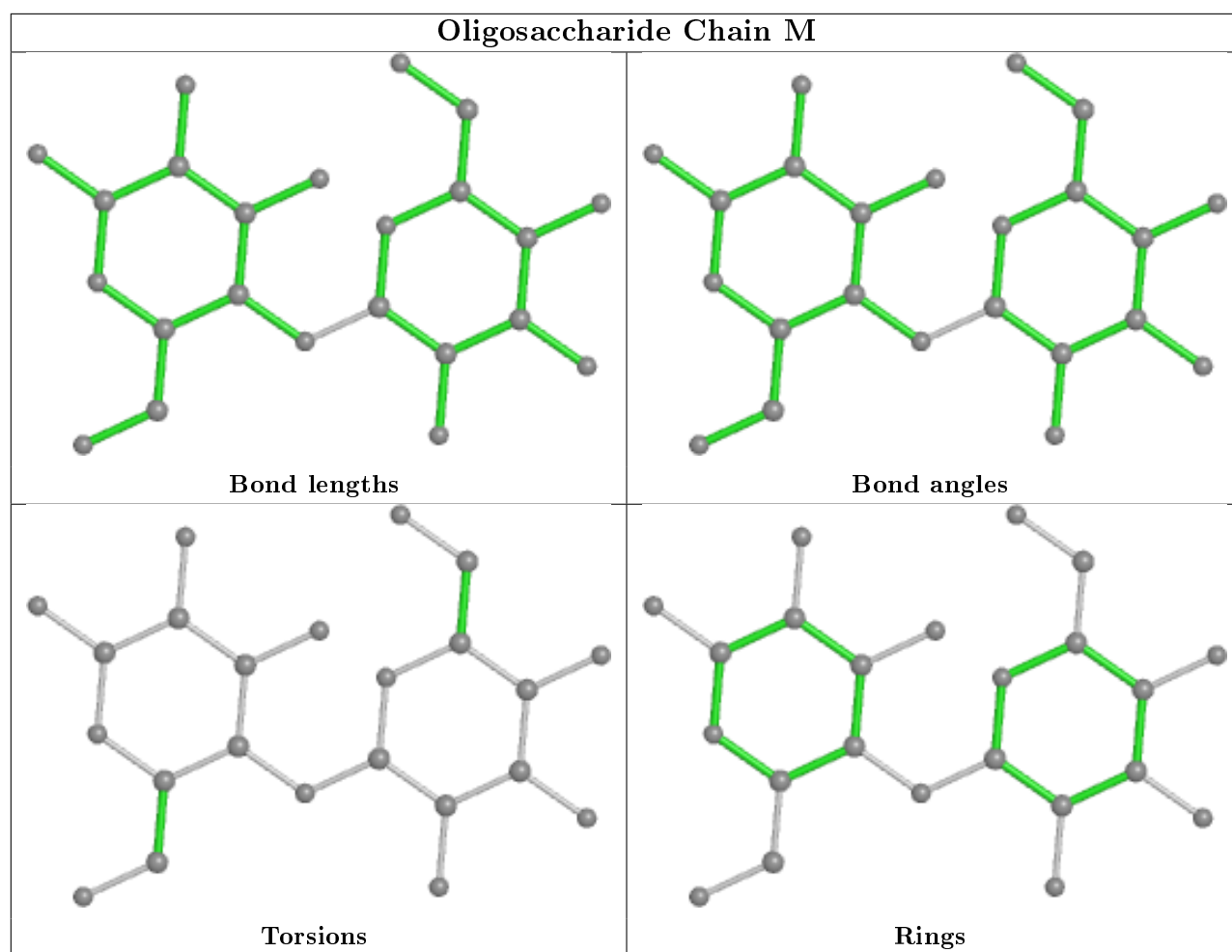


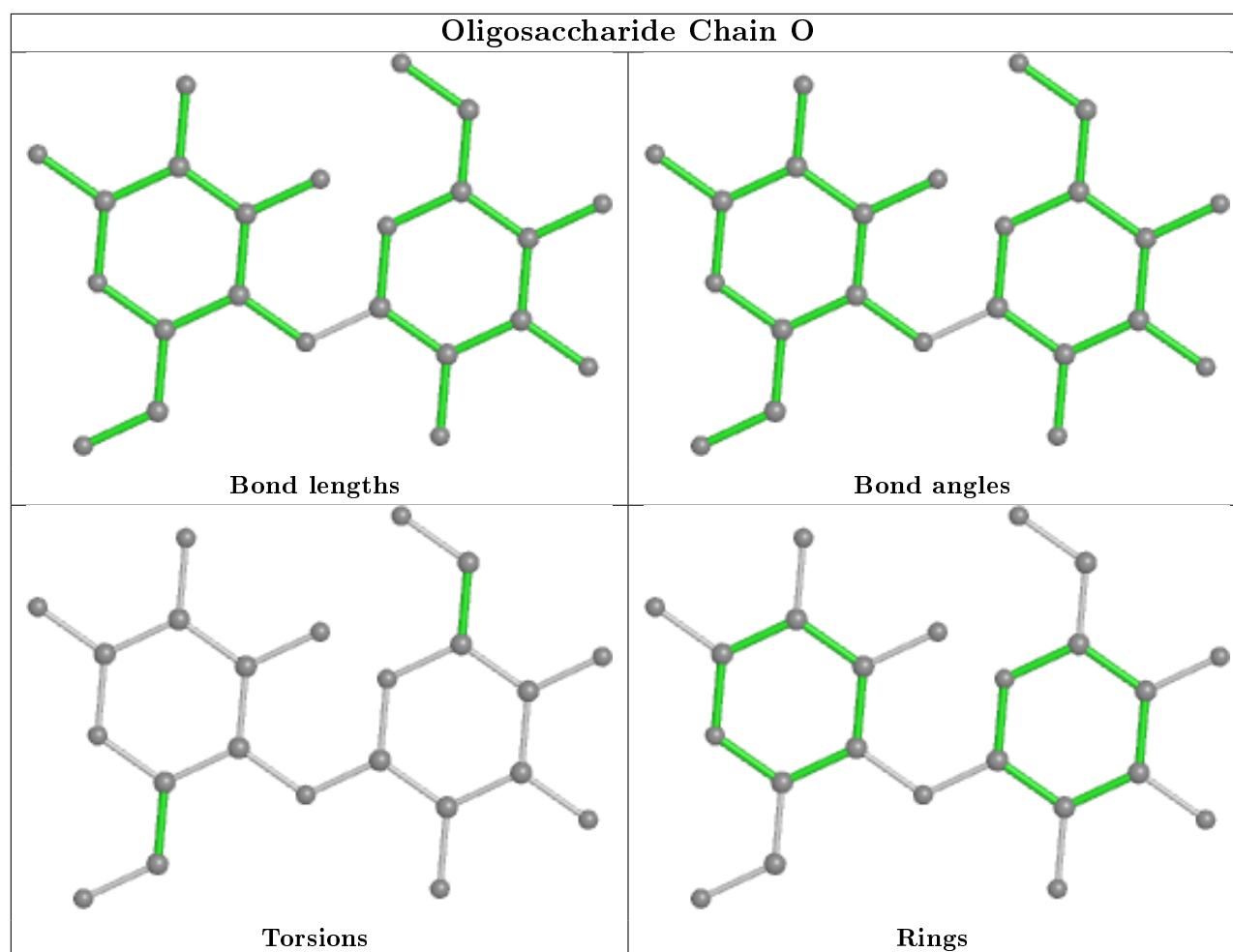












5.6 Ligand geometry [i](#)

10 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	A3P	D	2003	-	26,29,29	2.77	11 (42%)	31,45,45	1.16	1 (3%)
4	A3P	B	2003	-	26,29,29	2.74	11 (42%)	31,45,45	1.18	1 (3%)
4	A3P	F	2003	-	26,29,29	2.84	11 (42%)	31,45,45	1.16	2 (6%)
4	A3P	C	2003	-	26,29,29	2.82	10 (38%)	31,45,45	1.13	2 (6%)
4	A3P	A	2003	-	26,29,29	2.78	12 (46%)	31,45,45	1.14	2 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	A3P	E	2003	-	26,29,29	2.80	10 (38%)	31,45,45	1.18	1 (3%)
5	NPO	E	2011	2	9,10,10	0.40	0	11,13,13	0.58	0
5	NPO	D	2011	2	9,10,10	0.40	0	11,13,13	0.60	0
5	NPO	A	2011	2	9,10,10	0.48	0	11,13,13	0.95	1 (9%)
5	NPO	B	2011	2	9,10,10	0.45	0	11,13,13	0.79	1 (9%)

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	A3P	D	2003	-	-	0/11/31/31	0/3/3/3
4	A3P	B	2003	-	-	0/11/31/31	0/3/3/3
4	A3P	F	2003	-	-	1/11/31/31	0/3/3/3
4	A3P	C	2003	-	-	0/11/31/31	0/3/3/3
4	A3P	A	2003	-	-	0/11/31/31	0/3/3/3
4	A3P	E	2003	-	-	0/11/31/31	0/3/3/3
5	NPO	E	2011	2	-	0/2/4/4	0/1/1/1
5	NPO	D	2011	2	-	0/2/4/4	0/1/1/1
5	NPO	A	2011	2	-	0/2/4/4	0/1/1/1
5	NPO	B	2011	2	-	0/2/4/4	0/1/1/1

The worst 5 of 65 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	F	2003	A3P	C2-N3	7.26	1.43	1.32
4	E	2003	A3P	C2-N3	7.16	1.43	1.32
4	D	2003	A3P	C2-N3	7.09	1.43	1.32
4	C	2003	A3P	C2-N3	7.01	1.43	1.32
4	A	2003	A3P	C2-N3	6.95	1.43	1.32

The worst 5 of 11 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	E	2003	A3P	N3-C2-N1	-4.81	121.16	128.68
4	B	2003	A3P	N3-C2-N1	-4.69	121.34	128.68
4	D	2003	A3P	N3-C2-N1	-4.64	121.42	128.68
4	C	2003	A3P	N3-C2-N1	-4.58	121.52	128.68
4	F	2003	A3P	N3-C2-N1	-4.49	121.66	128.68

There are no chirality outliers.

