Full wwPDB X-ray Structure Validation Report ⓘ

Dec 4, 2023 – 04:40 pm GMT

PDB ID	:	2BGR
Title	:	Crystal structure of HIV-1 Tat derived nonapeptides Tat(1-9) bound to the active site of Dipeptidyl peptidase IV (CD26)
Authors	:	Weihofen, W.A.; Liu, J.; Reutter, W.; Saenger, W.; Fan, H.
Deposited on	:	2005-01-04
Resolution	:	2.00 Å(reported)

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

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

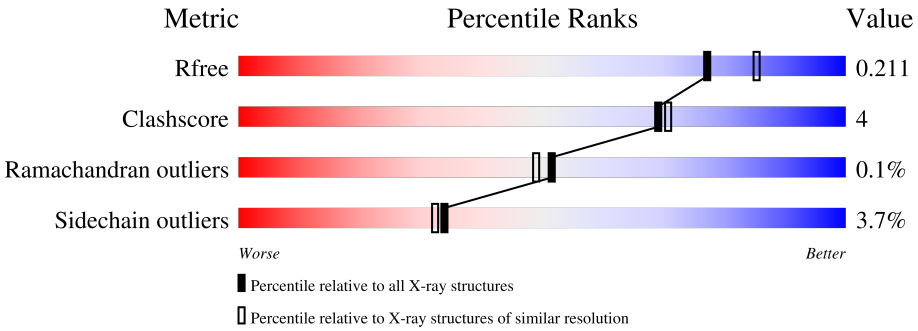
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	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 2.00 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	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)

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%

Mol	Chain	Length	Quality of chain
1	A	738	89% 8%
1	B	738	89% 8%
2	Y	9	11% 11% 11% 67%
2	Z	9	22% 11% 67%
3	C	3	100%
4	D	2	100%
4	F	2	100%

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Mol	Chain	Length	Quality of chain
4	G	2	100%
4	I	2	100%
4	K	2	100%
4	L	2	50% 50%
5	E	3	100%
5	J	3	33% 67%
6	H	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
3	FUC	C	3	X	-	-	-
6	FUC	H	2	X	-	-	-

2 Entry composition (i)

There are 8 unique types of molecules in this entry. The entry contains 14100 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.

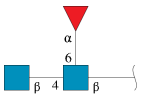
- Molecule 1 is a protein called DIPEPTIDYL PEPTIDASE IV.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	729	Total	C	N	O	S	0	0	0
			5972	3831	983	1132	26			
1	B	729	Total	C	N	O	S	0	0	0
			5972	3831	983	1132	26			

- Molecule 2 is a protein called HIV-1 TAT PROTEIN DERIVED N-TERMINAL NONAPEPTIDE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	Y	3	Total	C	N	O	S	0	0	0
			23	14	3	5	1			
2	Z	3	Total	C	N	O	S	0	0	0
			23	14	3	5	1			

- Molecule 3 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose.

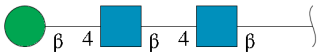


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	3	Total	C	N	O	0	0	0
			38	22	2	14			

- Molecule 4 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
4	D	2	Total C N O 28 16 2 10	0	0	0
4	F	2	Total C N O 28 16 2 10	0	0	0
4	G	2	Total C N O 28 16 2 10	0	0	0
4	I	2	Total C N O 28 16 2 10	0	0	0
4	K	2	Total C N O 28 16 2 10	0	0	0
4	L	2	Total C N O 28 16 2 10	0	0	0

- Molecule 5 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.

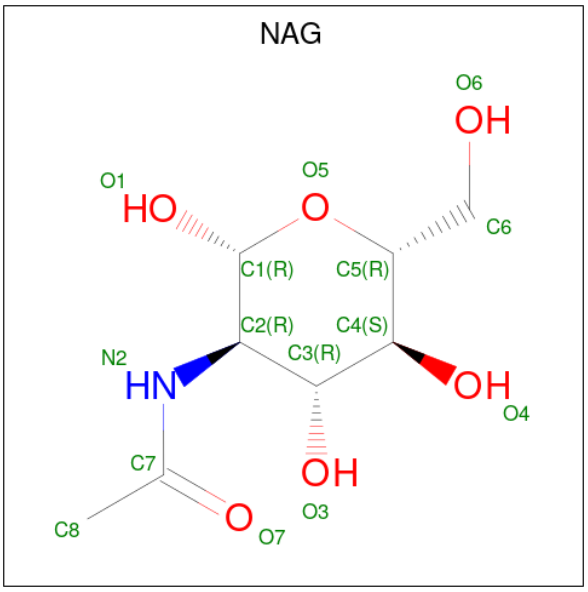


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf	Trace
5	E	3	Total C N O 39 22 2 15	0	0	0
5	J	3	Total C N O 39 22 2 15	0	0	0

- Molecule 6 is an oligosaccharide called alpha-L-fucopyranose-(1-6)-2-acetamido-2-deoxy-beta-D-glucopyranose.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf	Trace
6	H	2	Total C N O 24 14 1 9	0	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 C N O 14 8 1 5	0	0
7	A	1	Total C N O 14 8 1 5	0	0
7	B	1	Total C N O 14 8 1 5	0	0
7	B	1	Total C N O 14 8 1 5	0	0

- Molecule 8 is water.

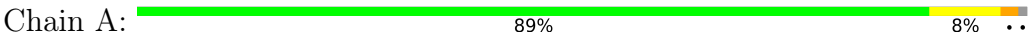
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	892	Total O 892 892	0	0
8	B	845	Total O 845 845	0	0
8	Y	6	Total O 6 6	0	0
8	Z	3	Total O 3 3	0	0

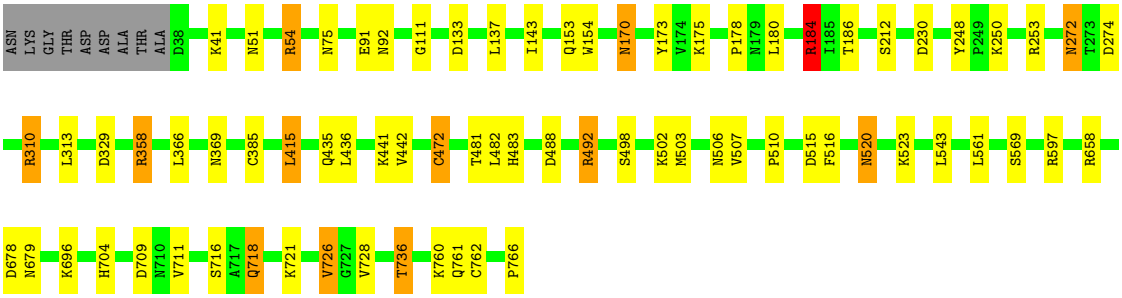
3 Residue-property plots (i)

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. 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. 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: DIPEPTIDYL PEPTIDASE IV

Chain A:



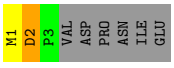


• Molecule 1: DIPEPTIDYL PEPTIDASE IV

Chain B:

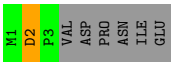
• Molecule 2: HIV-1 TAT PROTEIN DERIVED N-TERMINAL NONAPEPTIDE

Chain Y:



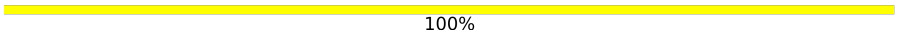
• Molecule 2: HIV-1 TAT PROTEIN DERIVED N-TERMINAL NONAPEPTIDE

Chain Z:



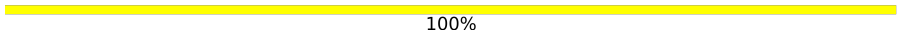
• Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain C:



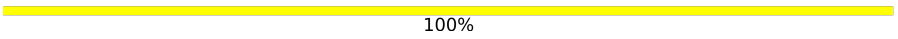
• Molecule 4: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain D:



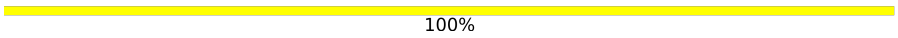
• Molecule 4: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain F:



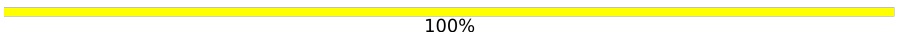
• Molecule 4: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain G:



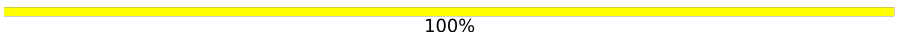
• Molecule 4: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain I:



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

Chain K:

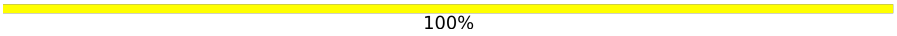


• Molecule 4: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain L:

• Molecule 5: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain E:



• Molecule 5: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain J:

• Molecule 6: alpha-L-fucopyranose-(1-6)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain H:

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	118.30Å 127.04Å 137.33Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.00 29.88 – 2.00	Depositor EDS
% Data completeness (in resolution range)	92.1 (30.00-2.00) 92.1 (29.88-2.00)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.35 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.1.1999	Depositor
R, R_{free}	0.160 , 0.203 0.170 , 0.211	Depositor DCC
R_{free} test set	1297 reflections (1.01%)	wwPDB-VP
Wilson B-factor (Å^2)	30.8	Xtriage
Anisotropy	0.145	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 39.8	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	14100	wwPDB-VP
Average B, all atoms (Å^2)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.62% 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: BMA, NAG, FUC

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	0.50	0/6144	0.86	15/8355 (0.2%)
1	B	0.49	0/6144	0.85	13/8355 (0.2%)
2	Y	0.52	0/23	1.51	1/30 (3.3%)
2	Z	0.45	0/23	1.30	0/30
All	All	0.50	0/12334	0.86	29/16770 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	1	0

There are no bond length outliers.

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	669	ARG	NE-CZ-NH2	-14.38	113.11	120.30
1	B	669	ARG	NE-CZ-NH1	11.74	126.17	120.30
1	A	492	ARG	NE-CZ-NH2	-10.91	114.85	120.30
1	A	658	ARG	NE-CZ-NH2	-8.26	116.17	120.30
1	A	658	ARG	NE-CZ-NH1	8.15	124.38	120.30
1	A	515	ASP	CB-CG-OD2	7.92	125.43	118.30
1	A	492	ARG	NE-CZ-NH1	7.88	124.24	120.30
1	A	310	ARG	NE-CZ-NH2	-7.75	116.42	120.30
1	A	184	ARG	NE-CZ-NH1	7.44	124.02	120.30
1	B	678	ASP	CB-CG-OD2	7.29	124.86	118.30
1	A	415	LEU	CA-CB-CG	7.24	131.95	115.30
1	B	691	ARG	NE-CZ-NH1	7.22	123.91	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	709	ASP	CB-CG-OD2	6.47	124.12	118.30
1	A	133	ASP	CB-CG-OD2	6.42	124.08	118.30
1	A	184	ARG	NE-CZ-NH2	-6.14	117.23	120.30
1	A	678	ASP	CB-CG-OD2	6.14	123.83	118.30
1	A	310	ARG	NE-CZ-NH1	6.07	123.33	120.30
1	A	415	LEU	CB-CG-CD1	5.92	121.07	111.00
1	B	358	ARG	NE-CZ-NH1	5.72	123.16	120.30
2	Y	2	ASP	CB-CG-OD2	5.60	123.34	118.30
1	A	709	ASP	CB-CG-OD2	5.59	123.33	118.30
1	B	535	ASP	CB-CG-OD2	5.50	123.25	118.30
1	B	691	ARG	NE-CZ-NH2	-5.45	117.57	120.30
1	B	501	ASP	CB-CG-OD2	5.29	123.06	118.30
1	B	358	ARG	NE-CZ-NH2	-5.27	117.67	120.30
1	B	60	LEU	CA-CB-CG	5.22	127.31	115.30
1	B	708	ASP	CB-CG-OD1	5.21	122.99	118.30
1	A	488	ASP	CB-CG-OD2	5.16	122.95	118.30
1	B	171	ASP	CB-CG-OD2	5.09	122.88	118.30

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	B	304	THR	CB

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	5972	0	5682	46	0
1	B	5972	0	5682	43	0
2	Y	23	0	22	1	0
2	Z	23	0	22	0	0
3	C	38	0	34	0	0
4	D	28	0	25	0	0
4	F	28	0	25	0	0
4	G	28	0	25	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	I	28	0	25	0	0
4	K	28	0	25	0	0
4	L	28	0	25	1	0
5	E	39	0	34	0	0
5	J	39	0	34	0	0
6	H	24	0	22	0	0
7	A	28	0	26	1	0
7	B	28	0	26	2	0
8	A	892	0	0	8	0
8	B	845	0	0	11	0
8	Y	6	0	0	1	0
8	Z	3	0	0	0	0
All	All	14100	0	11734	88	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.

