



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:42 AM UTC

PDB ID : 3C45 / pdb_00003c45
Title : Human dipeptidyl peptidase IV/CD26 in complex with a fluoroolefin inhibitor
Authors : Scapin, G.; Edmondson, S.D.; Weber, A.E.
Deposited on : 2008-01-29
Resolution : 2.05 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

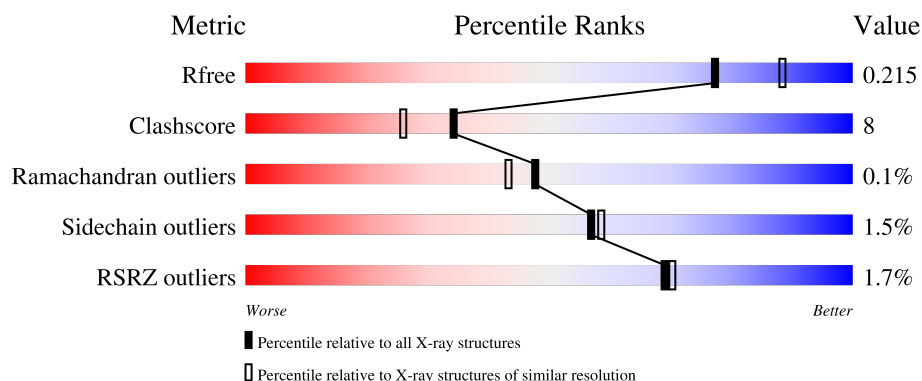
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.05 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	2260 (2.04-2.04)
Clashscore	190562	2333 (2.04-2.04)
Ramachandran outliers	187476	2318 (2.04-2.04)
Sidechain outliers	187428	2318 (2.04-2.04)
RSRZ outliers	180081	2260 (2.04-2.04)

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 $\leq 5\%$. 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	728	2% 82% 17% .
1	B	728	2% 82% 16% .
2	C	2	50% 50%
2	D	2	100%
2	E	2	100%

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Mol	Chain	Length	Quality of chain
2	F	2	 100%
2	G	2	 100%
2	H	2	 50% 50%
2	I	2	 100%
2	J	2	 50% 50%
2	K	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	NAG	C	2	X	-	-	-
2	NAG	D	1	X	-	-	-
2	NAG	F	2	X	-	-	-

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 13342 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 4 soluble form.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	728	Total	C	N	O	S	0	0	0
			5965	3828	982	1129	26			
1	B	728	Total	C	N	O	S	0	0	0
			5965	3828	982	1129	26			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	39	THR	SER	engineered mutation	UNP P27487
B	39	THR	SER	engineered mutation	UNP P27487

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



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

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

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

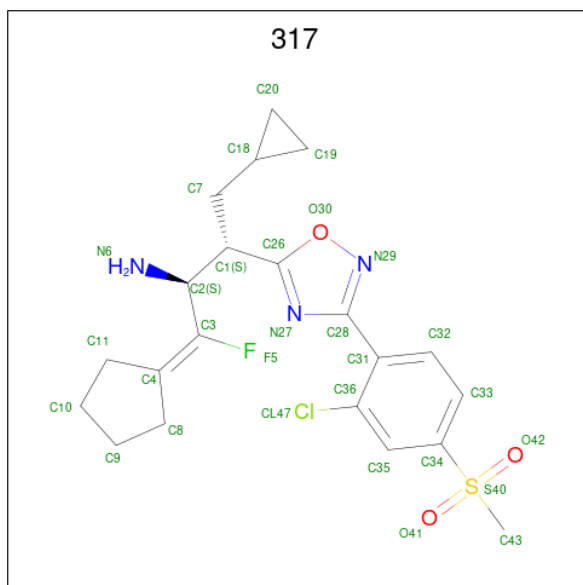


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	B	1	Total	C	N	O	0	0
			14	8	1	5		
3	B	1	Total	C	N	O	0	0
			14	8	1	5		
3	B	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 4 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total Na 1 1	0	0

- Molecule 5 is (2S,3S)-3-{3-[2-chloro-4-(methylsulfonyl)phenyl]-1,2,4-oxadiazol-5-yl}-1-cyclopentylidene-4-cyclopropyl-1-fluorobutan-2-amine (CCD ID: 317) (formula: C₂₁H₂₅ClFN₃O₃S).



Mol	Chain	Residues	Atoms							ZeroOcc	AltConf
5	A	1	Total	C	Cl	F	N	O	S	0	0
			30	21	1	1	3	3	1		
5	B	1	Total	C	Cl	F	N	O	S	0	0
			30	21	1	1	3	3	1		