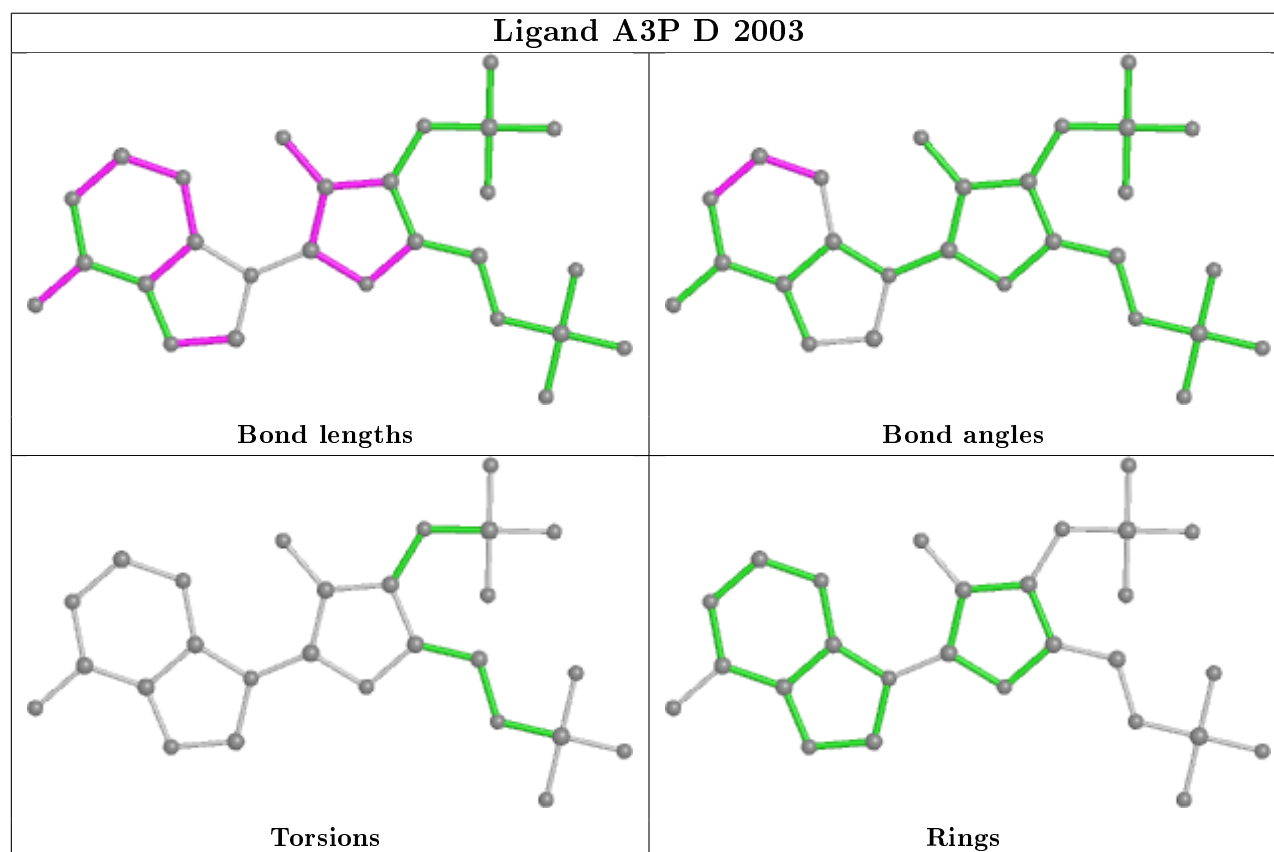
All (1) torsion outliers are listed below:

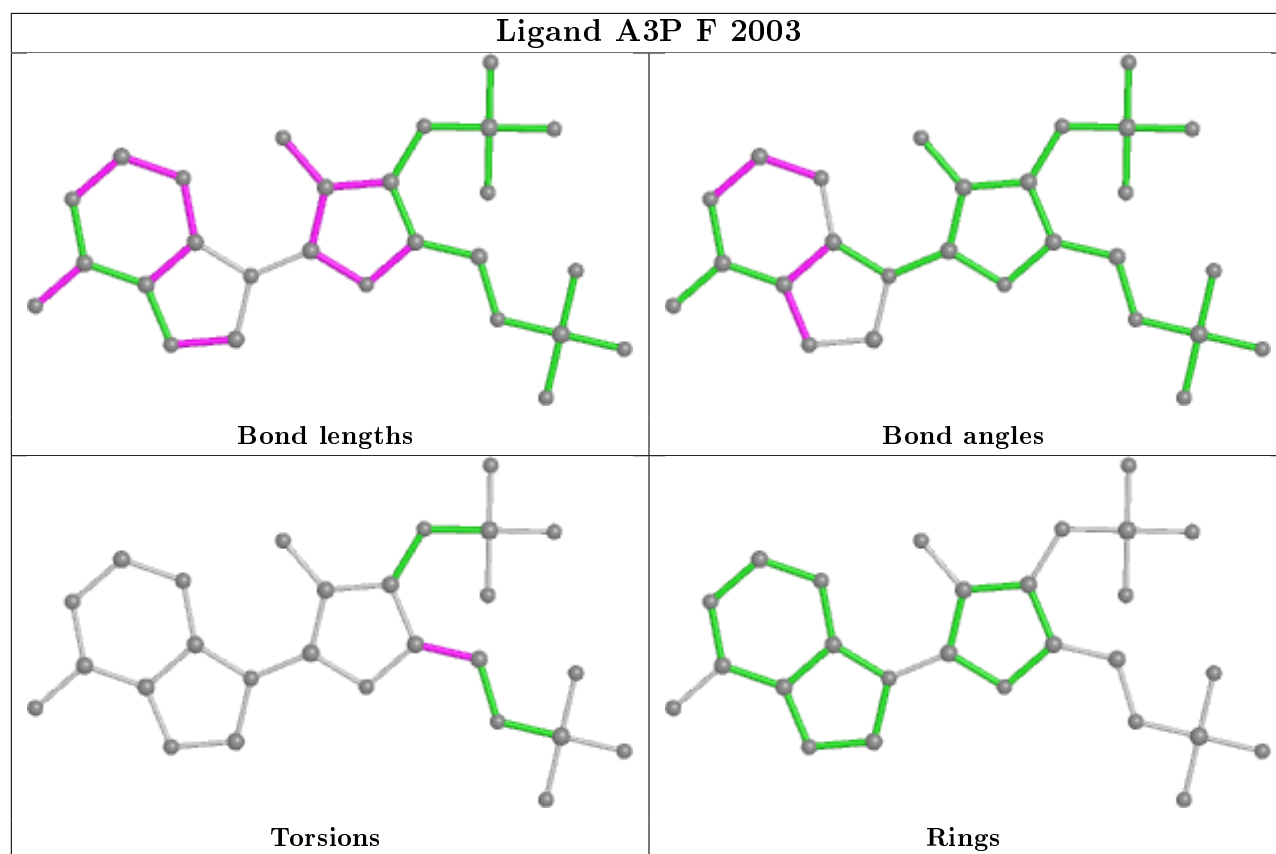
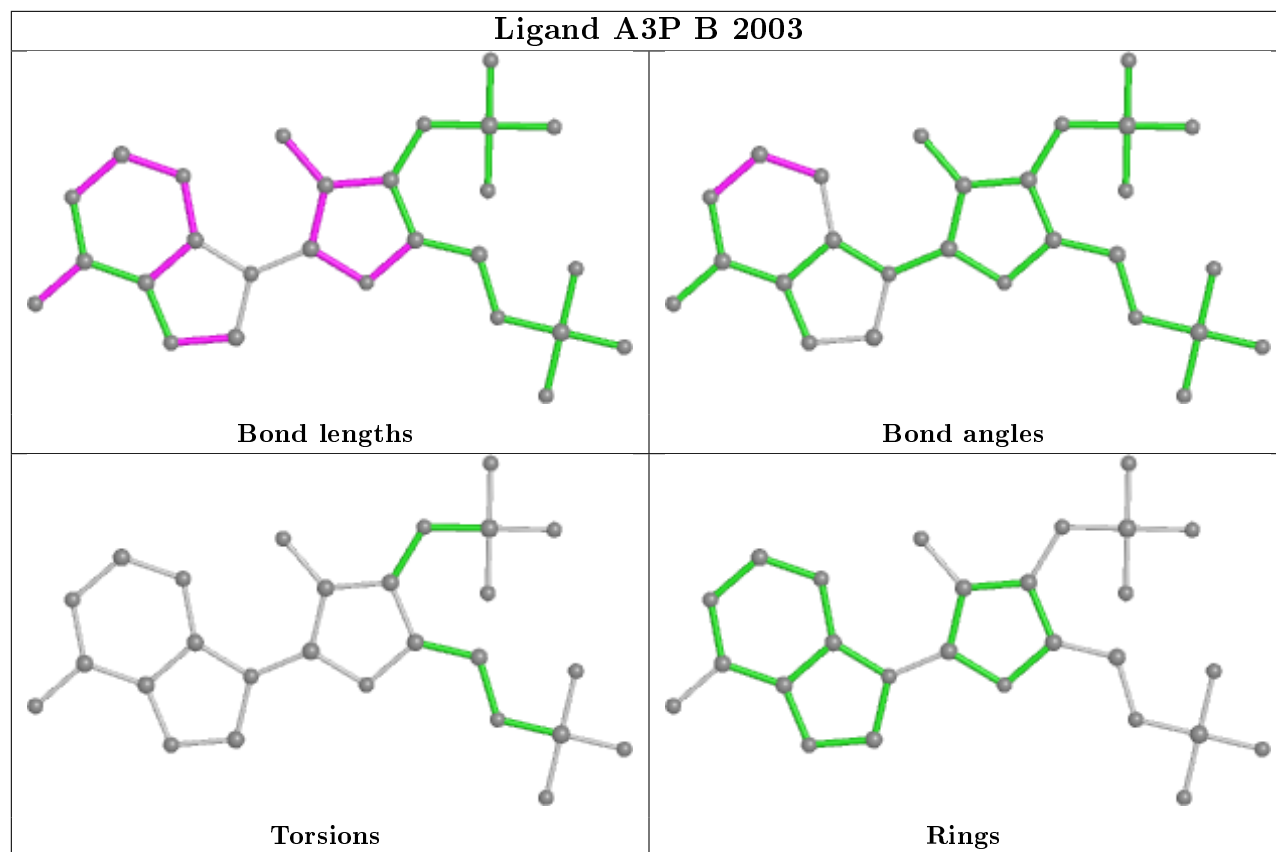
Mol	Chain	Res	Type	Atoms
4	F	2003	A3P	C3'-C4'-C5'-O5'