All (88) 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:472:CYS:SG	8:A:2625:HOH:O	2.01	1.15
1:A:762:CYS:HB2	8:A:2874:HOH:O	1.48	1.12
1:B:472:CYS:SG	8:B:2571:HOH:O	2.14	1.06
1:A:503:MET:HG3	8:A:2028:HOH:O	1.76	0.85
1:B:762:CYS:HB2	8:B:2766:HOH:O	1.75	0.85
1:A:310:ARG:HD3	1:A:329:ASP:OD1	1.77	0.84
1:A:736:THR:HG21	1:B:717:ALA:O	1.85	0.75
1:B:320:GLN:OE1	1:B:669:ARG:HD3	1.88	0.73
1:B:74:ASN:HD22	7:B:1779:NAG:H81	1.53	0.71
1:B:304:THR:HG22	8:B:2454:HOH:O	1.92	0.70
1:B:762:CYS:CB	8:B:2766:HOH:O	2.36	0.69
1:A:272:ASN:HD22	1:A:274:ASP:H	1.40	0.68
1:A:272:ASN:ND2	1:A:274:ASP:H	1.93	0.66
1:A:310:ARG:CD	1:A:329:ASP:OD1	2.43	0.66
1:A:704:HIS:HD2	1:A:716:SER:OG	1.78	0.66
1:A:173:TYR:CE2	1:A:184:ARG:HG3	2.32	0.65
1:B:760:LYS:HE2	1:B:766:PRO:OXT	1.99	0.62
1:B:272:ASN:HD21	1:B:274:ASP:HB2	1.64	0.62
1:A:153:GLN:HE22	1:A:170:ASN:ND2	1.98	0.62
1:B:258:LYS:NZ	1:B:712:HIS:HD2	1.99	0.61
1:B:377:ASN:C	1:B:377:ASN:HD22	2.03	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:435:GLN:NE2	1:A:441:LYS:HD2	2.16	0.60
1:B:272:ASN:C	1:B:272:ASN:HD22	2.05	0.60
1:A:184:ARG:HD3	1:A:186:THR:O	2.01	0.60
1:A:760:LYS:CE	1:A:766:PRO:OXT	2.51	0.59
1:A:760:LYS:NZ	1:A:766:PRO:OXT	2.35	0.58
1:A:435:GLN:HE22	1:A:441:LYS:HD2	1.69	0.58
1:A:696:LYS:HG2	1:A:728:VAL:HG22	1.85	0.58
1:B:203:TYR:HA	1:B:207:VAL:HG13	1.86	0.57
2:Y:1:MET:HA	8:Y:2004:HOH:O	2.04	0.57
1:B:65:ASP:OD2	1:B:466:LYS:HB2	2.06	0.56
1:A:111:GLY:O	1:A:137:LEU:HD12	2.05	0.56
1:A:272:ASN:HD21	1:A:274:ASP:HB2	1.70	0.56
1:B:472:CYS:CB	8:B:2571:HOH:O	2.49	0.56
1:A:726:VAL:HG13	1:A:728:VAL:HG23	1.87	0.55
1:A:272:ASN:HD22	1:A:272:ASN:C	2.12	0.54
1:B:276:LEU:O	8:B:2348:HOH:O	2.19	0.53
1:B:302:ASP:OD1	1:B:304:THR:HG23	2.08	0.53
1:B:760:LYS:CE	1:B:766:PRO:OXT	2.58	0.52
1:B:304:THR:HG21	1:B:362:PRO:HD2	1.91	0.51
1:B:452:GLU:H	1:B:452:GLU:CD	2.14	0.51
1:B:691:ARG:HD2	8:B:2752:HOH:O	2.10	0.51
1:B:726:VAL:HG23	1:B:728:VAL:HG23	1.91	0.50
1:A:310:ARG:HH22	1:A:369:ASN:ND2	2.09	0.50
1:A:516:PHE:CE2	1:A:523:LYS:HD2	2.47	0.49
1:A:696:LYS:CG	1:A:728:VAL:HG22	2.42	0.48
1:A:762:CYS:CB	8:A:2874:HOH:O	2.30	0.48
1:A:481:THR:OG1	1:A:483:HIS:HE1	1.97	0.48
1:A:92:ASN:ND2	7:A:1780:NAG:C7	2.77	0.48
1:A:248:TYR:CZ	1:B:234:PRO:HB2	2.49	0.47
1:B:693:GLU:OE2	1:B:696:LYS:NZ	2.38	0.47
1:A:250:LYS:HD3	8:A:2351:HOH:O	2.14	0.47
1:A:503:MET:HE3	8:A:2650:HOH:O	2.15	0.47
1:B:640:LEU:HD11	1:B:650:GLY:HA3	1.97	0.47
1:A:597:ARG:HH12	1:A:679:ASN:HD21	1.62	0.47
1:A:704:HIS:HE1	1:A:711:VAL:O	1.97	0.47
1:B:710:ASN:C	1:B:710:ASN:HD22	2.19	0.46
1:B:170:ASN:ND2	8:B:2190:HOH:O	2.48	0.46
1:A:154:TRP:CE2	1:A:212:SER:HB3	2.50	0.46
1:A:253:ARG:HH21	1:B:253:ARG:HH21	1.63	0.46
1:B:258:LYS:HZ1	1:B:712:HIS:HD2	1.62	0.46
1:A:175:LYS:NZ	1:A:180:LEU:O	2.47	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:377:ASN:ND2	1:B:381:TYR:H	2.14	0.45
1:B:175:LYS:NZ	1:B:180:LEU:O	2.49	0.45
1:B:74:ASN:ND2	7:B:1779:NAG:H81	2.28	0.45
1:B:126:HIS:HD2	8:B:2057:HOH:O	1.98	0.45
1:B:153:GLN:HE22	1:B:170:ASN:ND2	2.15	0.44
1:A:520:ASN:O	1:A:520:ASN:OD1	2.36	0.44
1:B:41:LYS:HE3	1:B:53:TYR:OH	2.17	0.44
1:B:55:LEU:HD23	1:B:500:LEU:CD2	2.48	0.44
1:A:704:HIS:CD2	1:A:716:SER:OG	2.66	0.43
8:B:2844:HOH:O	4:L:2:NAG:O7	2.21	0.43
1:B:105:TYR:HB2	1:B:114:ILE:HD11	2.00	0.43
1:B:114:ILE:HG23	1:B:135:TYR:HB3	2.00	0.43
1:B:718:GLN:HE22	1:B:721:LYS:NZ	2.17	0.42
1:A:51:ASN:HB3	8:A:2017:HOH:O	2.20	0.42
1:B:272:ASN:HD22	1:B:274:ASP:H	1.67	0.42
1:A:718:GLN:HE22	1:A:721:LYS:NZ	2.17	0.42
1:B:472:CYS:CA	8:B:2571:HOH:O	2.67	0.42
1:A:510:PRO:HD3	1:A:569:SER:HB2	2.01	0.42
1:B:741:GLY:O	1:B:742:ILE:C	2.57	0.42
1:A:75:ASN:ND2	1:A:92:ASN:H	2.17	0.41
1:A:143:ILE:HD13	1:A:178:PRO:HB2	2.03	0.41
1:A:54:ARG:HE	1:A:54:ARG:HB2	1.69	0.41
1:A:170:ASN:HD22	1:A:170:ASN:N	2.19	0.41
1:A:498:SER:O	1:A:502:LYS:HG2	2.21	0.41
1:A:358:ARG:HD3	8:A:2100:HOH:O	2.21	0.40
1:B:472:CYS:O	1:B:478:PRO:HA	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	727/738 (98%)	702 (97%)	24 (3%)	1 (0%)	51 49
1	B	727/738 (98%)	703 (97%)	24 (3%)	0	100 100
2	Y	1/9 (11%)	1 (100%)	0	0	100 100
2	Z	1/9 (11%)	0	0	1 (100%)	0 0
All	All	1456/1494 (98%)	1406 (97%)	48 (3%)	2 (0%)	51 49

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	520	ASN
2	Z	2	ASP

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	654/660 (99%)	629 (96%)	25 (4%)	33 31
1	B	654/660 (99%)	632 (97%)	22 (3%)	37 36
2	Y	3/9 (33%)	2 (67%)	1 (33%)	0 0
2	Z	3/9 (33%)	2 (67%)	1 (33%)	0 0
All	All	1314/1338 (98%)	1265 (96%)	49 (4%)	34 32

All (49) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	41	LYS
1	A	54	ARG
1	A	91	GLU
1	A	170	ASN
1	A	184	ARG
1	A	230	ASP
1	A	272	ASN
1	A	313	LEU
1	A	358	ARG

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Mol	Chain	Res	Type
1	A	366	LEU
1	A	385	CYS
1	A	415	LEU
1	A	436	LEU
1	A	442	VAL
1	A	472	CYS
1	A	482	LEU
1	A	492	ARG
1	A	506	ASN
1	A	507	VAL
1	A	543	LEU
1	A	561	LEU
1	A	718	GLN
1	A	726	VAL
1	A	736	THR
1	A	761	GLN
1	B	41	LYS
1	B	60	LEU
1	B	170	ASN
1	B	207	VAL
1	B	223	LEU
1	B	230	ASP
1	B	272	ASN
1	B	358	ARG
1	B	377	ASN
1	B	385	CYS
1	B	445	LEU
1	B	452	GLU
1	B	471	ARG
1	B	472	CYS
1	B	502	LYS
1	B	521	GLU
1	B	543	LEU
1	B	627	TRP
1	B	673	LEU
1	B	679	ASN
1	B	710	ASN
1	B	718	GLN
2	Y	2	ASP
2	Z	2	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (36) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	75	ASN
1	A	169	ASN
1	A	170	ASN
1	A	247	GLN
1	A	272	ASN
1	A	314	GLN
1	A	345	HIS
1	A	369	ASN
1	A	435	GLN
1	A	483	HIS
1	A	572	ASN
1	A	586	GLN
1	A	679	ASN
1	A	685	ASN
1	A	694	ASN
1	A	704	HIS
1	A	718	GLN
1	A	761	GLN
1	B	112	GLN
1	B	126	HIS
1	B	169	ASN
1	B	170	ASN
1	B	247	GLN
1	B	272	ASN
1	B	314	GLN
1	B	345	HIS
1	B	377	ASN
1	B	435	GLN
1	B	572	ASN
1	B	606	GLN
1	B	679	ASN
1	B	694	ASN
1	B	710	ASN
1	B	712	HIS
1	B	718	GLN
1	B	731	GLN

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)

23 monosaccharides are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
3	NAG	C	1	1,3	14,14,15	0.78	0	17,19,21	1.51	3 (17%)
3	NAG	C	2	3	14,14,15	0.59	0	17,19,21	1.78	6 (35%)
3	FUC	C	3	3	10,10,11	0.64	0	14,14,16	1.26	1 (7%)
4	NAG	D	1	1,4	14,14,15	0.66	0	17,19,21	1.31	2 (11%)
4	NAG	D	2	4	14,14,15	0.45	0	17,19,21	1.44	1 (5%)
5	NAG	E	1	5,1	14,14,15	0.75	0	17,19,21	1.01	1 (5%)
5	NAG	E	2	5	14,14,15	0.59	0	17,19,21	1.34	2 (11%)
5	BMA	E	3	5	11,11,12	0.69	0	15,15,17	1.42	3 (20%)
4	NAG	F	1	1,4	14,14,15	0.62	0	17,19,21	1.90	4 (23%)
4	NAG	F	2	4	14,14,15	0.51	0	17,19,21	1.10	2 (11%)
4	NAG	G	1	1,4	14,14,15	0.61	0	17,19,21	1.28	3 (17%)
4	NAG	G	2	4	14,14,15	0.49	0	17,19,21	1.66	3 (17%)
6	NAG	H	1	1,6	14,14,15	0.64	0	17,19,21	1.10	1 (5%)
6	FUC	H	2	6	10,10,11	0.68	0	14,14,16	0.89	0
4	NAG	I	1	1,4	14,14,15	0.45	0	17,19,21	1.37	3 (17%)
4	NAG	I	2	4	14,14,15	0.65	0	17,19,21	1.61	3 (17%)
5	NAG	J	1	5,1	14,14,15	0.52	0	17,19,21	1.48	2 (11%)
5	NAG	J	2	5	14,14,15	0.57	0	17,19,21	1.03	1 (5%)
5	BMA	J	3	5	11,11,12	0.63	0	15,15,17	1.02	0
4	NAG	K	1	1,4	14,14,15	0.62	0	17,19,21	1.27	1 (5%)
4	NAG	K	2	4	14,14,15	0.57	0	17,19,21	1.15	2 (11%)
4	NAG	L	1	1,4	14,14,15	0.59	0	17,19,21	1.52	3 (17%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
4	NAG	L	2	4	14,14,15	0.64	0	17,19,21	1.48	3 (17%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	C	1	1,3	-	1/6/23/26	0/1/1/1
3	NAG	C	2	3	-	4/6/23/26	0/1/1/1
3	FUC	C	3	3	1/1/4/5	-	0/1/1/1
4	NAG	D	1	1,4	-	2/6/23/26	0/1/1/1
4	NAG	D	2	4	-	2/6/23/26	0/1/1/1
5	NAG	E	1	5,1	-	0/6/23/26	0/1/1/1
5	NAG	E	2	5	-	2/6/23/26	0/1/1/1
5	BMA	E	3	5	-	2/2/19/22	0/1/1/1
4	NAG	F	1	1,4	-	1/6/23/26	0/1/1/1
4	NAG	F	2	4	-	5/6/23/26	0/1/1/1
4	NAG	G	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	G	2	4	-	1/6/23/26	0/1/1/1
6	NAG	H	1	1,6	-	2/6/23/26	0/1/1/1
6	FUC	H	2	6	1/1/4/5	-	0/1/1/1
4	NAG	I	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	I	2	4	-	4/6/23/26	0/1/1/1
5	NAG	J	1	5,1	-	0/6/23/26	0/1/1/1
5	NAG	J	2	5	-	1/6/23/26	0/1/1/1
5	BMA	J	3	5	-	1/2/19/22	0/1/1/1
4	NAG	K	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	K	2	4	-	0/6/23/26	0/1/1/1
4	NAG	L	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	L	2	4	-	4/6/23/26	0/1/1/1

There are no bond length outliers.