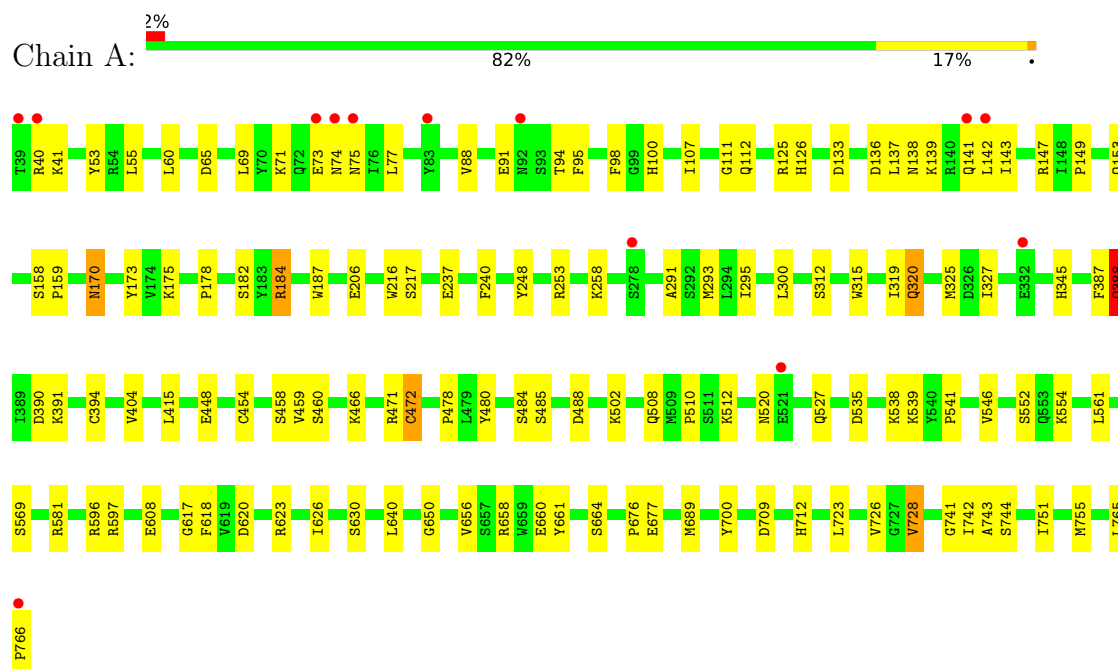
- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	504	Total O 504 504	0	0
6	B	511	Total O 511 511	0	0

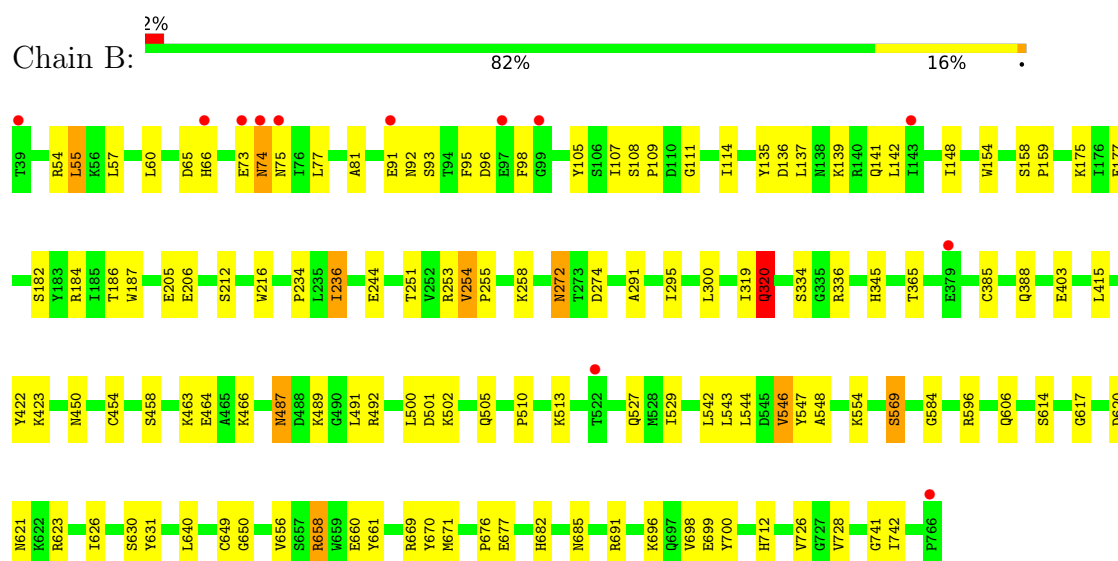
3 Residue-property plots

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: Dipeptidyl peptidase 4 soluble form



• Molecule 1: Dipeptidyl peptidase 4 soluble form



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

Chain C:  50% 50%

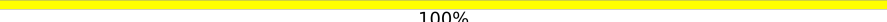
MAG1
MAG2

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

Chain D:  100%

MAG1
MAG2

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

Chain E:  100%

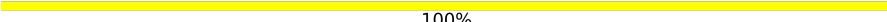
MAG1
MAG2

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

Chain F:  100%

MAG1
MAG2

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

Chain G:  100%

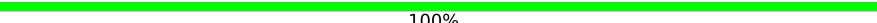
MAG1
MAG2

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

Chain H:  50% 50%

MAG1
MAG2

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

Chain I:  100%



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

Chain J:  50%  50%



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

Chain K:  50%  50%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	117.89Å 125.91Å 136.75Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.05 30.00 – 2.05	Depositor EDS
% Data completeness (in resolution range)	99.7 (30.00-2.05) 99.7 (30.00-2.05)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.99 (at 2.04Å)	Xtriage
Refinement program	CNX	Depositor
R, R_{free}	0.191 , 0.218 0.188 , 0.215	Depositor DCC
R_{free} test set	6493 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	27.2	Xtriage
Anisotropy	0.426	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 40.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	13342	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.74% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NAG, NA, 317

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	0.49	0/6137	0.93	17/8346 (0.2%)
1	B	0.48	0/6137	0.94	24/8346 (0.3%)
All	All	0.48	0/12274	0.94	41/16692 (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	A	0	1
1	B	0	1
All	All	0	2

There are no bond length outliers.