There are no ring outliers.

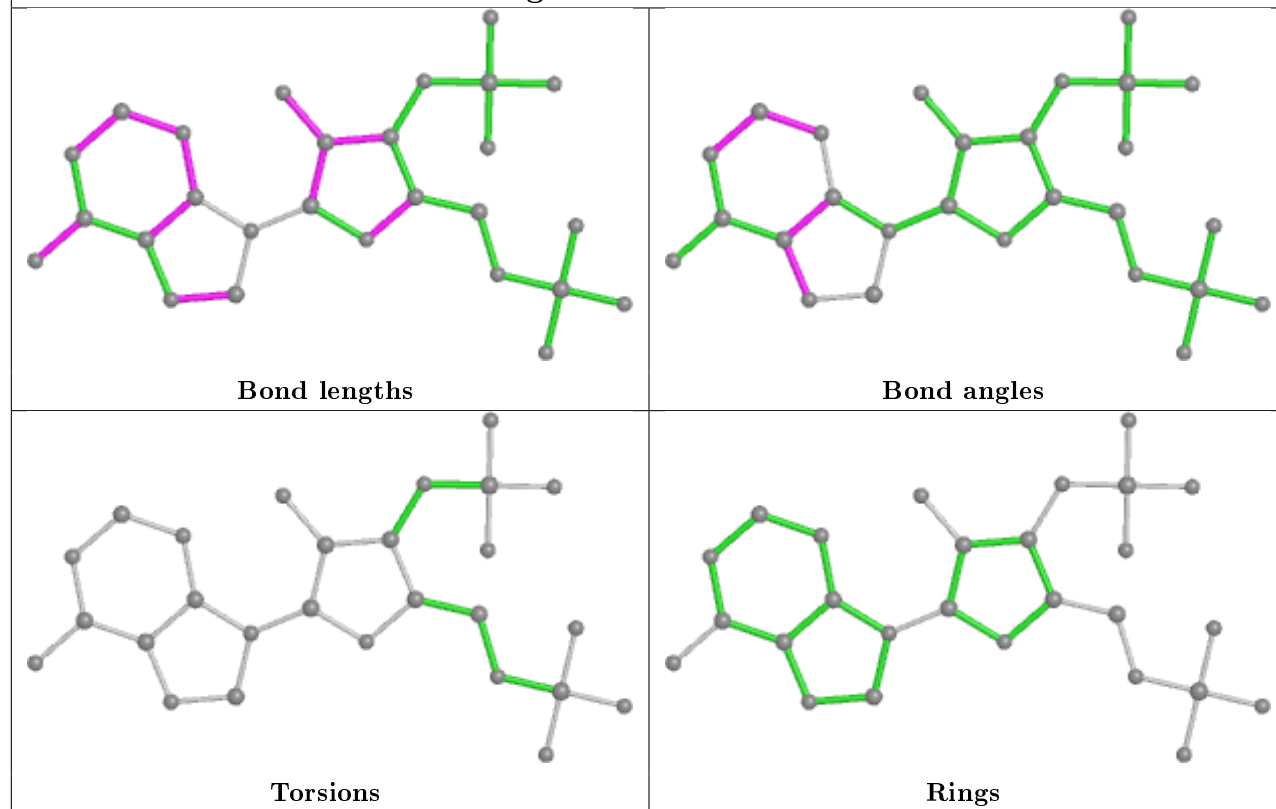
No monomer is involved in short contacts.

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 than 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.