All (50) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	F	1	NAG	C1-O5-C5	5.08	119.08	112.19
4	I	2	NAG	C4-C3-C2	4.61	117.77	111.02
4	D	2	NAG	C1-O5-C5	4.51	118.30	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	L	1	NAG	O5-C1-C2	-4.10	104.81	111.29
3	C	2	NAG	O5-C1-C2	-3.89	105.15	111.29
4	G	2	NAG	C1-O5-C5	3.85	117.41	112.19
4	G	2	NAG	C4-C3-C2	-3.71	105.58	111.02
5	J	1	NAG	O5-C1-C2	-3.70	105.44	111.29
4	L	2	NAG	C2-N2-C7	3.50	127.89	122.90
4	I	2	NAG	C3-C4-C5	3.44	116.38	110.24
3	C	1	NAG	O5-C1-C2	-3.42	105.88	111.29
3	C	2	NAG	C3-C4-C5	3.35	116.22	110.24
5	E	2	NAG	C8-C7-N2	3.33	121.73	116.10
4	I	1	NAG	O5-C1-C2	-3.29	106.09	111.29
3	C	1	NAG	C1-O5-C5	3.17	116.48	112.19
4	K	2	NAG	C1-O5-C5	3.14	116.45	112.19
4	G	1	NAG	O5-C1-C2	-3.14	106.33	111.29
5	E	3	BMA	C1-C2-C3	3.10	113.47	109.67
5	E	3	BMA	C2-C3-C4	3.00	116.09	110.89
4	K	1	NAG	C1-C2-N2	2.96	115.54	110.49
4	F	1	NAG	O5-C1-C2	-2.88	106.74	111.29
4	L	2	NAG	C4-C3-C2	2.83	115.17	111.02
4	L	1	NAG	C1-O5-C5	2.82	116.01	112.19
4	F	2	NAG	O5-C1-C2	-2.81	106.86	111.29
4	G	2	NAG	C1-C2-N2	2.65	115.02	110.49
3	C	3	FUC	O5-C5-C6	2.65	113.03	107.33
4	F	1	NAG	O5-C5-C6	-2.65	103.05	107.20
4	L	2	NAG	C1-C2-N2	-2.59	106.06	110.49
4	I	1	NAG	O5-C5-C6	2.48	111.09	107.20
3	C	2	NAG	C2-N2-C7	-2.47	119.39	122.90
3	C	2	NAG	C1-O5-C5	2.43	115.49	112.19
5	E	3	BMA	C3-C4-C5	2.42	114.56	110.24
4	D	1	NAG	O5-C1-C2	-2.41	107.48	111.29
3	C	1	NAG	O4-C4-C5	-2.37	103.40	109.30
4	G	1	NAG	C1-O5-C5	2.35	115.38	112.19
4	F	1	NAG	O4-C4-C3	-2.33	104.95	110.35
4	F	2	NAG	C2-N2-C7	2.32	126.21	122.90
5	E	2	NAG	O7-C7-C8	-2.32	117.75	122.06
3	C	2	NAG	O5-C5-C4	2.31	116.45	110.83
4	G	1	NAG	O7-C7-N2	2.29	126.16	121.95
4	D	1	NAG	C4-C3-C2	2.27	114.35	111.02
3	C	2	NAG	C1-C2-N2	2.26	114.35	110.49
4	K	2	NAG	C1-C2-N2	2.24	114.31	110.49
5	J	2	NAG	C4-C3-C2	2.23	114.28	111.02
5	J	1	NAG	O5-C5-C6	2.20	110.65	107.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	L	1	NAG	O7-C7-C8	-2.17	118.02	122.06
4	I	1	NAG	O7-C7-N2	2.16	125.93	121.95
5	E	1	NAG	O5-C5-C6	2.08	110.46	107.20
4	I	2	NAG	C1-O5-C5	2.03	114.94	112.19
6	H	1	NAG	O7-C7-C8	-2.01	118.32	122.06

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	C	3	FUC	C1
6	H	2	FUC	C1

All (32) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	2	NAG	C8-C7-N2-C2
3	C	2	NAG	O7-C7-N2-C2
4	F	2	NAG	C3-C2-N2-C7
4	L	2	NAG	C8-C7-N2-C2
4	L	2	NAG	O7-C7-N2-C2
5	E	2	NAG	C8-C7-N2-C2
5	E	2	NAG	O7-C7-N2-C2
3	C	2	NAG	O5-C5-C6-O6
4	D	1	NAG	C8-C7-N2-C2
4	F	2	NAG	C8-C7-N2-C2
4	I	2	NAG	C4-C5-C6-O6
4	I	2	NAG	O5-C5-C6-O6
6	H	1	NAG	O5-C5-C6-O6
4	D	2	NAG	C4-C5-C6-O6
4	F	2	NAG	O7-C7-N2-C2
4	F	2	NAG	C4-C5-C6-O6
5	E	3	BMA	C4-C5-C6-O6
3	C	2	NAG	C4-C5-C6-O6
6	H	1	NAG	C4-C5-C6-O6
4	D	1	NAG	O7-C7-N2-C2
5	E	3	BMA	O5-C5-C6-O6
4	D	2	NAG	O5-C5-C6-O6
4	F	2	NAG	O5-C5-C6-O6
4	I	2	NAG	C8-C7-N2-C2
4	L	2	NAG	O5-C5-C6-O6
4	F	1	NAG	O5-C5-C6-O6
4	I	2	NAG	O7-C7-N2-C2

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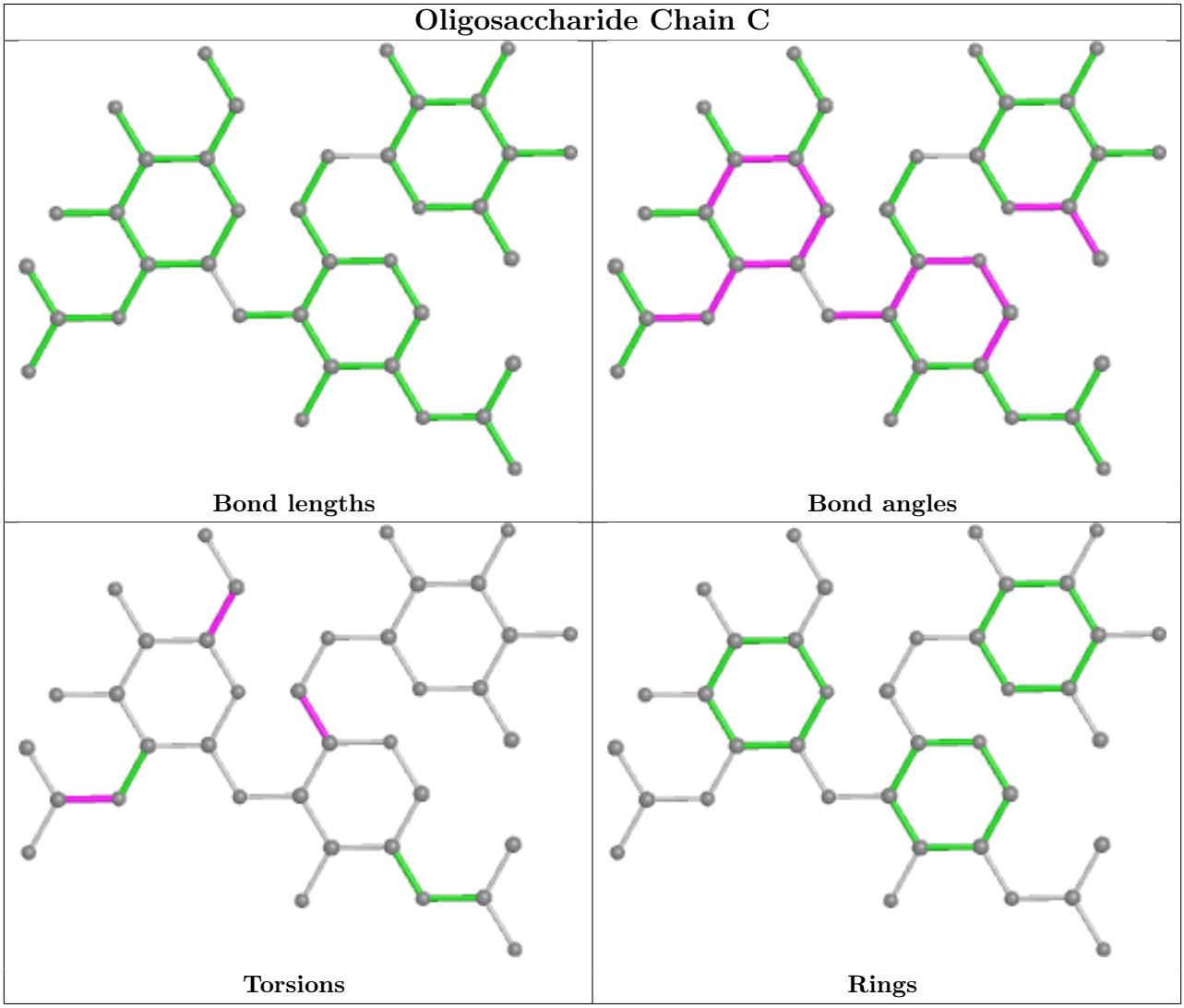
Mol	Chain	Res	Type	Atoms
4	L	2	NAG	C4-C5-C6-O6
5	J	3	BMA	C4-C5-C6-O6
3	C	1	NAG	O5-C5-C6-O6
4	G	2	NAG	C3-C2-N2-C7
5	J	2	NAG	C4-C5-C6-O6

There are no ring outliers.

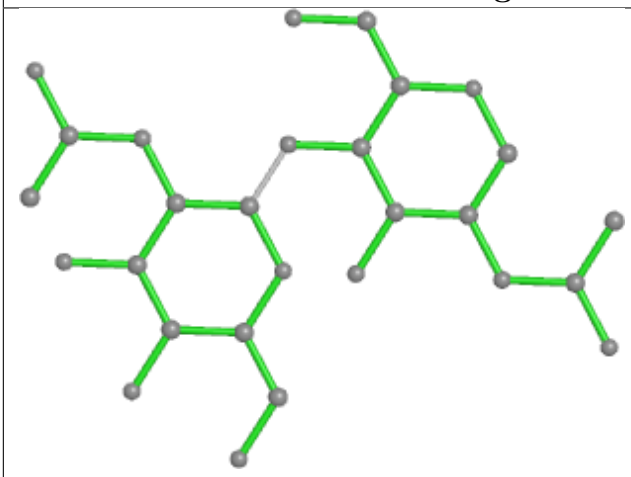
1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	L	2	NAG	1	0

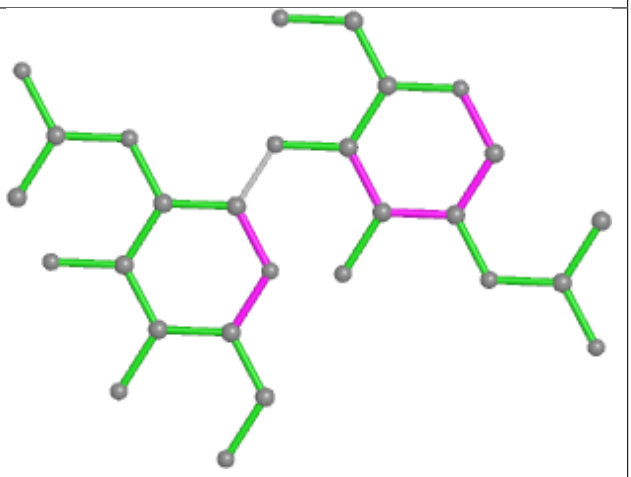
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.