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	388	GLN	N-CA-C	-8.20	96.32	109.76
1	A	388	GLN	N-CA-C	-7.92	96.78	109.76
1	A	458	SER	N-CA-C	-7.53	97.92	109.14
1	A	300	LEU	N-CA-C	-7.45	97.09	109.24
1	A	206	GLU	N-CA-C	7.17	122.20	113.38
1	B	656	VAL	N-CA-C	-7.15	99.68	109.55
1	A	656	VAL	N-CA-C	-6.98	98.09	109.12
1	A	485	SER	N-CA-C	6.86	118.84	111.36
1	B	458	SER	N-CA-C	-6.82	98.95	109.24
1	B	548	ALA	N-CA-C	6.68	121.38	113.23
1	B	463	LYS	N-CA-C	6.52	119.22	111.33
1	B	529	ILE	N-CA-C	-6.51	98.27	108.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	319	ILE	N-CA-C	-6.43	97.31	107.15
1	B	617	GLY	N-CA-C	6.36	122.64	115.08
1	B	300	LEU	N-CA-C	-6.35	98.38	108.73
1	B	148	ILE	N-CA-C	-6.34	103.12	109.02
1	A	664	SER	N-CA-C	6.25	117.76	111.07
1	B	569	SER	N-CA-C	6.25	117.89	111.14
1	B	177	GLU	CA-C-N	6.16	126.34	119.32
1	B	177	GLU	C-N-CA	6.16	126.34	119.32
1	B	546	VAL	N-CA-C	6.15	118.32	108.85
1	B	319	ILE	N-CA-C	-6.06	97.88	107.15
1	B	454	CYS	N-CA-C	6.02	118.55	108.02
1	B	320	GLN	N-CA-C	5.99	123.56	110.80
1	A	546	VAL	N-CA-C	5.96	118.57	108.86
1	A	744	SER	N-CA-C	-5.82	102.44	110.35
1	A	454	CYS	N-CA-C	5.80	118.34	108.02
1	A	320	GLN	N-CA-C	5.79	123.14	110.80
1	A	158	SER	N-CA-C	-5.72	101.97	110.20
1	B	403	GLU	N-CA-C	5.68	118.47	109.50
1	B	205	GLU	N-CA-C	5.67	117.54	111.36
1	B	206	GLU	N-CA-C	5.50	120.87	113.72
1	B	584	GLY	N-CA-C	5.37	121.06	112.58
1	A	541	PRO	N-CA-C	-5.35	103.39	111.41
1	B	698	VAL	N-CA-C	5.30	116.34	108.65
1	B	631	TYR	N-CA-C	-5.29	105.62	111.71
1	A	404	VAL	N-CA-C	-5.17	102.06	108.89
1	B	236	ILE	N-CA-C	-5.09	101.13	108.87
1	A	743	ALA	N-CA-C	5.05	119.53	113.17
1	A	240	PHE	N-CA-C	-5.04	100.08	108.34
1	B	365	THR	N-CA-C	-5.03	103.90	110.53

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	700	TYR	Sidechain
1	B	700	TYR	Sidechain

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen

atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5965	0	5676	94	0
1	B	5965	0	5679	90	0
2	C	28	0	25	0	0
2	D	28	0	25	3	0
2	E	28	0	25	0	0
2	F	28	0	25	2	0
2	G	28	0	25	1	0
2	H	28	0	25	0	0
2	I	28	0	25	0	0
2	J	28	0	25	0	0
2	K	28	0	25	0	0
3	A	42	0	39	4	0
3	B	42	0	39	1	0
4	A	1	0	0	0	0
5	A	30	0	25	4	0
5	B	30	0	25	2	0
6	A	504	0	0	1	0
6	B	511	0	0	4	0
All	All	13342	0	11708	183	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 8.