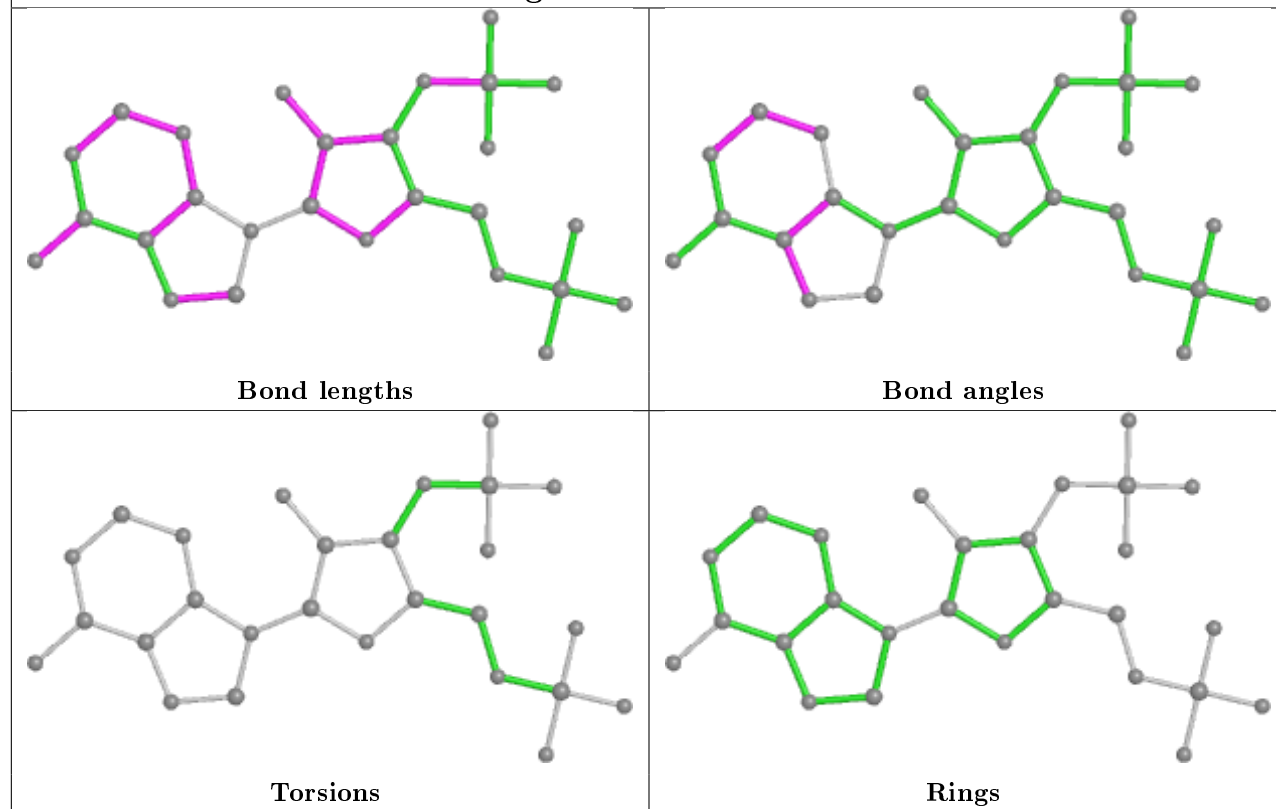


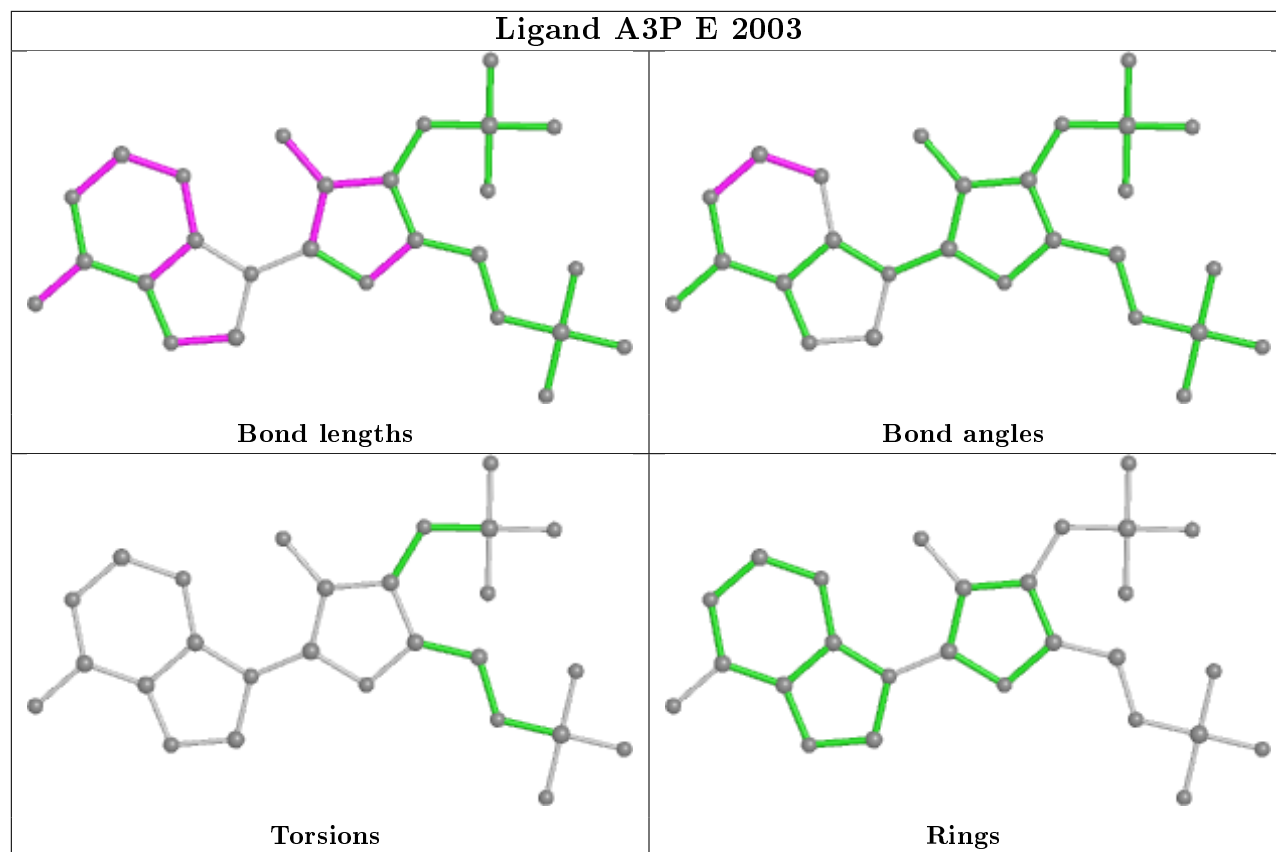


Ligand A3P C 2003



Ligand A3P A 2003





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	287/658 (43%)	-0.09	1 (0%) 94 92	48, 83, 133, 185	0
1	B	654/658 (99%)	1.41	180 (27%) 0 0	54, 196, 294, 399	0
1	C	655/658 (99%)	0.46	68 (10%) 6 8	63, 129, 207, 293	0
1	D	654/658 (99%)	0.34	41 (6%) 20 20	54, 128, 195, 252	0
1	E	654/658 (99%)	0.25	31 (4%) 31 31	58, 122, 191, 274	0
1	F	654/658 (99%)	1.06	150 (22%) 0 1	57, 170, 263, 333	0
All	All	3558/3948 (90%)	0.64	471 (13%) 3 5	48, 127, 250, 399	0

The worst 5 of 471 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	266	ILE	24.8
1	B	7	LEU	17.5
1	B	51	ALA	16.0
1	B	35	VAL	11.6
1	F	111	GLU	9.9

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.

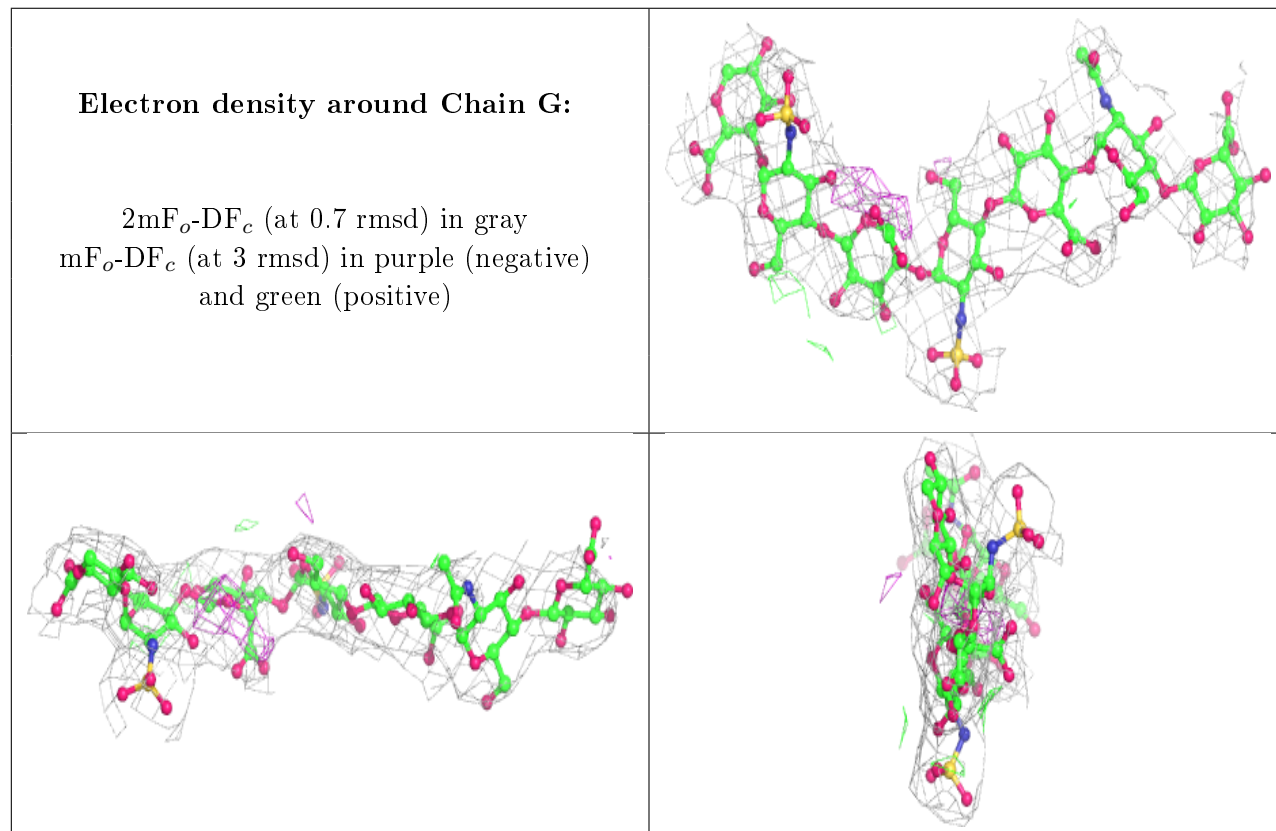
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	BDP	P	5	12/13	0.68	0.24	136,206,212,214	0
2	BDP	P	1	13/13	0.69	0.39	120,143,173,179	0
3	GLC	H	2	11/12	0.72	0.35	218,235,256,257	0
2	BDP	P	7	12/13	0.72	0.39	195,218,230,230	0
2	BDP	N	5	12/13	0.75	0.44	145,194,213,217	0
2	BDP	N	7	12/13	0.77	0.45	186,205,213,218	0
3	GLC	H	1	12/12	0.77	0.27	197,218,230,239	0
2	BDP	G	7	12/13	0.78	0.39	179,197,225,229	0
2	BDP	L	7	12/13	0.81	0.25	172,199,213,219	0
2	BDP	L	5	12/13	0.83	0.20	126,144,163,173	0
2	NDG	P	6	14/15	0.83	0.22	145,173,202,210	0
2	BDP	N	1	12/13	0.85	0.18	146,191,216,218	0
2	NDG	N	6	14/15	0.87	0.30	139,188,202,212	0
2	IDR	P	3	12/13	0.87	0.23	142,162,187,200	0
2	BDP	I	7	12/13	0.88	0.16	148,161,172,176	0
2	BDP	G	5	12/13	0.88	0.15	95,130,153,167	0
3	GLC	O	2	11/12	0.89	0.31	149,166,181,185	0
2	GNS	P	4	15/16	0.90	0.28	112,150,186,198	0
2	NDG	L	6	14/15	0.90	0.18	104,145,162,178	0
3	GLC	K	1	12/12	0.90	0.28	135,142,153,154	0
2	GNS	P	2	15/16	0.91	0.53	177,183,197,203	0
3	GLC	O	1	12/12	0.91	0.24	165,176,181,186	0
2	BDP	I	5	12/13	0.91	0.14	108,137,152,153	0
2	NDG	G	6	14/15	0.92	0.17	107,132,150,170	0
2	NDG	I	6	14/15	0.93	0.13	109,128,158,160	0
3	GLC	K	2	11/12	0.93	0.25	92,109,127,127	0
2	BDP	L	1	12/13	0.93	0.14	103,127,150,155	0
2	GNS	L	2	15/16	0.94	0.18	100,120,139,152	0
2	GNS	N	2	15/16	0.94	0.28	150,164,191,196	0
2	IDR	I	3	12/13	0.94	0.20	100,112,121,124	0
2	IDR	N	3	12/13	0.95	0.31	128,148,209,213	0
3	GLC	M	1	12/12	0.96	0.33	88,98,112,113	0
2	GNS	L	4	15/16	0.96	0.19	88,117,142,145	0
2	BDP	G	1	12/13	0.96	0.15	68,93,105,108	0
2	GNS	I	2	15/16	0.96	0.18	84,89,108,118	0
2	GNS	N	4	15/16	0.96	0.24	123,144,168,187	0
3	GLC	J	1	12/12	0.96	0.30	95,124,137,149	0
2	IDR	L	3	12/13	0.97	0.19	103,126,151,165	0
2	GNS	G	2	15/16	0.97	0.17	63,75,99,100	0
2	IDR	G	3	12/13	0.97	0.24	68,88,103,113	0
2	BDP	I	1	12/13	0.97	0.18	77,94,107,109	0
3	GLC	J	2	11/12	0.98	0.13	93,97,110,112	0
2	GNS	I	4	15/16	0.98	0.18	79,93,110,133	0