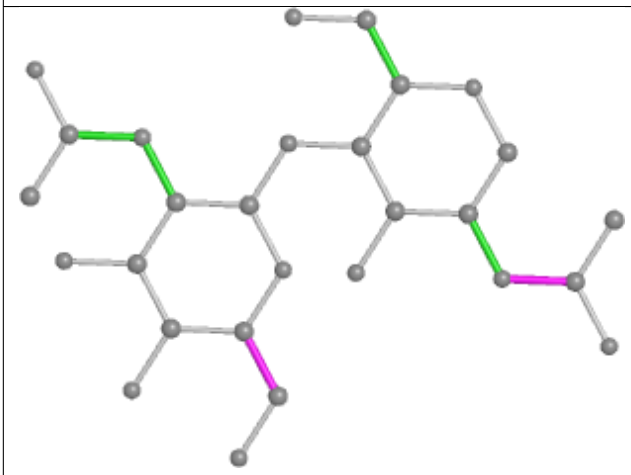
Oligosaccharide Chain D



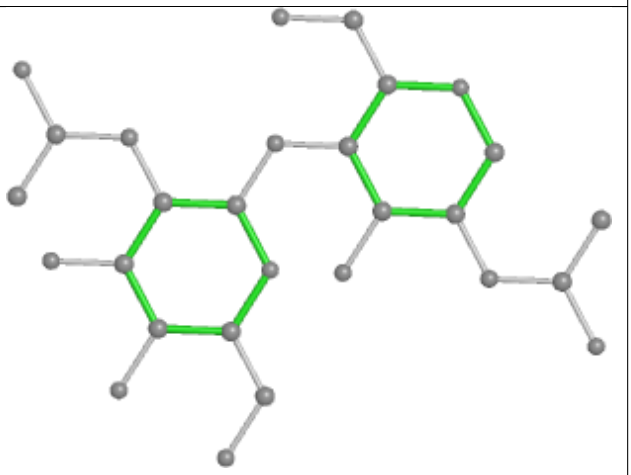
Bond lengths



Bond angles

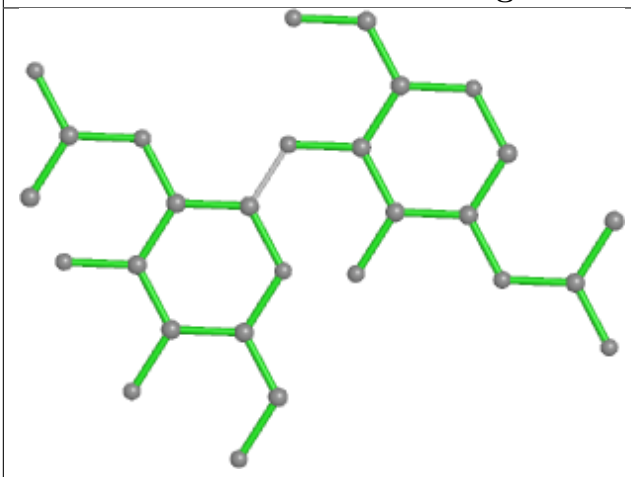


Torsions

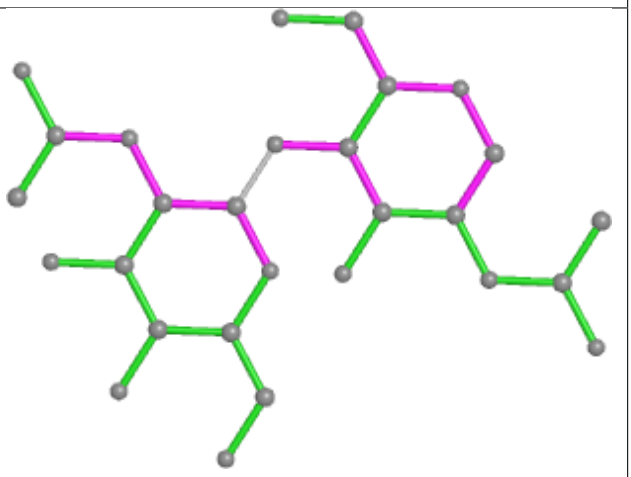


Rings

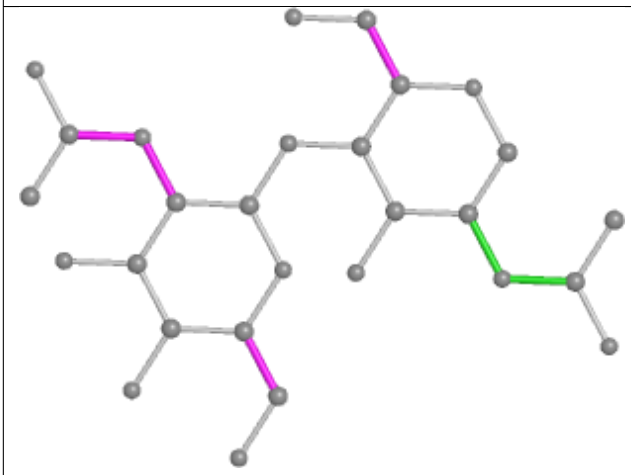
Oligosaccharide Chain F



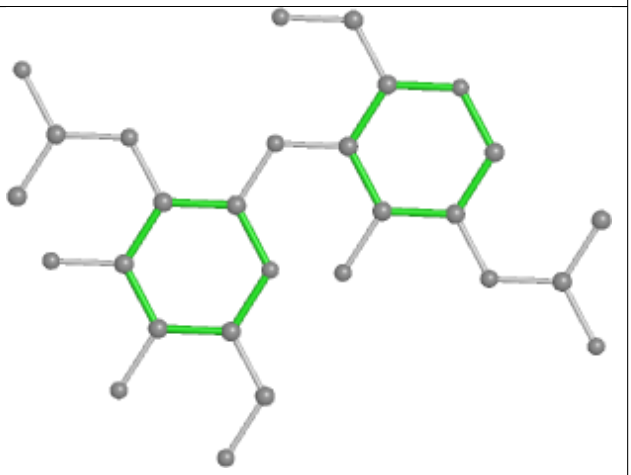
Bond lengths



Bond angles

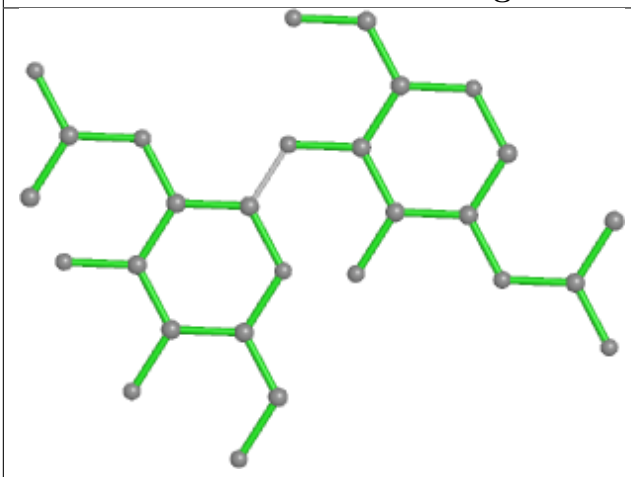


Torsions

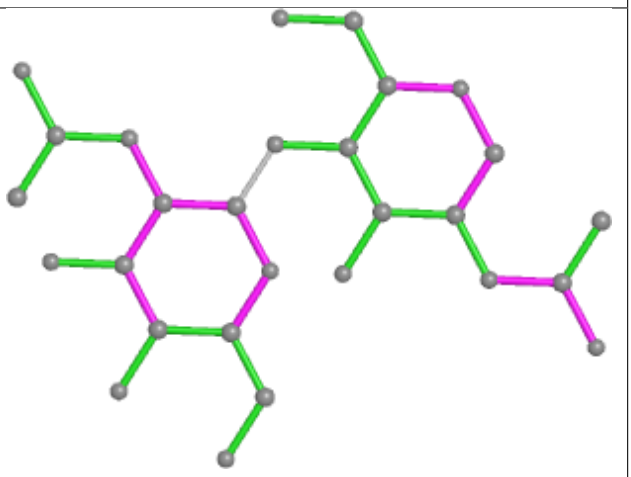


Rings

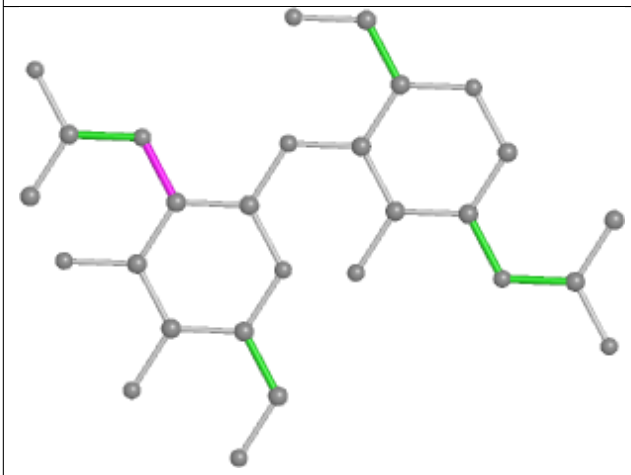
Oligosaccharide Chain G



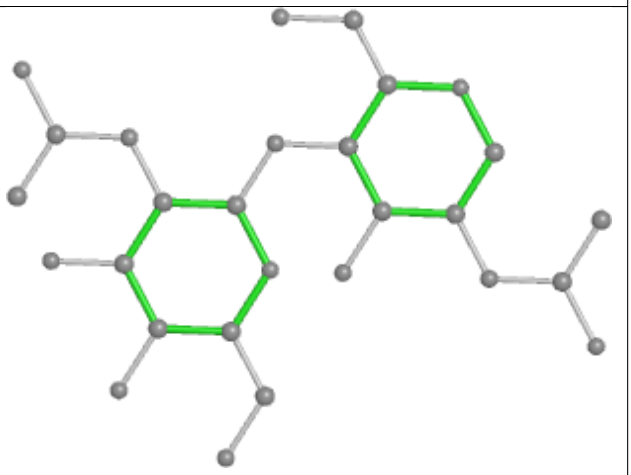
Bond lengths



Bond angles

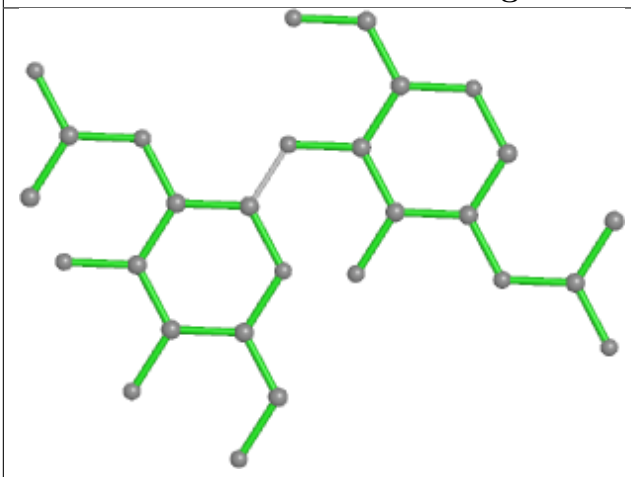


Torsions

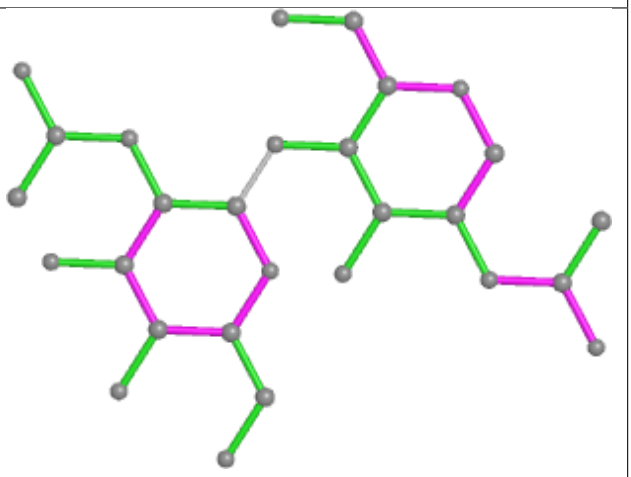


Rings

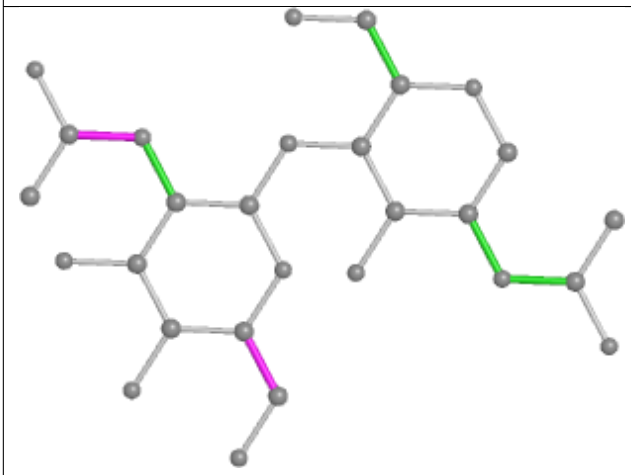
Oligosaccharide Chain I



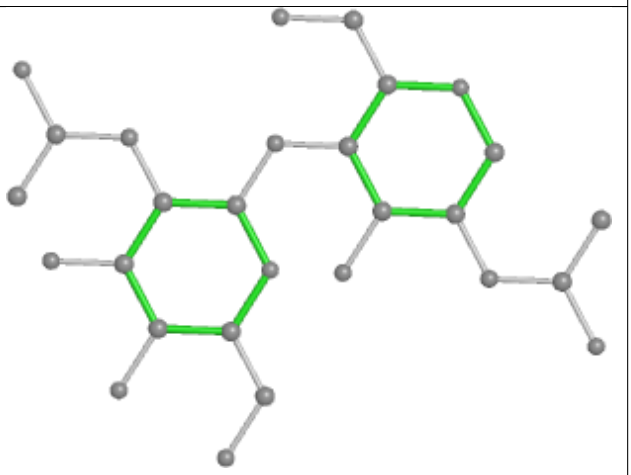
Bond lengths



Bond angles

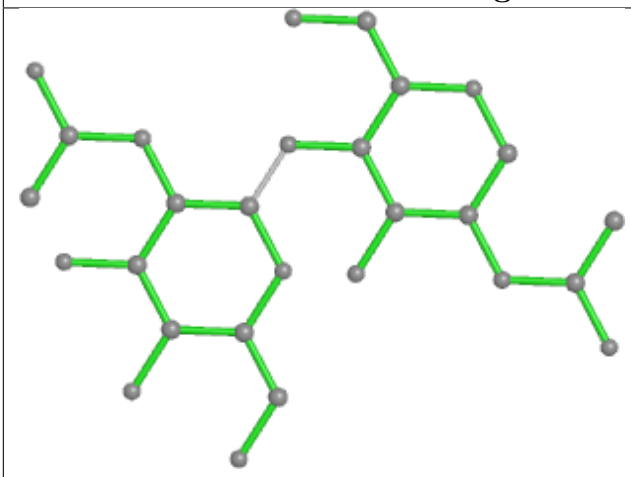


Torsions

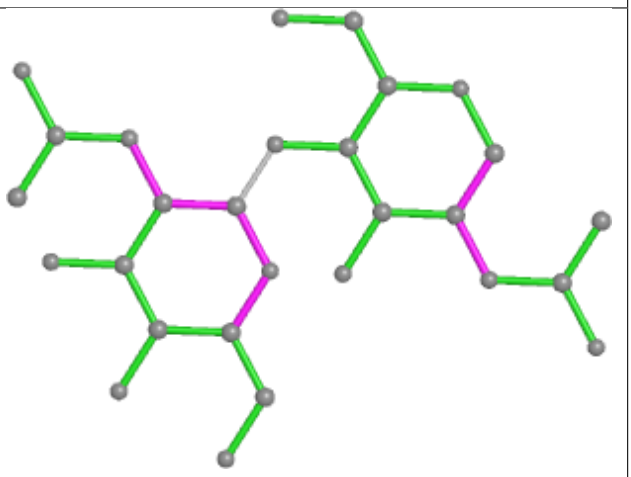


Rings

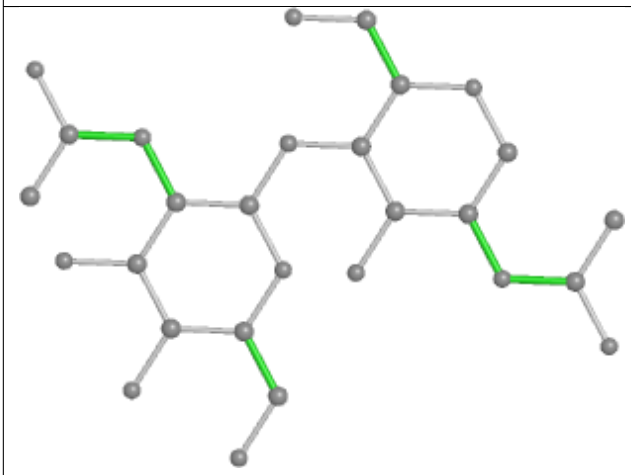
Oligosaccharide Chain K



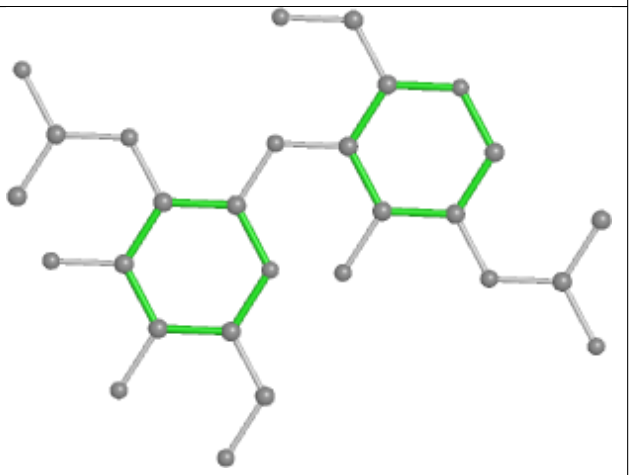
Bond lengths



Bond angles

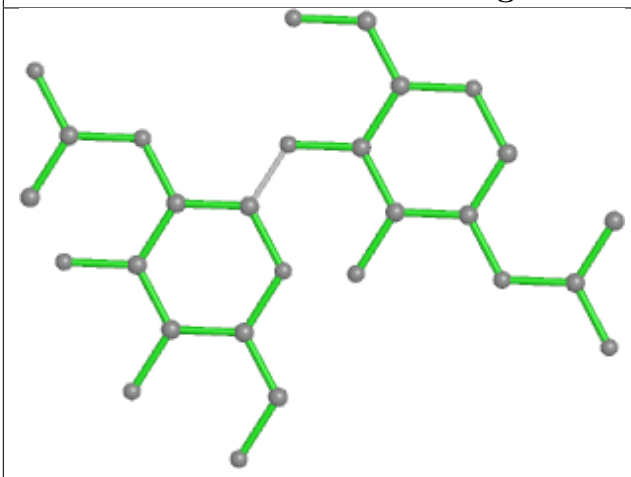


Torsions

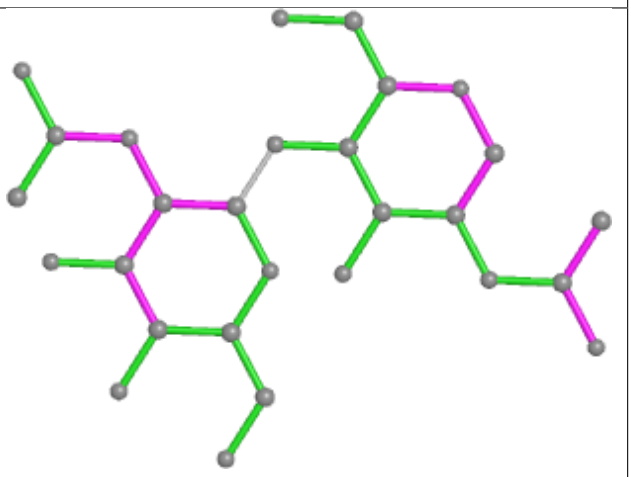


Rings

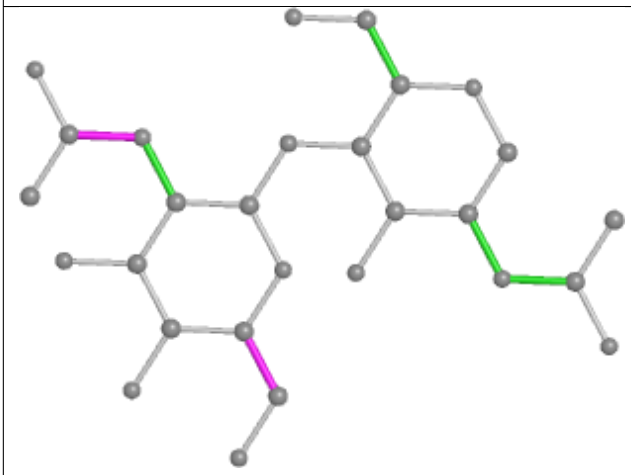
Oligosaccharide Chain L



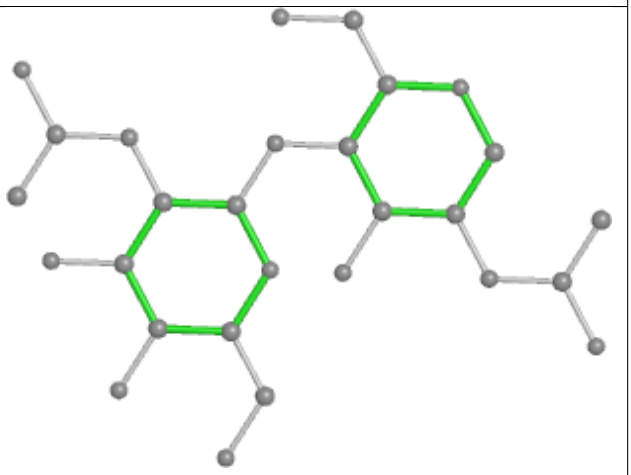