All (183) 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:149:PRO:HA	2:D:1:NAG:H82	1.42	0.99
1:A:581:ARG:CZ	3:A:803:NAG:H62	2.07	0.85
1:A:75:ASN:HD21	3:A:801:NAG:HN2	1.29	0.79
1:B:184:ARG:HD3	1:B:186:THR:O	1.84	0.78
1:A:581:ARG:NH1	3:A:803:NAG:H62	1.98	0.77
1:A:554:LYS:HG2	5:A:805:317:H43	1.67	0.77
1:A:258:LYS:NZ	1:A:712:HIS:HD2	1.82	0.76
1:A:312:SER:HB2	1:A:325:MET:HE3	1.66	0.75
1:B:258:LYS:NZ	1:B:712:HIS:HD2	1.84	0.75
1:A:596:ARG:O	1:A:597:ARG:HD2	1.88	0.73
1:B:726:VAL:HG23	1:B:728:VAL:HG23	1.70	0.73
1:B:450:ASN:HB2	6:B:1230:HOH:O	1.89	0.73
1:B:258:LYS:HZ1	1:B:712:HIS:HD2	1.38	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:272:ASN:C	1:B:272:ASN:HD22	1.97	0.72
1:B:184:ARG:HD2	1:B:187:TRP:CE2	2.27	0.70
1:B:696:LYS:HG3	1:B:728:VAL:HG22	1.72	0.70
1:A:676:PRO:HG2	1:A:677:GLU:OE2	1.92	0.70
1:B:320:GLN:OE1	1:B:669:ARG:HD3	1.92	0.69
1:B:676:PRO:HG2	1:B:677:GLU:OE2	1.92	0.69
1:A:640:LEU:HD11	1:A:650:GLY:HA3	1.72	0.69
1:B:81:ALA:O	1:B:492:ARG:NH2	2.25	0.68
1:B:272:ASN:HD21	1:B:274:ASP:HB2	1.57	0.68
1:A:184:ARG:NH1	1:A:187:TRP:HA	2.08	0.68
1:B:75:ASN:HB3	1:B:92:ASN:N	2.08	0.68
1:A:471:ARG:HG3	1:A:480:TYR:CE2	2.28	0.68
1:A:139:LYS:HG3	1:A:141:GLN:HB3	1.75	0.68
1:A:658:ARG:HG2	1:A:661:TYR:CE2	2.29	0.67
1:B:691:ARG:HD2	6:B:1096:HOH:O	1.94	0.67
1:A:153:GLN:HE22	1:A:170:ASN:ND2	1.93	0.66
1:A:253:ARG:HH21	1:B:253:ARG:NH1	1.94	0.66
1:A:175:LYS:HG3	1:A:182:SER:HB3	1.78	0.64
1:A:184:ARG:HH11	1:A:187:TRP:HA	1.61	0.64
1:A:77:LEU:HD23	1:A:88:VAL:HA	1.79	0.63
1:A:173:TYR:CE2	1:A:184:ARG:HG2	2.34	0.63
1:A:88:VAL:HG11	1:A:91:GLU:OE2	1.99	0.63
1:B:114:ILE:HG23	1:B:135:TYR:HB3	1.81	0.63
2:F:1:NAG:H62	2:F:2:NAG:O5	1.98	0.62
1:B:554:LYS:HG2	5:B:804:317:H43	1.81	0.62
1:A:98:PHE:CE2	1:A:100:HIS:HB2	2.35	0.62
1:B:614:SER:HB2	1:B:621:ASN:OD1	2.00	0.62
1:A:258:LYS:HZ1	1:A:712:HIS:HD2	1.45	0.61
1:B:65:ASP:OD2	1:B:466:LYS:HB2	2.00	0.60
1:B:658:ARG:HG2	1:B:661:TYR:CE2	2.38	0.59
1:A:630:SER:OG	5:A:805:317:H8A	2.03	0.58
1:A:620:ASP:OD2	1:A:623:ARG:HD3	2.03	0.58
1:B:640:LEU:HD11	1:B:650:GLY:HA3	1.84	0.58
1:B:741:GLY:O	1:B:742:ILE:C	2.46	0.58
1:B:60:LEU:HD12	1:B:60:LEU:C	2.28	0.58
1:A:133:ASP:HB3	1:A:142:LEU:HD11	1.85	0.57
1:B:139:LYS:HB3	1:B:141:GLN:HE21	1.68	0.57
1:B:544:LEU:HD21	1:B:606:GLN:HG3	1.84	0.57
1:A:723:LEU:HB3	1:A:728:VAL:HG13	1.87	0.57
1:A:170:ASN:HD22	1:A:170:ASN:N	2.03	0.57
1:B:73:GLU:OE1	3:B:801:NAG:H83	2.03	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:248:TYR:CZ	1:B:234:PRO:HB2	2.39	0.57
1:B:510:PRO:HD3	1:B:569:SER:HB2	1.87	0.57
1:B:251:THR:HG21	1:B:253:ARG:NH1	2.20	0.56
1:B:75:ASN:HB3	1:B:92:ASN:H	1.70	0.56
1:A:65:ASP:OD2	1:A:466:LYS:HB2	2.06	0.55
1:A:136:ASP:HB3	1:A:139:LYS:HG2	1.88	0.55
1:B:107:ILE:HD12	1:B:107:ILE:N	2.22	0.55