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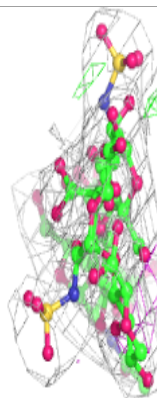
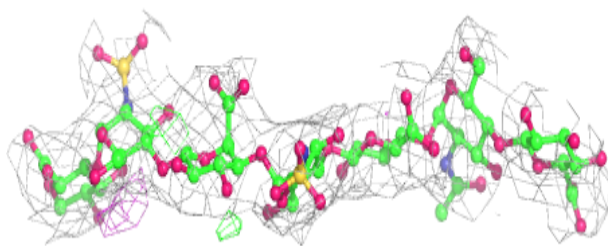
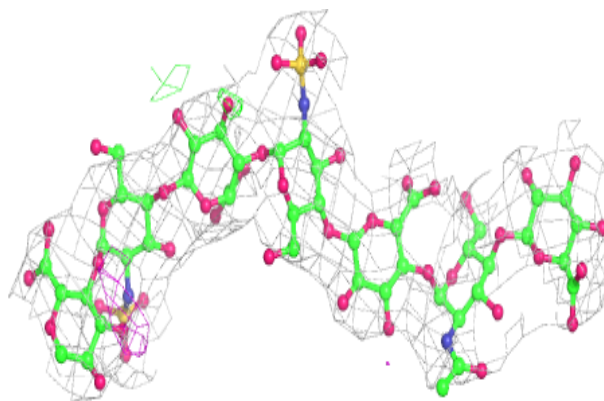
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	GNS	G	4	15/16	0.98	0.20	72,78,98,101	0
3	GLC	M	2	11/12	0.98	0.20	79,95,124,133	0

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.

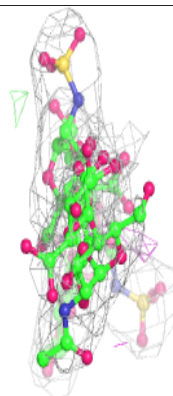
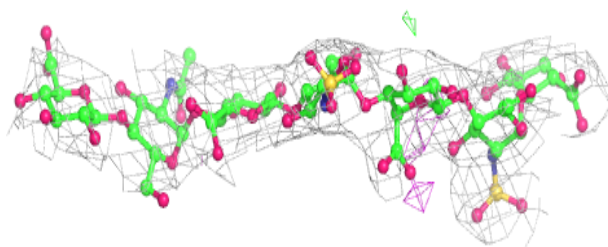
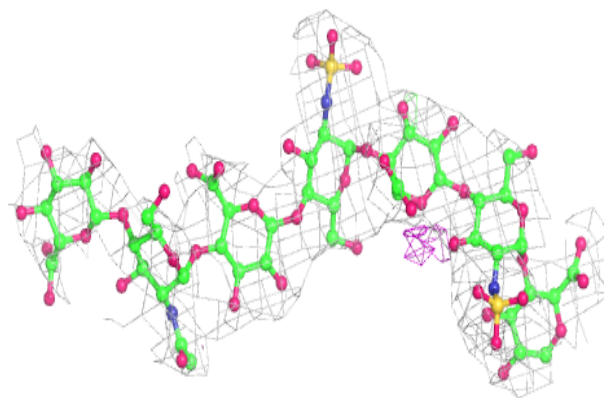


Electron density around Chain I:

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

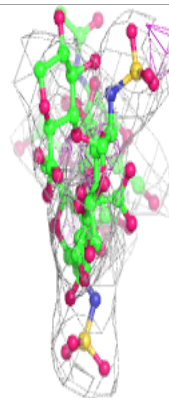
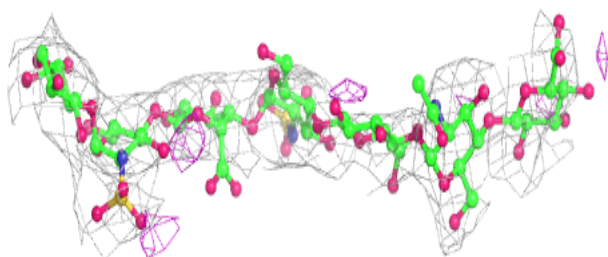
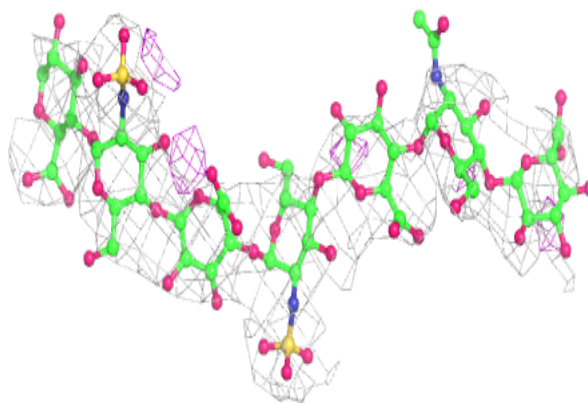
**Electron density around Chain L:**

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

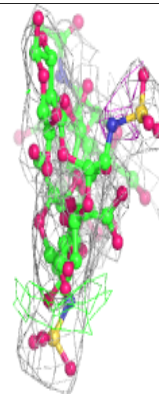
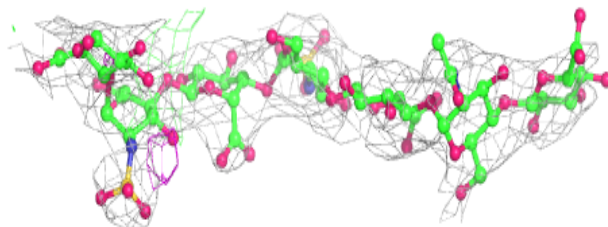
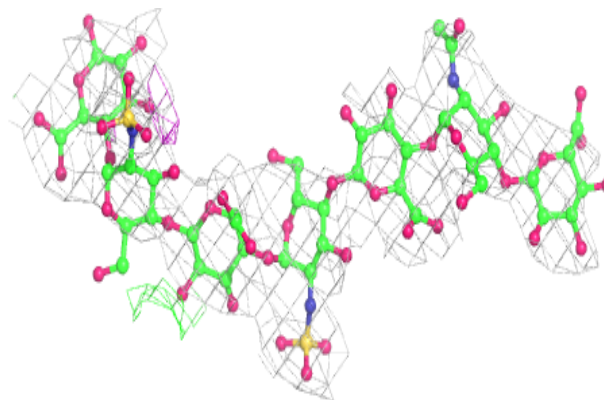


Electron density around Chain N:

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

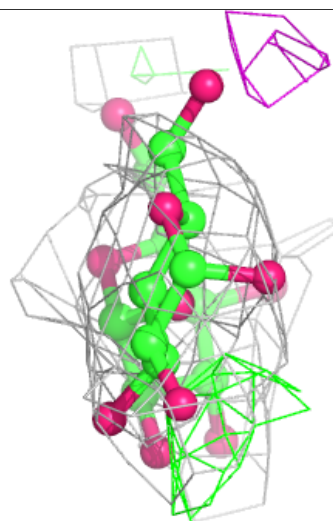
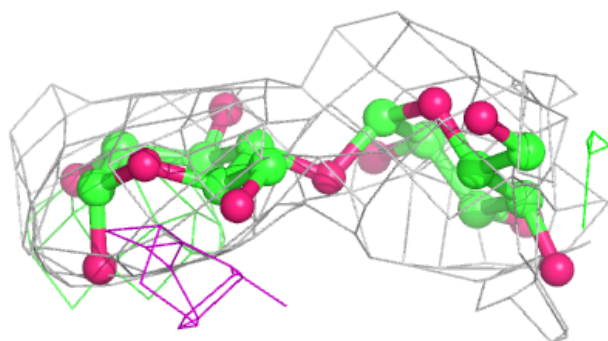
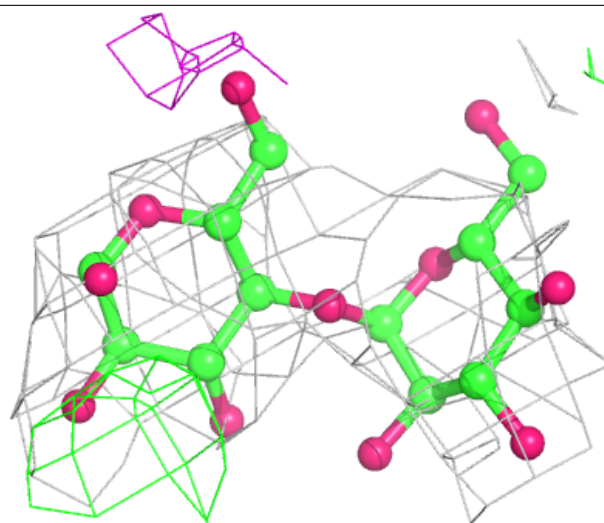
**Electron density around Chain P:**

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



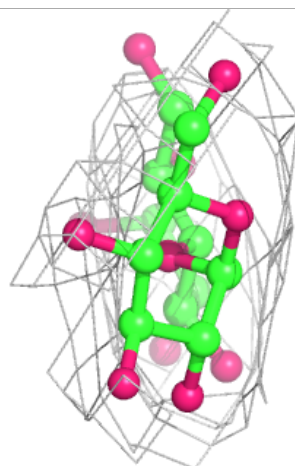
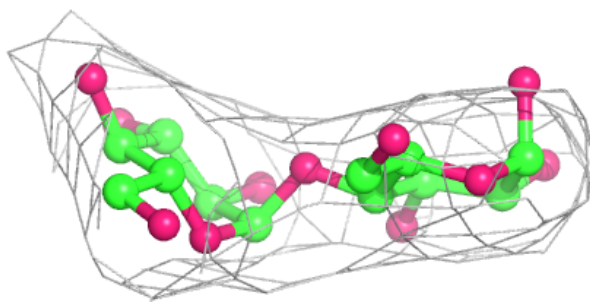
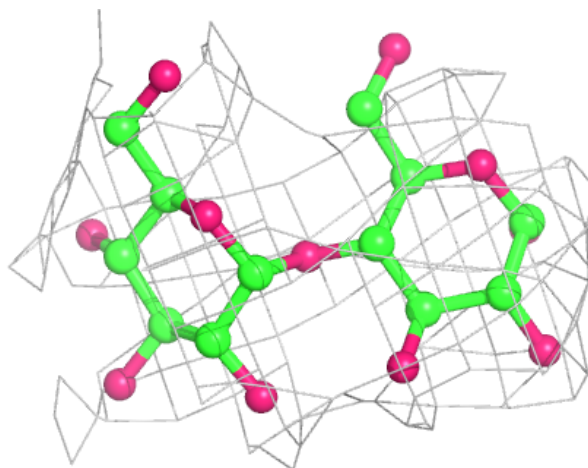
Electron density around Chain H:

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



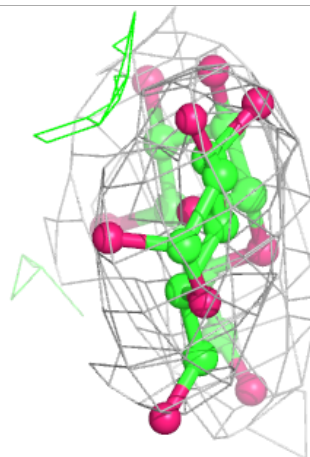
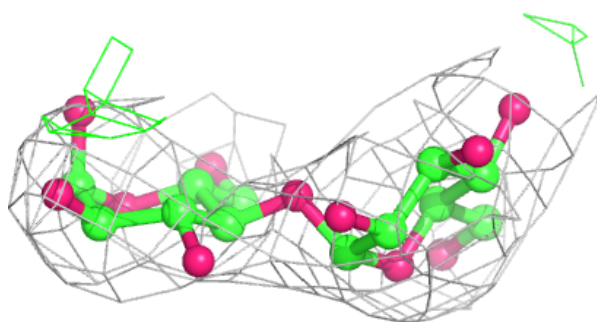
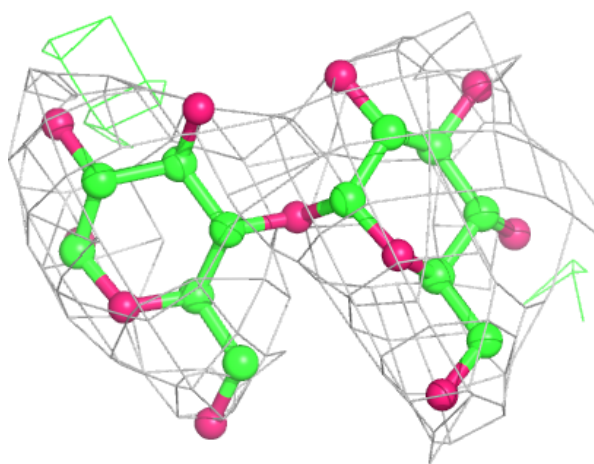
Electron density around Chain J:

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



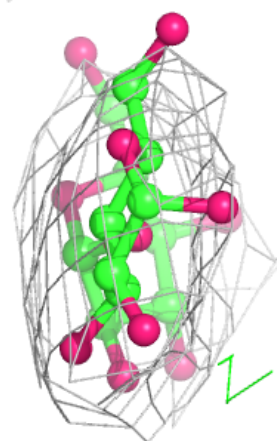
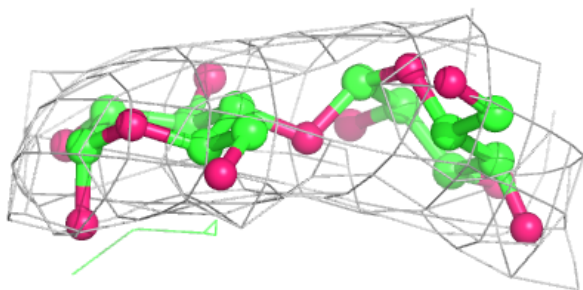
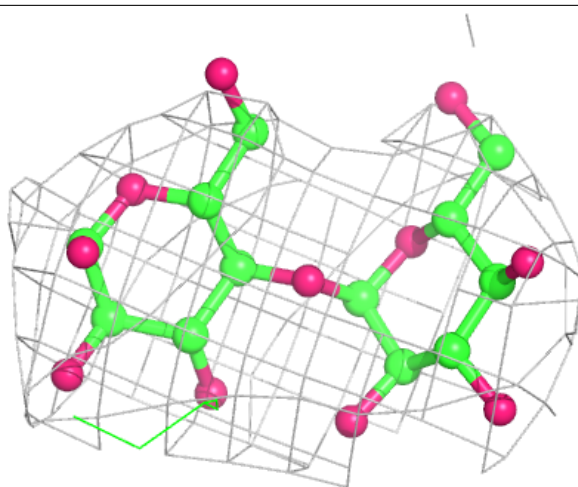
Electron density around Chain K:

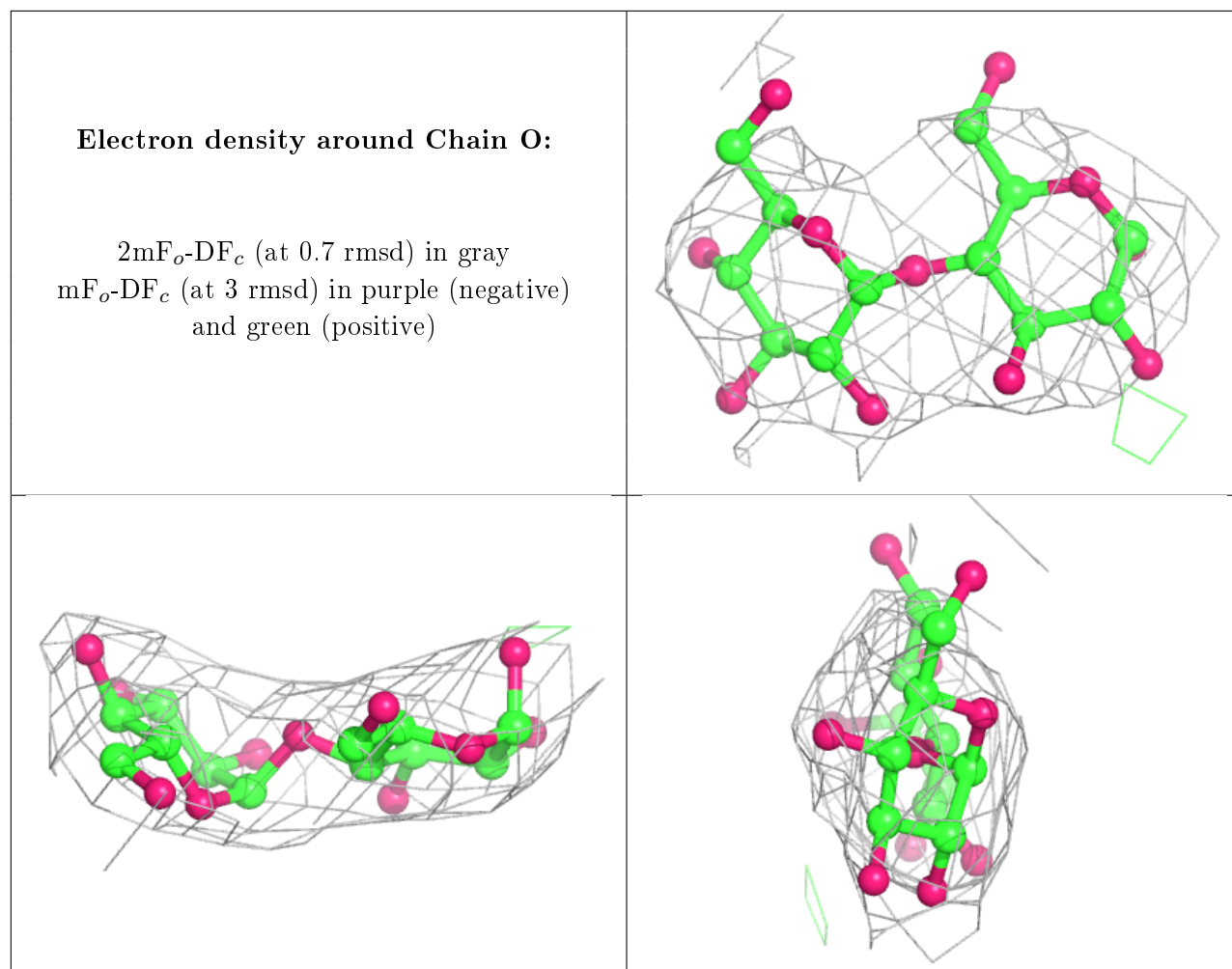
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain M:

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



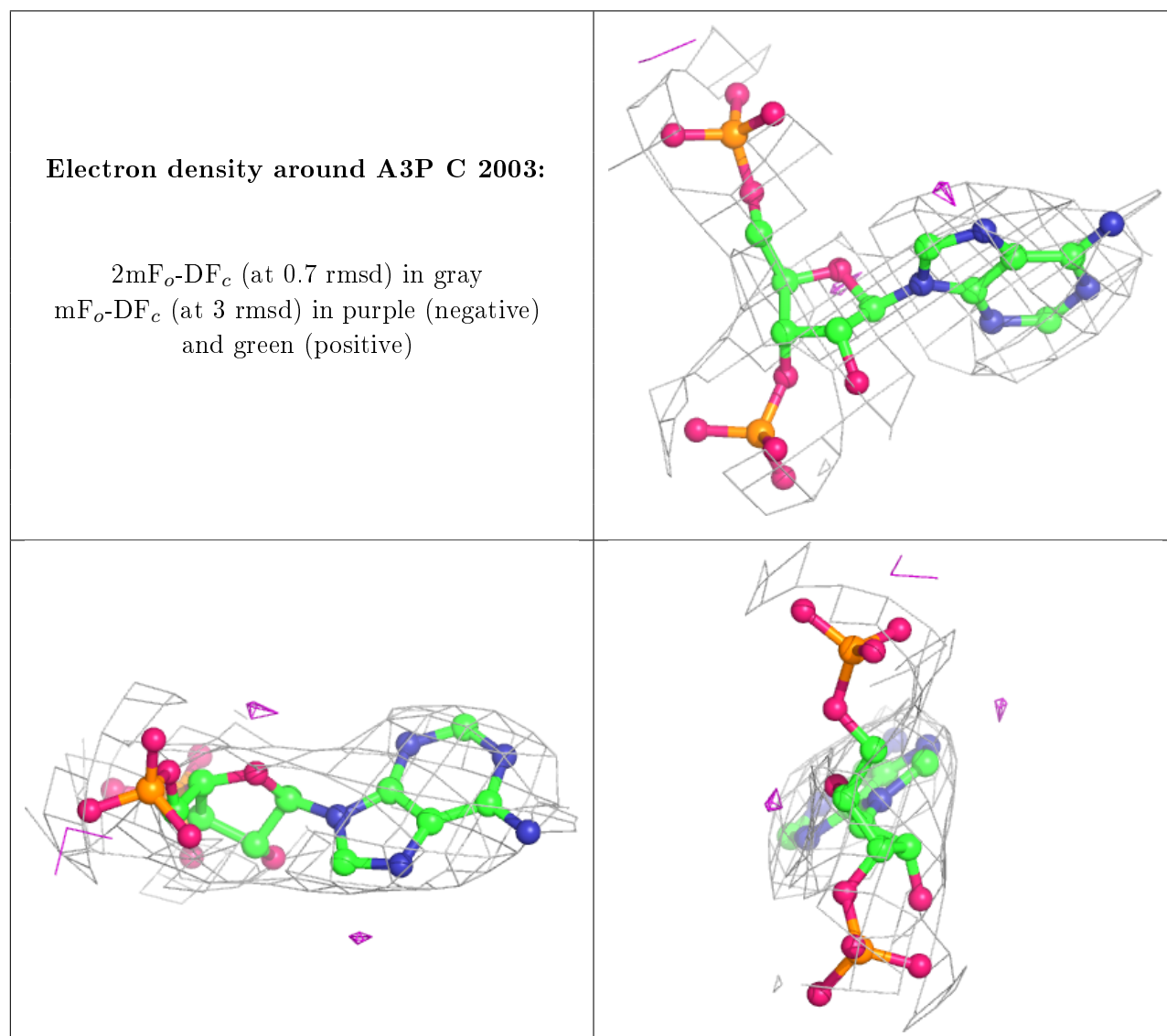


6.4 Ligands [i](#)

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.

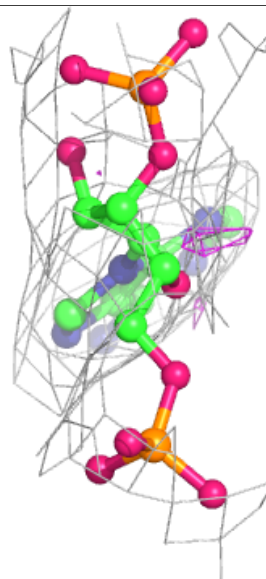
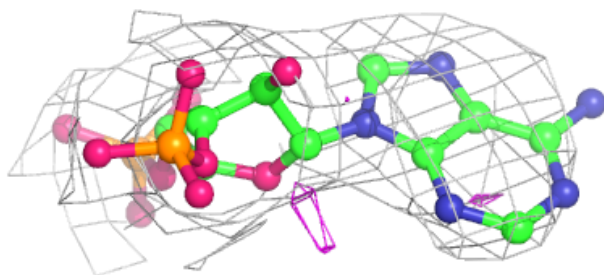
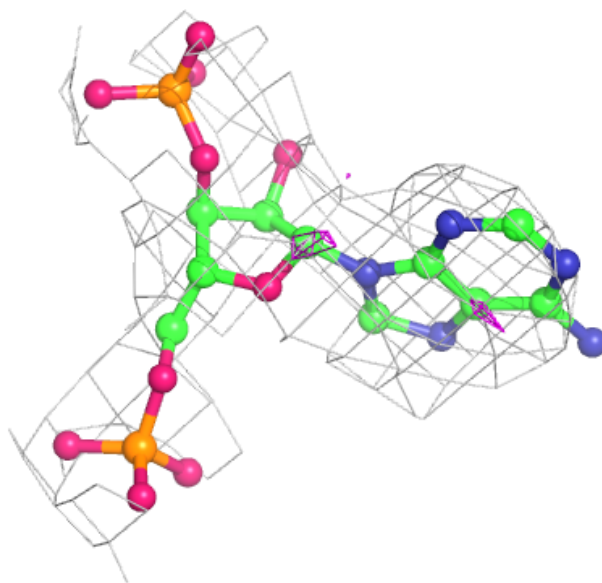
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	NPO	E	2011	10/10	0.73	0.37	157,172,185,189	0
5	NPO	D	2011	10/10	0.89	0.29	159,170,199,200	0
5	NPO	A	2011	10/10	0.92	0.31	103,138,146,154	0
4	A3P	C	2003	27/27	0.93	0.21	74,106,135,145	0
4	A3P	E	2003	27/27	0.94	0.27	80,104,114,119	0
4	A3P	F	2003	27/27	0.96	0.20	67,89,104,115	0
5	NPO	B	2011	10/10	0.96	0.30	105,120,169,173	0
4	A3P	A	2003	27/27	0.97	0.18	53,68,87,93	0
4	A3P	D	2003	27/27	0.97	0.20	57,83,108,112	0
4	A3P	B	2003	27/27	0.97	0.18	66,80,103,148	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.