Bond lengths



Bond angles

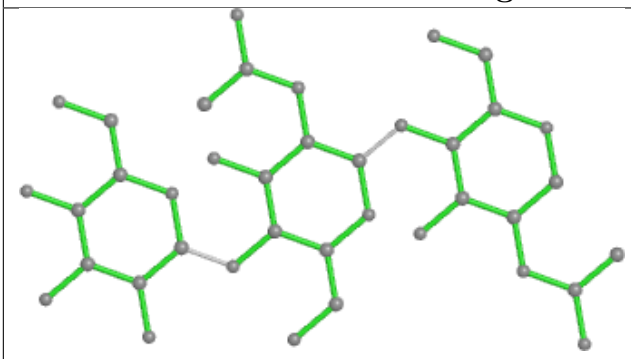


Torsions

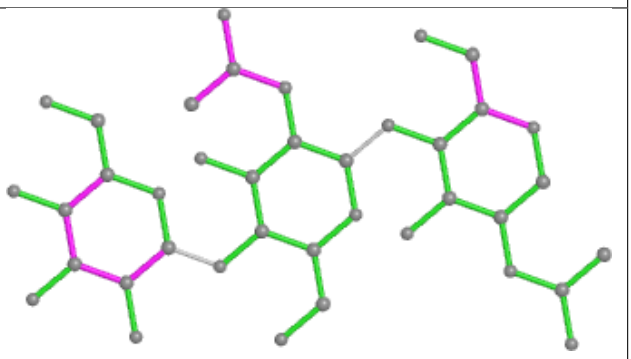


Rings

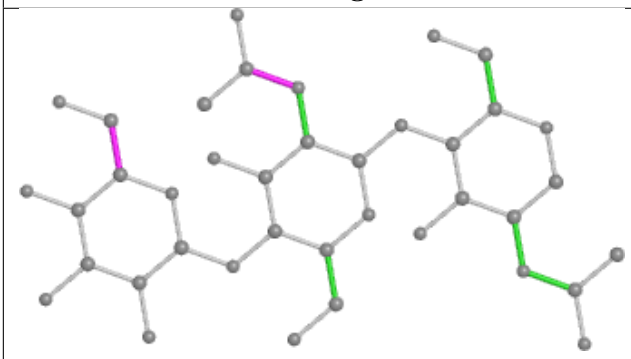
Oligosaccharide Chain E



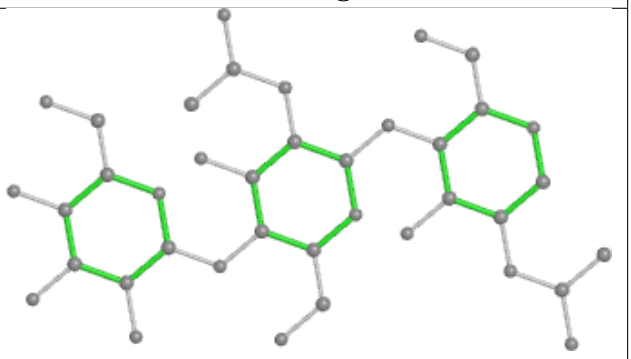
Bond lengths



Bond angles

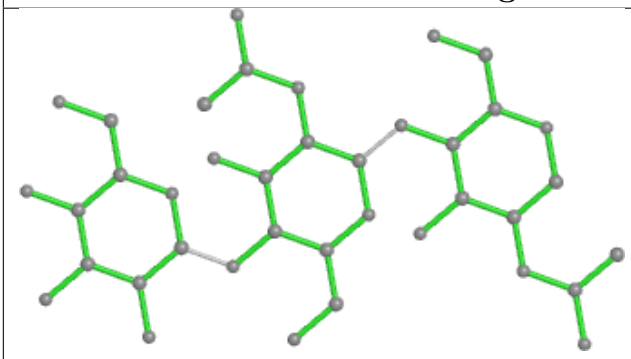


Torsions

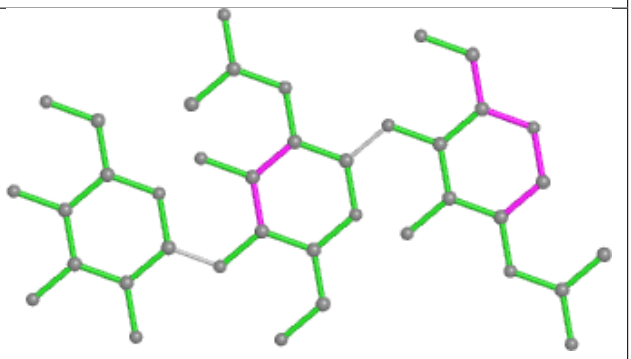


Rings

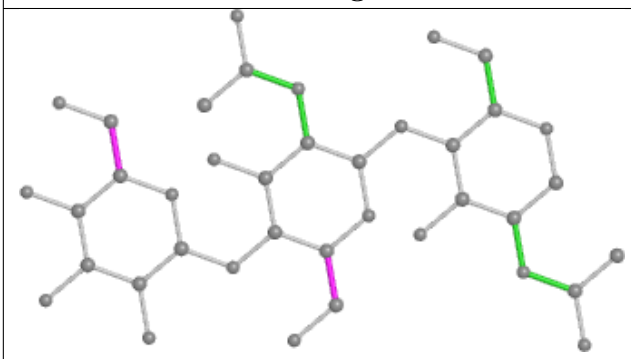
Oligosaccharide Chain J



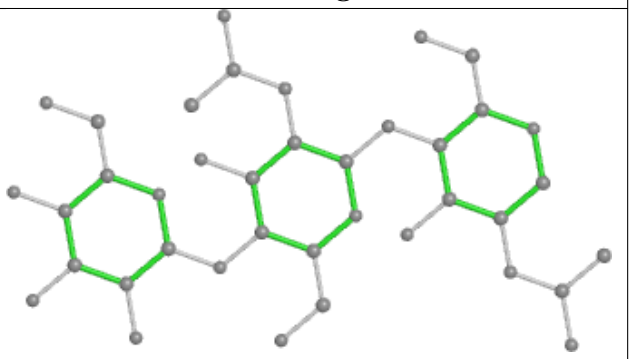
Bond lengths



Bond angles

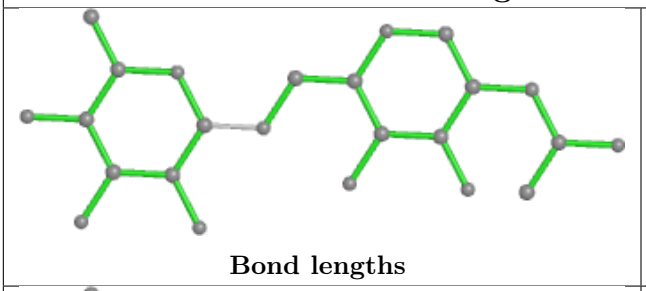


Torsions

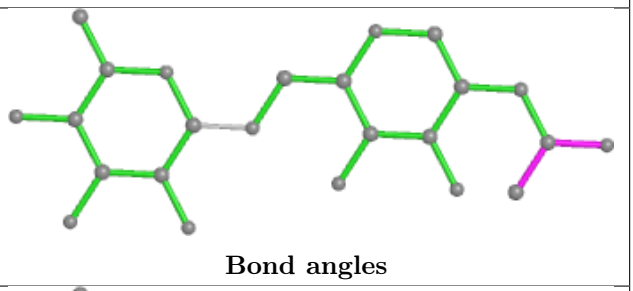


Rings

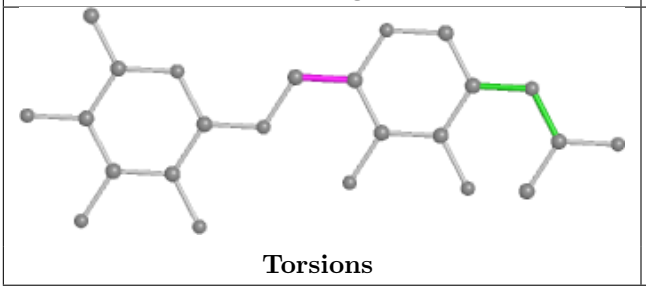
Oligosaccharide Chain H



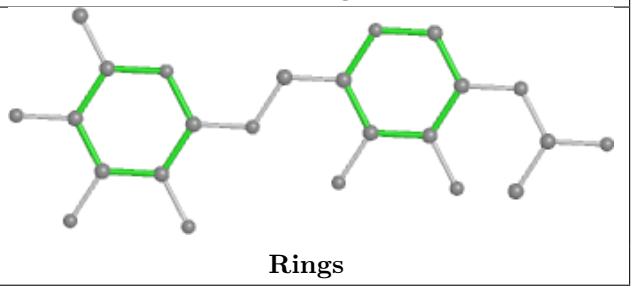
Bond lengths



Bond angles



Torsions



Rings

5.6 Ligand geometry (i)

4 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	$\#\|Z\| > 2$
7	NAG	A	1770	1	14,14,15	0.45	0	17,19,21	1.24	3 (17%)
7	NAG	B	1779	1	14,14,15	0.48	0	17,19,21	1.25	2 (11%)
7	NAG	B	1769	1	14,14,15	0.49	0	17,19,21	1.02	0
7	NAG	A	1780	1	14,14,15	0.53	0	17,19,21	1.82	5 (29%)

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
7	NAG	A	1770	1	-	0/6/23/26	0/1/1/1
7	NAG	B	1779	1	-	4/6/23/26	0/1/1/1
7	NAG	B	1769	1	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
7	NAG	A	1780	1	-	6/6/23/26	0/1/1/1