1:B:658:ARG:HD3	1:B:660:GLU:HB2	1.88	0.55
1:A:159:PRO:HD3	1:A:216:TRP:CB	2.37	0.55
1:B:415:LEU:C	1:B:415:LEU:HD23	2.31	0.55
1:A:312:SER:CB	1:A:325:MET:HE3	2.37	0.54
1:B:154:TRP:CE2	1:B:212:SER:HB3	2.41	0.54
1:B:620:ASP:OD2	1:B:623:ARG:HD3	2.07	0.54
1:A:111:GLY:O	1:A:137:LEU:HD12	2.07	0.54
1:B:109:PRO:HG2	1:B:158:SER:O	2.07	0.54
1:B:272:ASN:ND2	1:B:274:ASP:H	2.06	0.54
1:A:95:PHE:HB3	1:A:98:PHE:HB2	1.90	0.54
1:B:91:GLU:HA	1:B:91:GLU:OE1	2.08	0.54
1:B:345:HIS:HD2	6:B:959:HOH:O	1.91	0.53
1:A:677:GLU:H	1:A:677:GLU:CD	2.17	0.52
1:B:175:LYS:CG	1:B:182:SER:HB3	2.39	0.52
1:A:159:PRO:HG3	1:A:217:SER:O	2.09	0.51
1:B:272:ASN:HD22	1:B:274:ASP:H	1.58	0.51
1:A:741:GLY:O	1:A:742:ILE:C	2.54	0.51
2:F:1:NAG:H62	2:F:2:NAG:C1	2.41	0.51
1:B:542:LEU:C	1:B:542:LEU:HD23	2.36	0.51
1:A:325:MET:HE2	1:A:327:ILE:CG1	2.41	0.51
1:A:41:LYS:HE2	1:A:53:TYR:OH	2.11	0.51
1:B:74:ASN:HA	6:B:1304:HOH:O	2.10	0.50
1:A:159:PRO:HD3	1:A:216:TRP:HB3	1.93	0.50
1:B:77:LEU:HD12	1:B:77:LEU:N	2.26	0.50
1:B:114:ILE:CG2	1:B:135:TYR:HB3	2.40	0.50
1:B:513:LYS:O	1:B:527:GLN:HA	2.12	0.50
1:A:60:LEU:HD12	1:A:60:LEU:C	2.37	0.50
1:B:107:ILE:HG22	1:B:108:SER:O	2.11	0.50
1:A:535:ASP:HB3	1:A:538:LYS:HD2	1.92	0.50
1:B:272:ASN:C	1:B:272:ASN:ND2	2.69	0.50
1:A:147:ARG:HB2	6:A:1168:HOH:O	2.12	0.49
1:A:112:GLN:HG2	1:A:138:ASN:HD21	1.76	0.49
1:B:502:LYS:HD3	1:B:502:LYS:C	2.37	0.49
1:A:253:ARG:NH2	1:B:253:ARG:NH1	2.60	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:658:ARG:O	1:B:658:ARG:HG3	2.12	0.49
1:B:105:TYR:HB2	1:B:114:ILE:HD11	1.95	0.48
1:B:630:SER:OG	5:B:804:317:H8A	2.13	0.48
1:A:69:LEU:HD13	1:A:107:ILE:HD12	1.95	0.48
1:B:491:LEU:O	1:B:492:ARG:HB3	2.12	0.48
1:B:422:TYR:CE2	1:B:423:LYS:HE2	2.48	0.48
1:B:658:ARG:HG2	1:B:661:TYR:CD2	2.48	0.48
1:A:726:VAL:HG23	1:A:728:VAL:HG12	1.96	0.48
1:A:258:LYS:NZ	1:A:712:HIS:CD2	2.72	0.48
1:A:71:LYS:HE3	1:A:74:ASN:HA	1.96	0.47
1:B:139:LYS:O	1:B:141:GLN:HG3	2.14	0.47
1:A:415:LEU:HD23	1:A:415:LEU:C	2.39	0.47
1:A:258:LYS:HZ3	1:A:712:HIS:HD2	1.59	0.47
1:B:139:LYS:HE2	1:B:141:GLN:NE2	2.30	0.47
1:B:175:LYS:HG3	1:B:182:SER:HB3	1.96	0.47
1:A:91:GLU:HB2	1:A:94:THR:OG1	2.15	0.47
1:A:689:MET:HE3	1:B:244:GLU:HG3	1.97	0.47
1:A:237:GLU:HG2	1:A:253:ARG:HG2	1.97	0.47
1:B:65:ASP:CG	1:B:464:GLU:HB2	2.40	0.47
1:B:98:PHE:HE1	1:B:142:LEU:HD21	1.80	0.47
1:B:542:LEU:HD23	1:B:543:LEU:N	2.29	0.47
1:B:93:SER:HA	1:B:96:ASP:OD1	2.15	0.47
1:A:258:LYS:HZ3	1:A:712:HIS:CD2	2.31	0.46
1:B:487:ASN:C	1:B:487:ASN:HD22	2.22	0.46
2:D:1:NAG:H4	2:D:2:NAG:H2	1.68	0.46
1:B:254:VAL:HA	1:B:255:PRO:HD3	1.83	0.46
1:A:40:ARG:HG3	1:A:508:GLN:HG3	1.98	0.46
1:A:510:PRO:HD3	1:A:569:SER:HB2	1.97	0.45
1:B:236:ILE:CG2	1:B:254:VAL:HG13	2.46	0.45
1:A:658:ARG:HG2	1:A:661:TYR:CD2	2.52	0.45