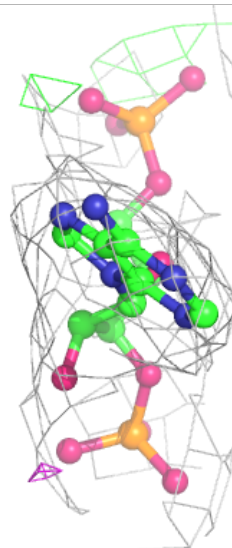
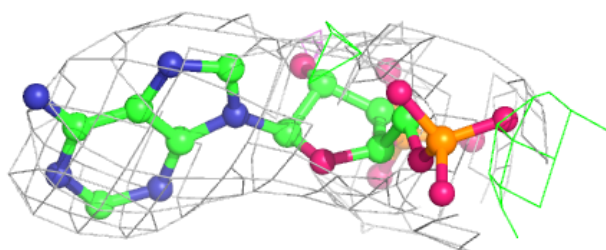
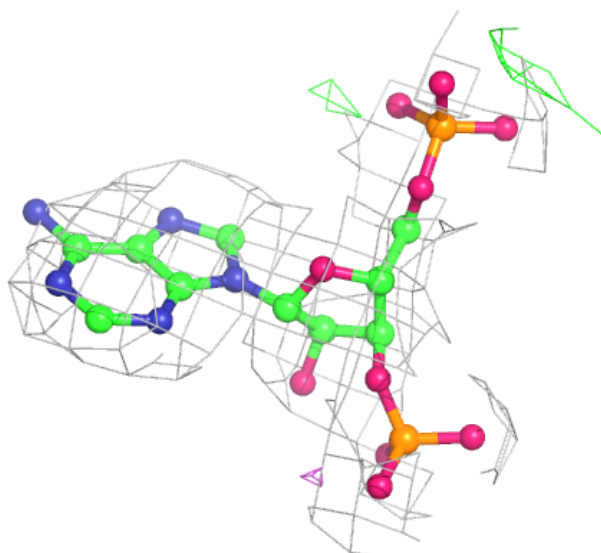
Electron density around A3P E 2003:

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



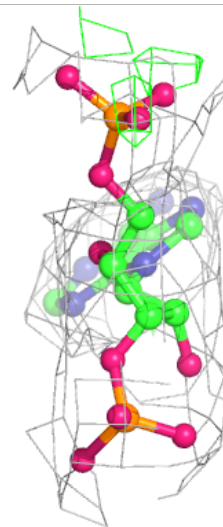
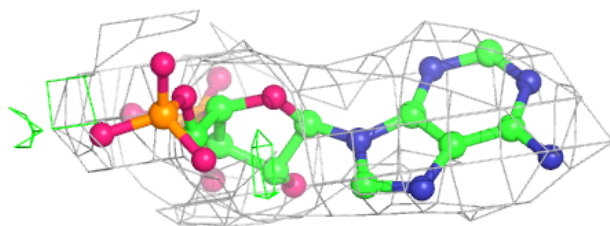
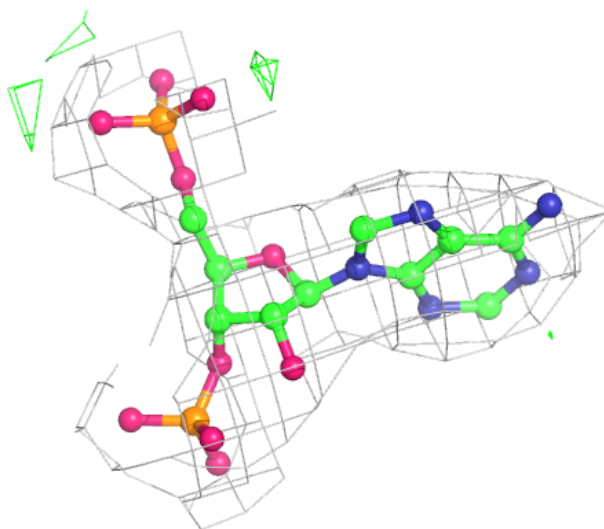
Electron density around A3P F 2003:

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



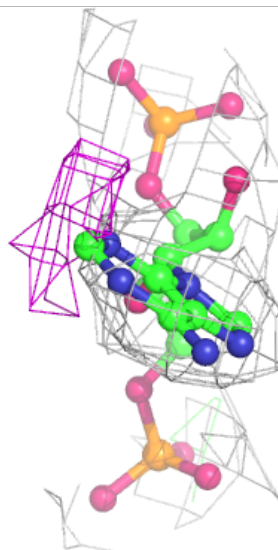
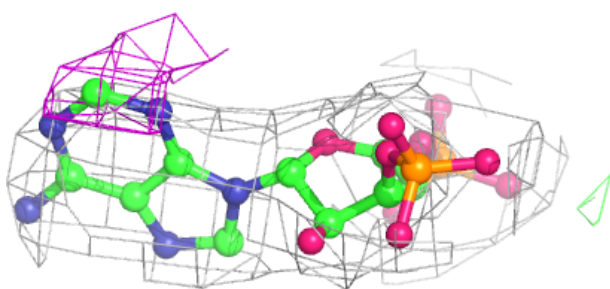
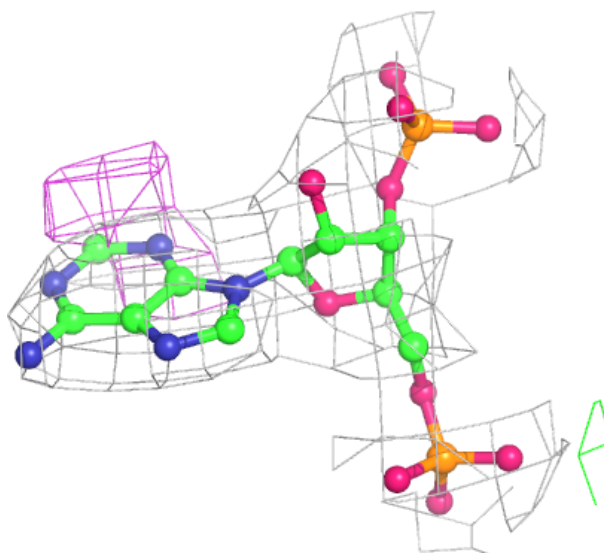
Electron density around A3P A 2003:

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



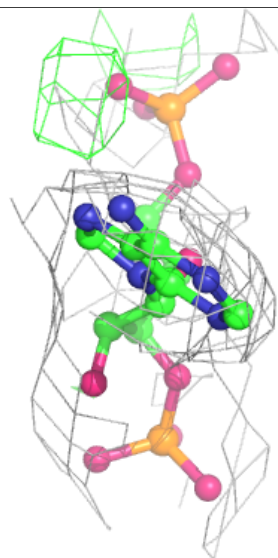
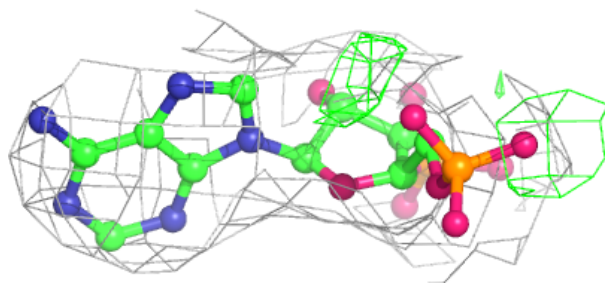
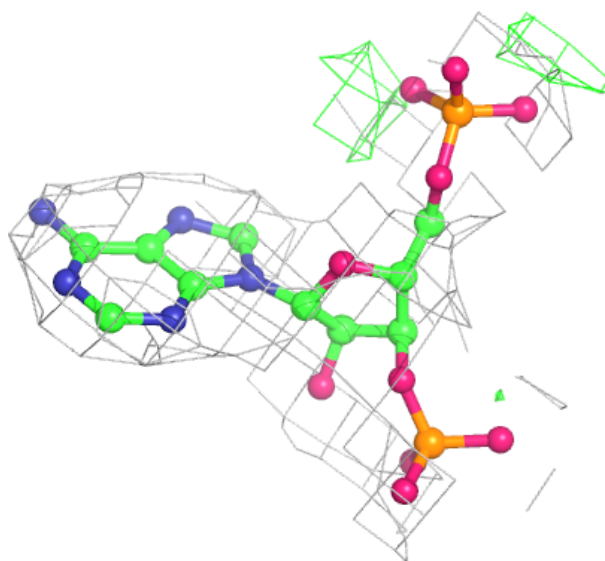
Electron density around A3P D 2003:

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



Electron density around A3P B 2003:

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.