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
7	A	1780	NAG	C4-C3-C2	-3.98	105.18	111.02
7	B	1779	NAG	C1-O5-C5	3.14	116.45	112.19
7	A	1780	NAG	O5-C5-C6	2.99	111.90	107.20
7	A	1780	NAG	O5-C5-C4	-2.69	104.27	110.83
7	A	1780	NAG	C1-O5-C5	2.61	115.73	112.19
7	A	1770	NAG	C1-O5-C5	2.60	115.72	112.19
7	A	1780	NAG	C3-C4-C5	-2.51	105.76	110.24
7	A	1770	NAG	O5-C5-C6	2.38	110.93	107.20
7	B	1779	NAG	C3-C4-C5	2.33	114.39	110.24
7	A	1770	NAG	O5-C1-C2	-2.00	108.12	111.29

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	A	1780	NAG	C3-C2-N2-C7
7	A	1780	NAG	C8-C7-N2-C2
7	A	1780	NAG	O7-C7-N2-C2
7	B	1779	NAG	O7-C7-N2-C2
7	A	1780	NAG	C4-C5-C6-O6
7	B	1779	NAG	C8-C7-N2-C2
7	A	1780	NAG	O5-C5-C6-O6
7	B	1779	NAG	O5-C5-C6-O6
7	B	1779	NAG	C4-C5-C6-O6
7	A	1780	NAG	C1-C2-N2-C7

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	B	1779	NAG	2	0
7	A	1780	NAG	1	0

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

Unable to reproduce the depositors R factor - this section is therefore empty.

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

Unable to reproduce the depositors R factor - this section is therefore empty.

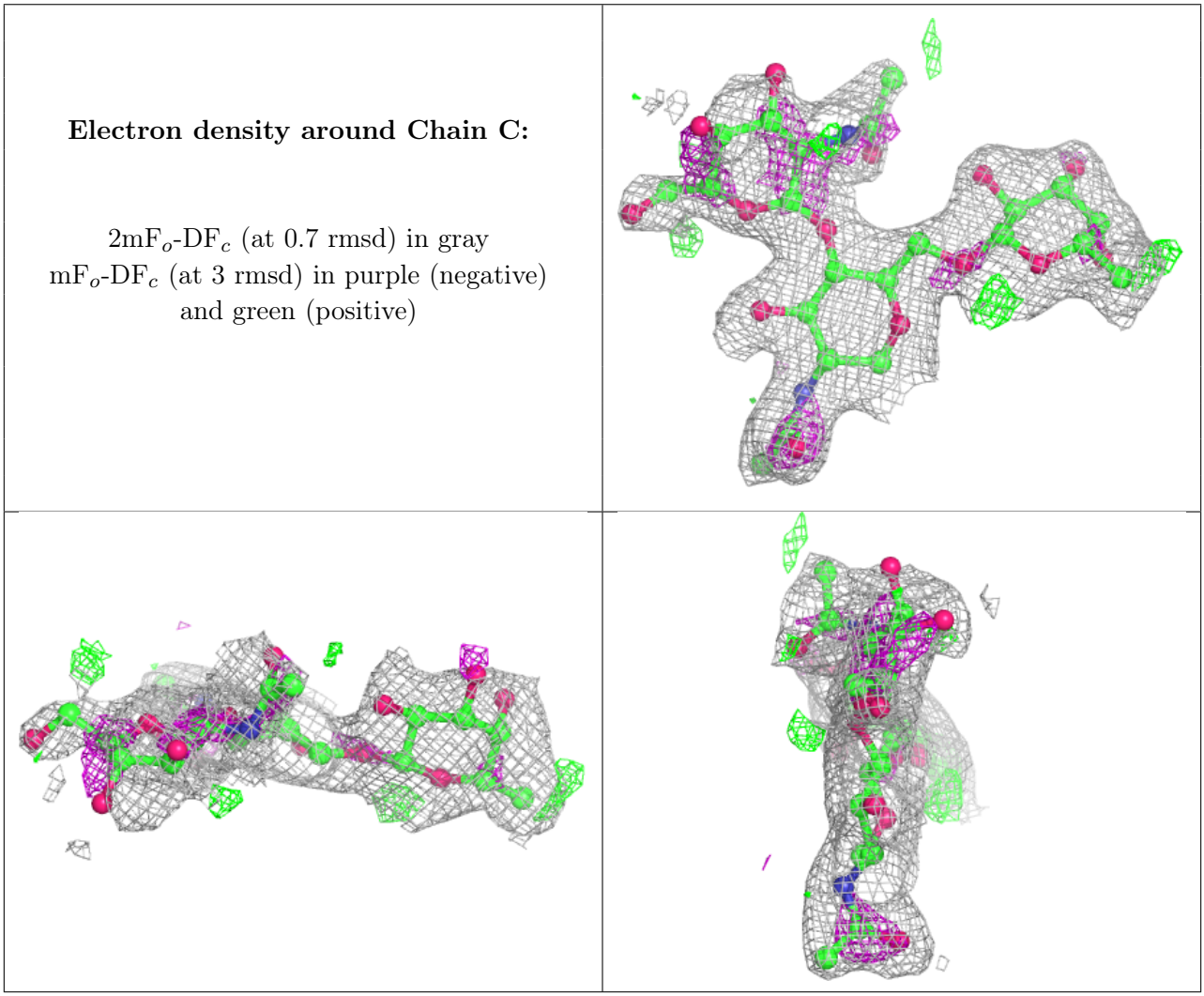
6.3 Carbohydrates

Unable to reproduce the depositors R factor - this section is therefore empty.

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.