1:A:325:MET:HE2	1:A:327:ILE:HG12	1.98	0.44
1:A:539:LYS:HD3	1:A:617:GLY:O	2.18	0.44
1:A:596:ARG:C	1:A:597:ARG:HD2	2.42	0.44
1:B:649:CYS:HB3	1:B:699:GLU:HB2	1.99	0.44
1:A:98:PHE:CD2	1:A:100:HIS:HB2	2.53	0.44
1:A:538:LYS:O	1:A:618:PHE:HA	2.18	0.44
1:A:253:ARG:HH21	1:B:253:ARG:HH12	1.65	0.43
2:G:1:NAG:H61	2:G:2:NAG:O5	2.18	0.43
1:B:95:PHE:HB3	1:B:98:PHE:HB2	2.00	0.43
1:B:334:SER:O	1:B:336:ARG:HG2	2.18	0.43
1:A:626:ILE:O	1:A:650:GLY:HA2	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:658:ARG:HD3	1:A:660:GLU:HB2	2.00	0.43
1:B:159:PRO:HD3	1:B:216:TRP:CB	2.48	0.43
1:B:291:ALA:O	1:B:295:ILE:HG23	2.18	0.43
1:A:388:GLN:HB3	1:A:391:LYS:HB2	2.01	0.43
1:A:459:VAL:HG22	1:A:460:SER:N	2.34	0.43
1:B:111:GLY:O	1:B:137:LEU:HD12	2.19	0.43
1:B:184:ARG:HD2	1:B:187:TRP:CD2	2.53	0.43
1:B:272:ASN:HD22	1:B:274:ASP:N	2.17	0.43
1:A:73:GLU:O	1:A:74:ASN:HB2	2.19	0.43
1:A:139:LYS:HG3	1:A:141:GLN:CB	2.47	0.42
1:B:671:MET:HE1	1:B:682:HIS:CD2	2.54	0.42
1:A:387:PHE:CD1	1:A:394:CYS:HB3	2.54	0.42
1:A:175:LYS:CG	1:A:182:SER:HB3	2.47	0.42
1:B:546:VAL:HG22	1:B:547:TYR:N	2.35	0.42
1:A:55:LEU:CD1	1:A:561:LEU:HD12	2.49	0.42
2:D:2:NAG:O7	2:D:2:NAG:H3	2.20	0.42
1:B:258:LYS:NZ	1:B:712:HIS:CD2	2.76	0.42
1:B:60:LEU:HD12	1:B:60:LEU:O	2.20	0.42
1:A:484:SER:O	1:A:488:ASP:HA	2.20	0.41
1:B:501:ASP:O	1:B:505:GLN:HG2	2.20	0.41
1:A:709:ASP:O	1:A:712:HIS:HE1	2.02	0.41
1:A:390:ASP:O	1:A:391:LYS:HD2	2.21	0.41
1:A:184:ARG:HD2	1:A:187:TRP:CE2	2.56	0.41
1:A:658:ARG:O	1:A:658:ARG:HG3	2.19	0.41
1:B:487:ASN:ND2	1:B:489:LYS:H	2.18	0.41
1:A:74:ASN:HD22	3:A:801:NAG:H5	1.85	0.41
1:A:153:GLN:HE22	1:A:170:ASN:HD21	1.65	0.41
1:A:143:ILE:CD1	1:A:178:PRO:HB2	2.50	0.41
1:A:253:ARG:HH21	1:B:253:ARG:CZ	2.33	0.41
1:A:293:MET:HG2	1:A:315:TRP:HB3	2.02	0.41
1:A:554:LYS:CG	5:A:805:317:H43	2.44	0.41
1:B:626:ILE:O	1:B:650:GLY:HA2	2.20	0.41
1:A:291:ALA:O	1:A:295:ILE:HG23	2.20	0.41
1:A:552:SER:HB2	5:A:805:317:C43	2.51	0.41
1:A:765:LEU:HA	1:A:766:PRO:HD3	1.95	0.41
1:A:512:LYS:HE3	1:A:527:GLN:CD	2.45	0.40
1:B:136:ASP:OD2	1:B:139:LYS:HD3	2.21	0.40
1:B:487:ASN:ND2	1:B:487:ASN:H	2.20	0.40
1:A:751:ILE:HG12	1:A:755:MET:HE2	2.04	0.40
1:B:55:LEU:HD12	1:B:500:LEU:CD2	2.51	0.40
1:B:74:ASN:HA	1:B:74:ASN:HD22	1.69	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:125:ARG:HG2	1:A:126:HIS:NE2	2.37	0.40
1:A:136:ASP:CB	1:A:139:LYS:HG2	2.51	0.40
1:B:596:ARG:N	1:B:670:TYR:O	2.53	0.40
1:A:345:HIS:HE1	1:A:391:LYS:O	2.04	0.40
1:A:472:CYS:O	1:A:478:PRO:HA	2.22	0.40
1:B:54:ARG:HE	1:B:54:ARG:HB2	1.73	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	726/728 (100%)	702 (97%)	23 (3%)	1 (0%)	48	43
1	B	726/728 (100%)	701 (97%)	24 (3%)	1 (0%)	48	43
All	All	1452/1456 (100%)	1403 (97%)	47 (3%)	2 (0%)	48	43

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	320	GLN
1	A	320	GLN

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	653/653 (100%)	644 (99%)	9 (1%)	59	60
1	B	653/653 (100%)	643 (98%)	10 (2%)	57	58
All	All	1306/1306 (100%)	1287 (98%)	19 (2%)	57	58

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

Mol	Chain	Res	Type
1	A	170	ASN
1	A	184	ARG
1	A	388	GLN
1	A	448	GLU
1	A	472	CYS
1	A	502	LYS
1	A	520	ASN
1	A	608	GLU
1	A	728	VAL
1	B	55	LEU
1	B	57	LEU
1	B	66	HIS
1	B	74	ASN
1	B	254	VAL
1	B	272	ASN
1	B	385	CYS
1	B	487	ASN
1	B	658	ARG
1	B	685	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (31) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	74	ASN
1	A	75	ASN
1	A	100	HIS
1	A	138	ASN
1	A	169	ASN
1	A	170	ASN
1	A	227	GLN
1	A	247	GLN
1	A	388	GLN
1	A	505	GLN
1	A	572	ASN

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Mol	Chain	Res	Type
1	A	712	HIS
1	A	731	GLN
1	B	51	ASN
1	B	74	ASN
1	B	141	GLN
1	B	169	ASN
1	B	227	GLN
1	B	247	GLN
1	B	272	ASN
1	B	345	HIS
1	B	430	ASN
1	B	435	GLN
1	B	487	ASN
1	B	533	HIS
1	B	586	GLN
1	B	606	GLN
1	B	685	ASN
1	B	712	HIS
1	B	731	GLN
1	B	761	GLN

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 ⓘ

18 monosaccharides are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NAG	C	1	2,1	14,14,15	0.48	0	17,19,21	1.15	1 (5%)
2	NAG	C	2	2	14,14,15	0.63	0	17,19,21	1.13	0
2	NAG	D	1	2,1	14,14,15	0.81	1 (7%)	17,19,21	0.95	1 (5%)
2	NAG	D	2	2	14,14,15	0.84	1 (7%)	17,19,21	0.75	0
2	NAG	E	1	2,1	14,14,15	0.53	0	17,19,21	0.74	1 (5%)
2	NAG	E	2	2	14,14,15	0.67	0	17,19,21	0.79	1 (5%)
2	NAG	F	1	2,1	14,14,15	0.70	0	17,19,21	0.83	1 (5%)
2	NAG	F	2	2	14,14,15	0.84	1 (7%)	17,19,21	0.89	0
2	NAG	G	1	2,1	14,14,15	0.74	0	17,19,21	0.92	0
2	NAG	G	2	2	14,14,15	0.60	0	17,19,21	0.68	0
2	NAG	H	1	2,1	14,14,15	0.66	0	17,19,21	0.66	0
2	NAG	H	2	2	14,14,15	0.78	0	17,19,21	1.49	3 (17%)
2	NAG	I	1	2,1	14,14,15	0.57	0	17,19,21	0.69	0
2	NAG	I	2	2	14,14,15	0.48	0	17,19,21	0.69	0
2	NAG	J	1	2,1	14,14,15	0.62	0	17,19,21	0.78	0
2	NAG	J	2	2	14,14,15	0.60	0	17,19,21	0.80	1 (5%)
2	NAG	K	1	2,1	14,14,15	0.40	0	17,19,21	0.82	1 (5%)
2	NAG	K	2	2	14,14,15	0.59	0	17,19,21	0.71	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	C	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	C	2	2	1/1/5/7	2/6/23/26	0/1/1/1
2	NAG	D	1	2,1	1/1/5/7	0/6/23/26	0/1/1/1
2	NAG	D	2	2	-	4/6/23/26	0/1/1/1
2	NAG	E	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	E	2	2	-	2/6/23/26	0/1/1/1
2	NAG	F	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	F	2	2	1/1/5/7	2/6/23/26	0/1/1/1
2	NAG	G	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	G	2	2	-	2/6/23/26	0/1/1/1
2	NAG	H	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	H	2	2	-	3/6/23/26	0/1/1/1
2	NAG	I	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	I	2	2	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	J	1	2,1	-	2/6/23/26	0/1/1/1
2	NAG	J	2	2	-	0/6/23/26	0/1/1/1
2	NAG	K	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	K	2	2	-	2/6/23/26	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	2	NAG	C1-C2	2.54	1.55	1.52
2	D	1	NAG	C1-C2	2.39	1.55	1.52
2	D	2	NAG	C1-C2	2.31	1.55	1.52

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	2	NAG	C4-C3-C2	4.02	116.91	111.02
2	H	2	NAG	C8-C7-N2	2.54	120.33	116.12
2	E	1	NAG	C2-N2-C7	-2.40	119.68	122.90
2	F	1	NAG	C2-N2-C7	-2.39	119.70	122.90
2	K	1	NAG	C2-N2-C7	-2.36	119.74	122.90
2	C	1	NAG	C4-C3-C2	-2.33	107.61	111.02
2	D	1	NAG	C2-N2-C7	-2.20	119.95	122.90
2	E	2	NAG	C2-N2-C7	-2.20	119.95	122.90
2	H	2	NAG	C1-C2-N2	-2.09	107.14	110.43
2	J	2	NAG	C2-N2-C7	-2.08	120.11	122.90