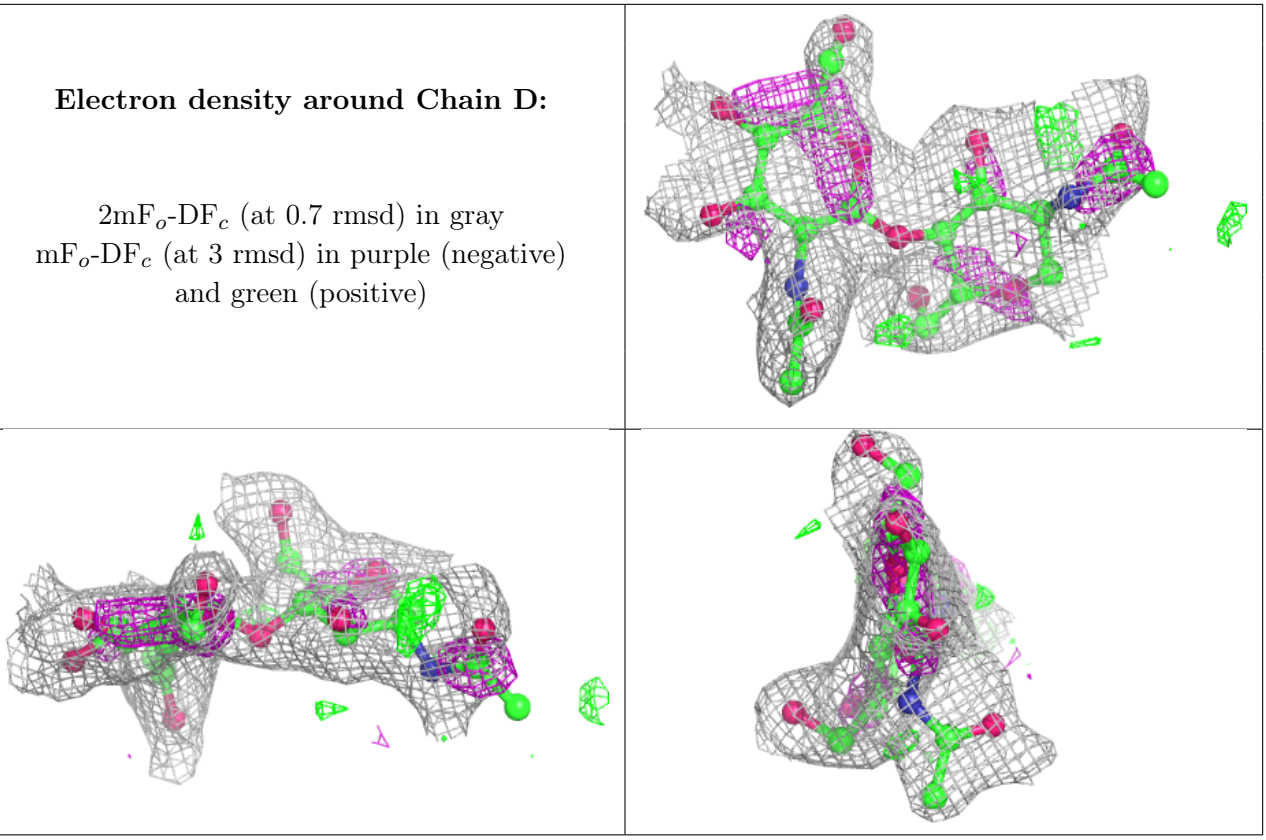
Electron density around Chain C:

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



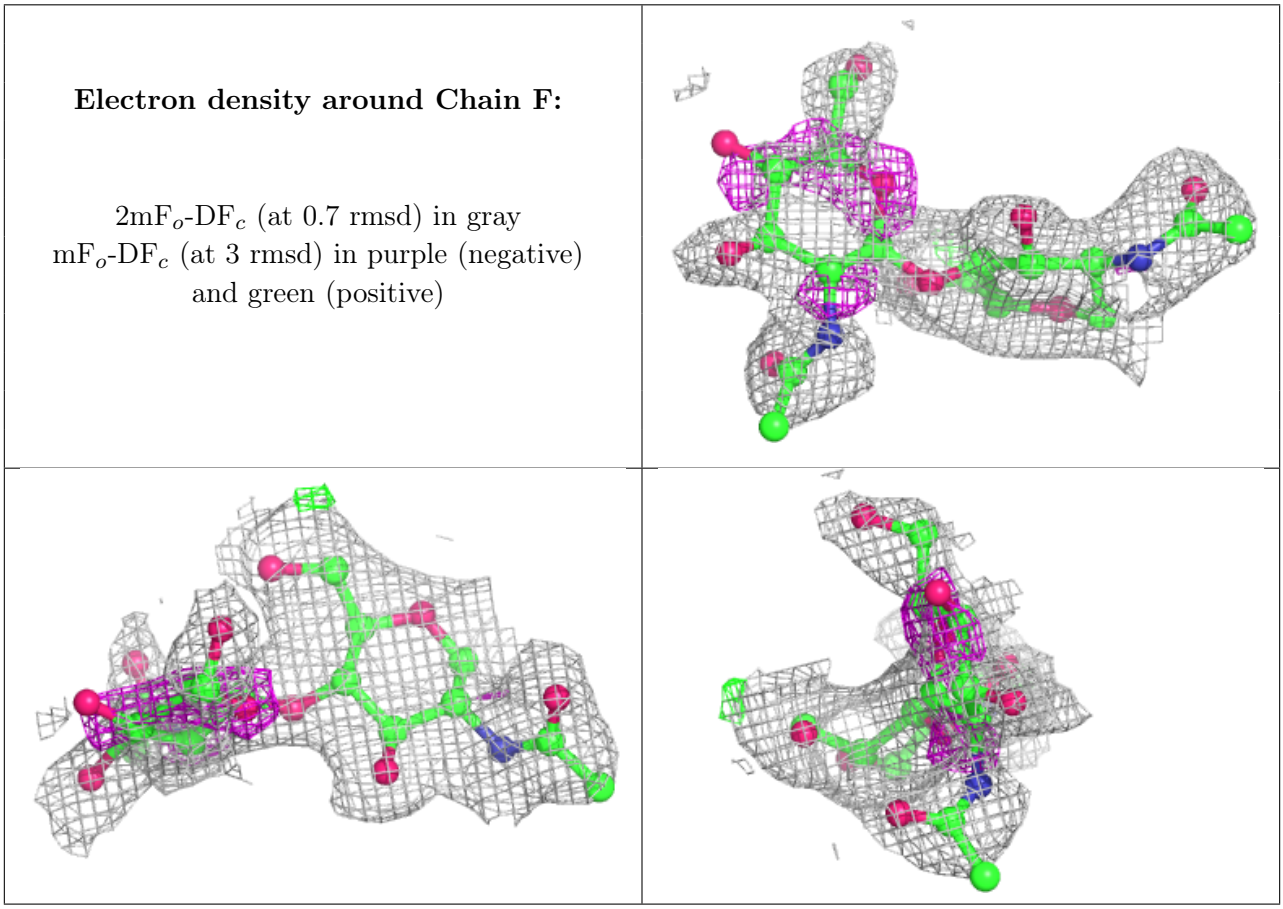
Electron density around Chain D:

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



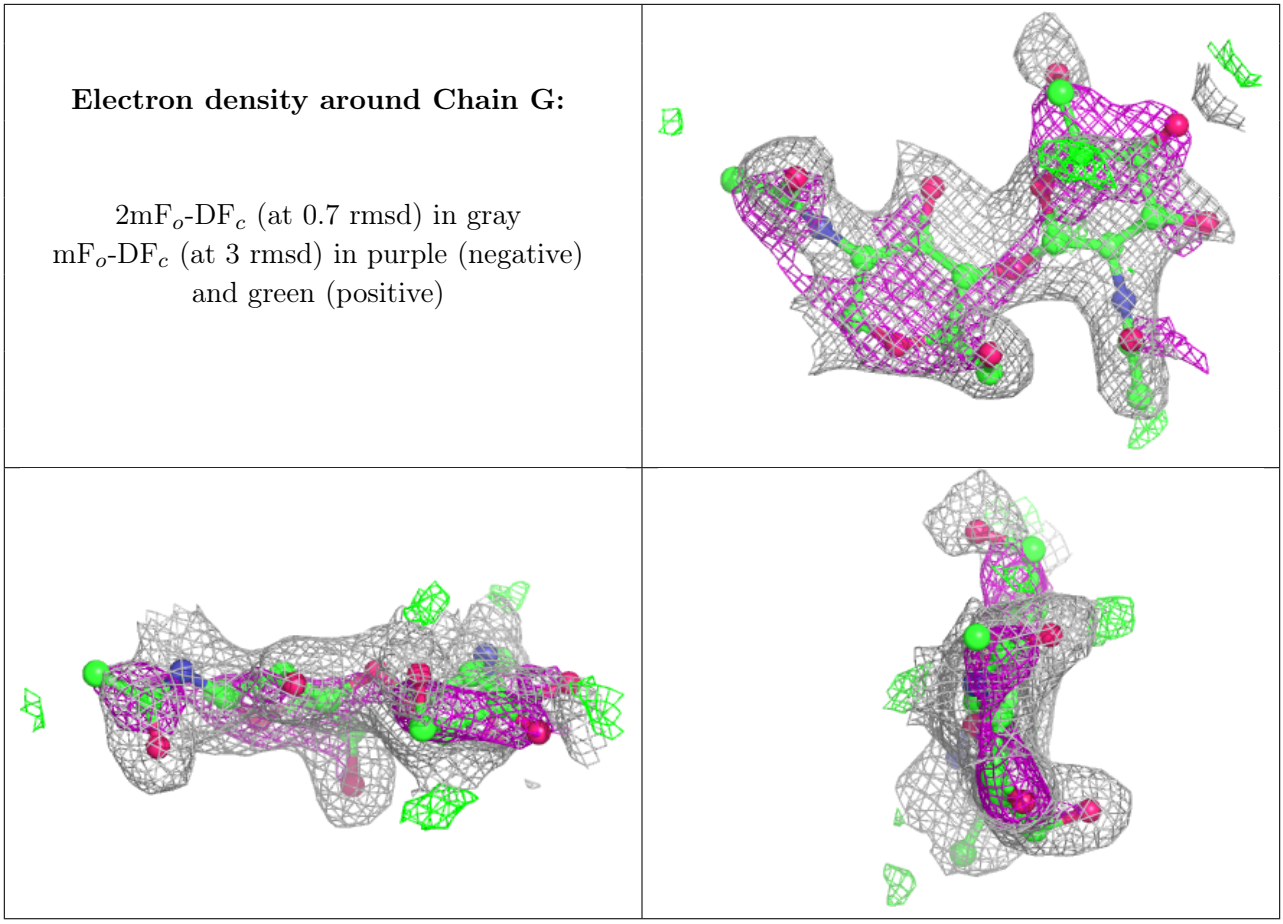
Electron density around Chain F:

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



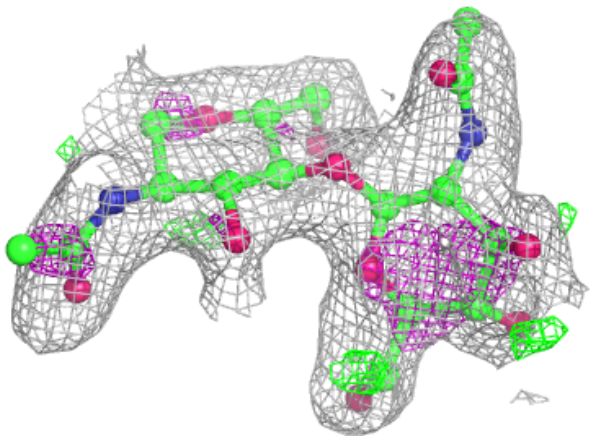
Electron density around Chain G:

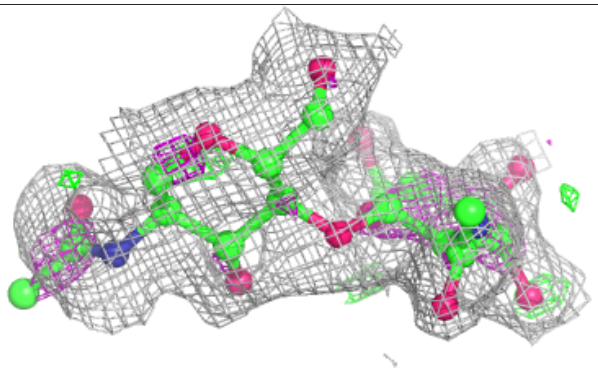
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

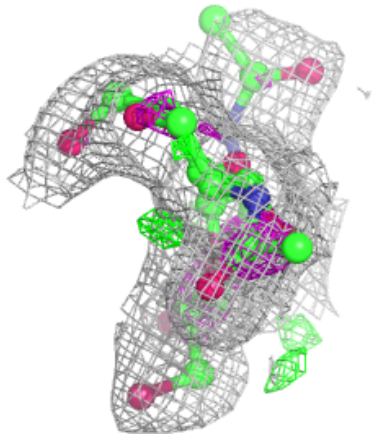


Electron density around Chain I:

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

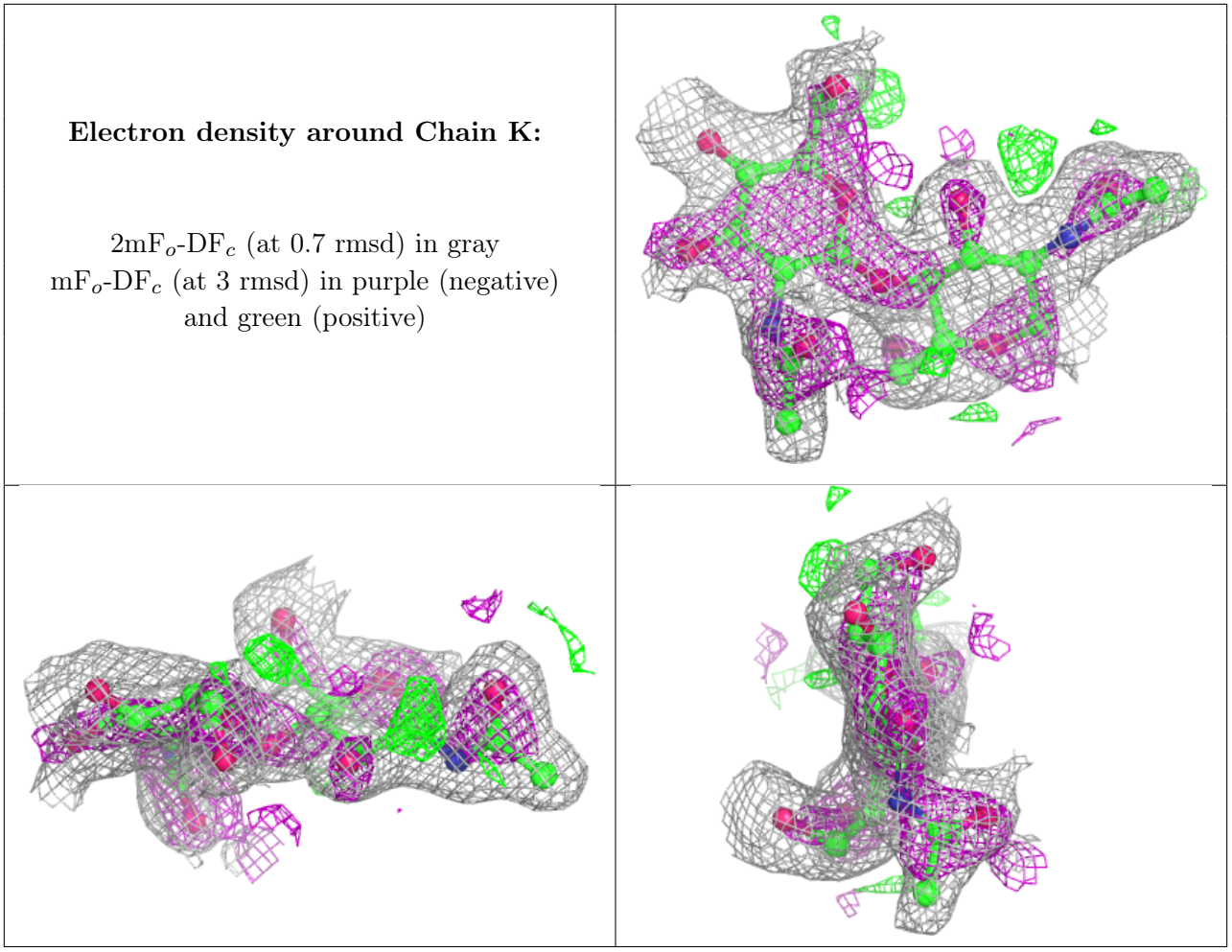






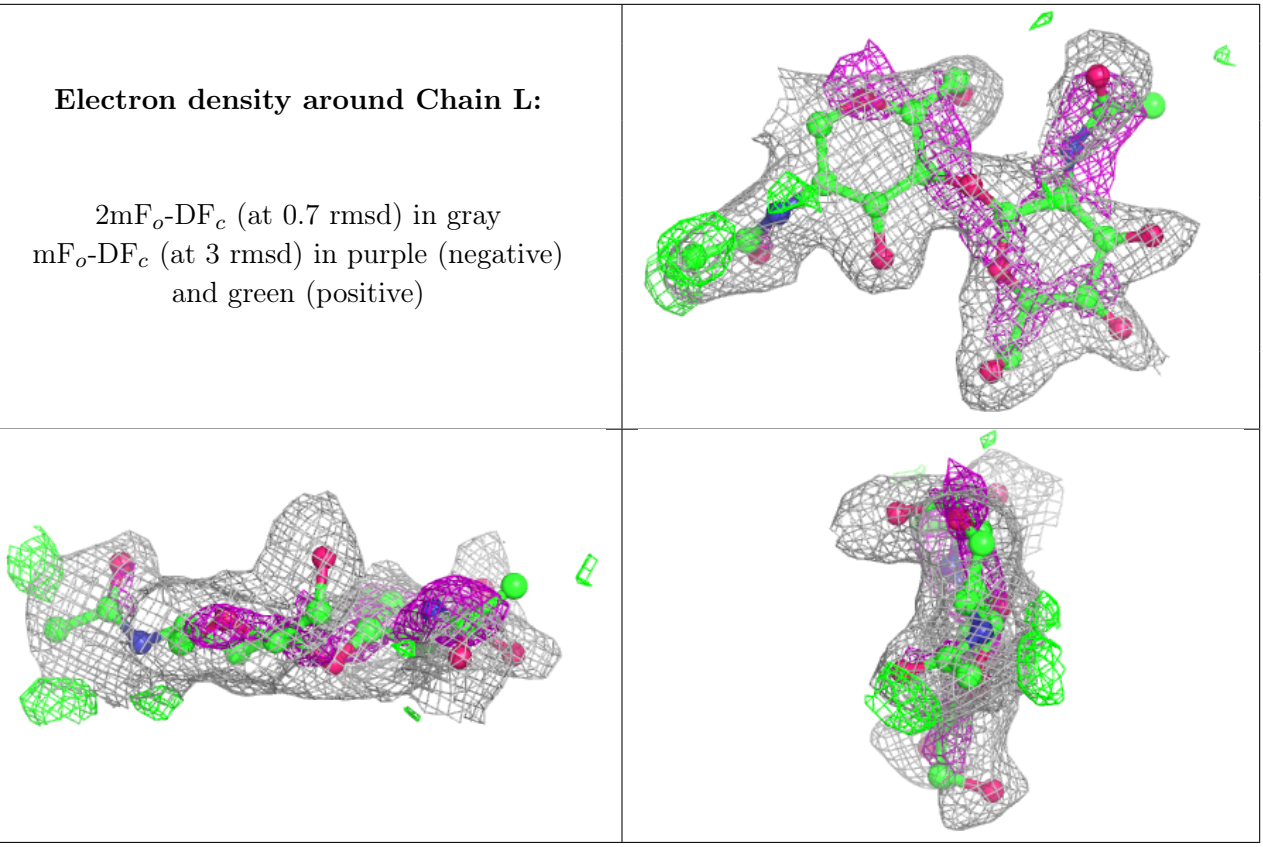
Electron density around Chain K:

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



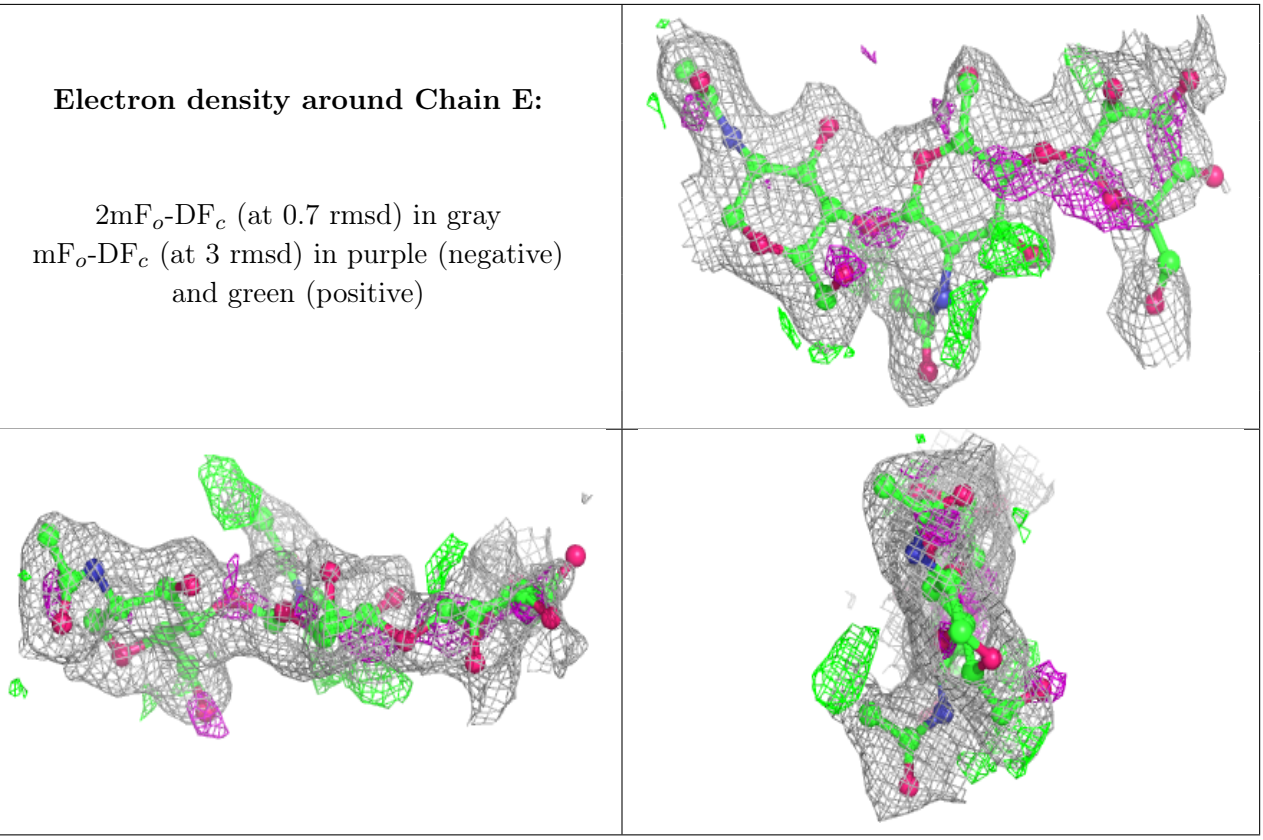
Electron density around Chain L:

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



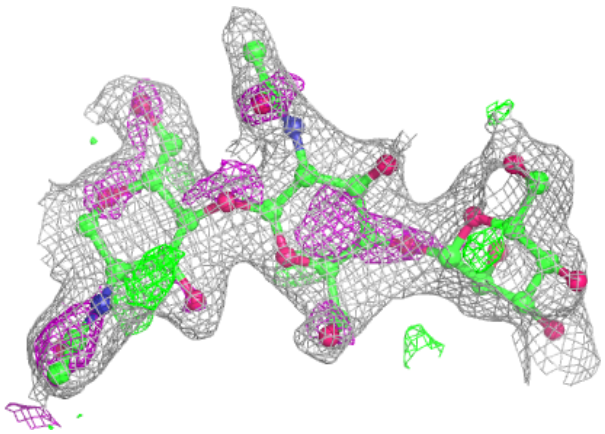
Electron density around Chain E:

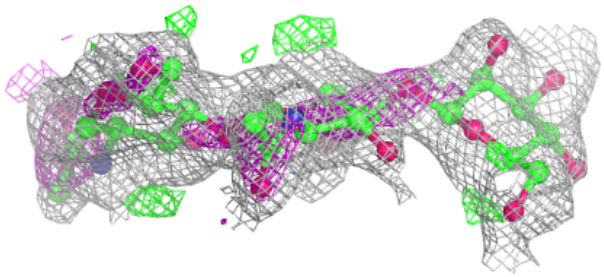
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

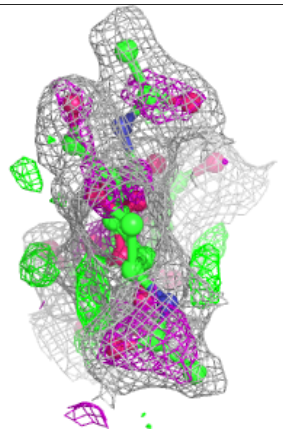


Electron density around Chain J:

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

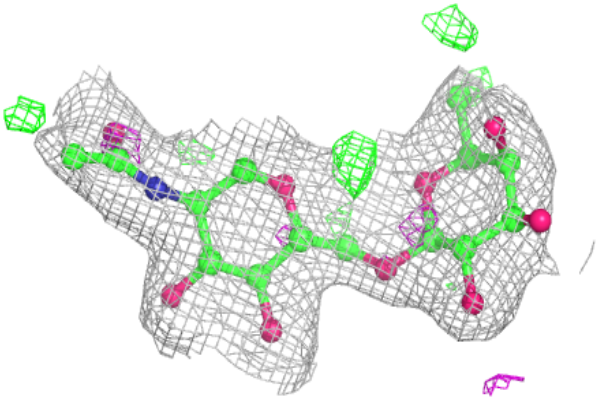


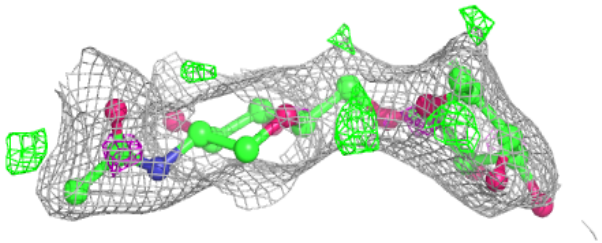


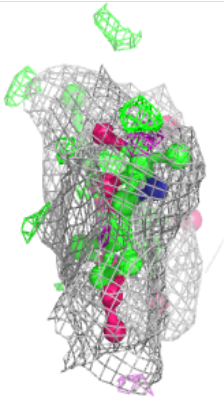


Electron density around Chain H:

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







6.4 Ligands

Unable to reproduce the depositors R factor - this section is therefore empty.

6.5 Other polymers

Unable to reproduce the depositors R factor - this section is therefore empty.