All (3) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	C	2	NAG	C1
2	D	1	NAG	C1
2	F	2	NAG	C1

All (19) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	2	NAG	O5-C5-C6-O6
2	K	2	NAG	O5-C5-C6-O6
2	D	2	NAG	O5-C5-C6-O6
2	K	2	NAG	C4-C5-C6-O6
2	C	2	NAG	C4-C5-C6-O6

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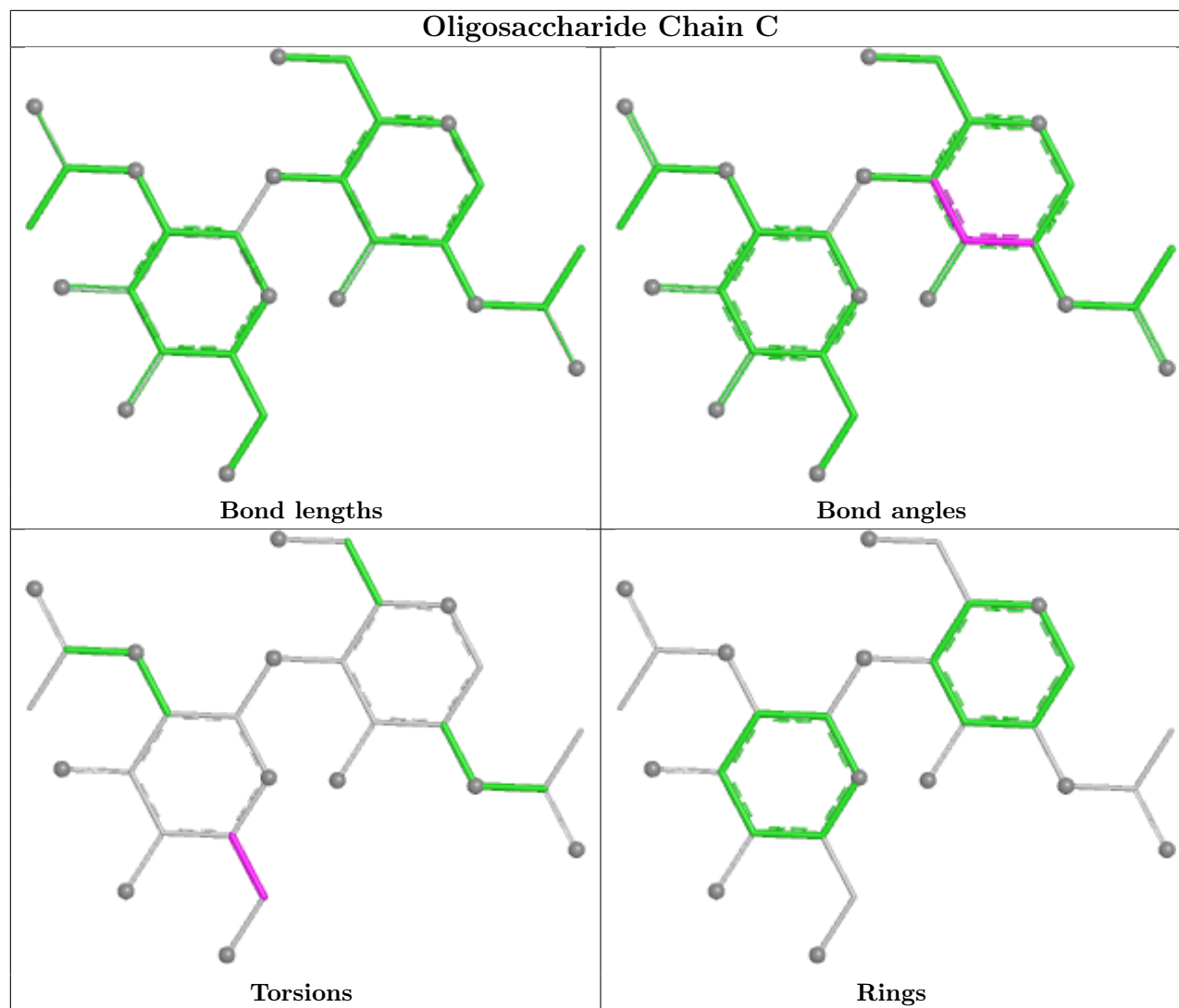
Mol	Chain	Res	Type	Atoms
2	F	2	NAG	C4-C5-C6-O6
2	D	2	NAG	C4-C5-C6-O6
2	F	2	NAG	O5-C5-C6-O6
2	J	1	NAG	C4-C5-C6-O6
2	H	2	NAG	C3-C2-N2-C7
2	J	1	NAG	O5-C5-C6-O6
2	G	2	NAG	C1-C2-N2-C7
2	H	2	NAG	C4-C5-C6-O6
2	D	2	NAG	C3-C2-N2-C7
2	G	2	NAG	C3-C2-N2-C7
2	E	2	NAG	C4-C5-C6-O6
2	H	2	NAG	O5-C5-C6-O6
2	D	2	NAG	C1-C2-N2-C7
2	E	2	NAG	O5-C5-C6-O6

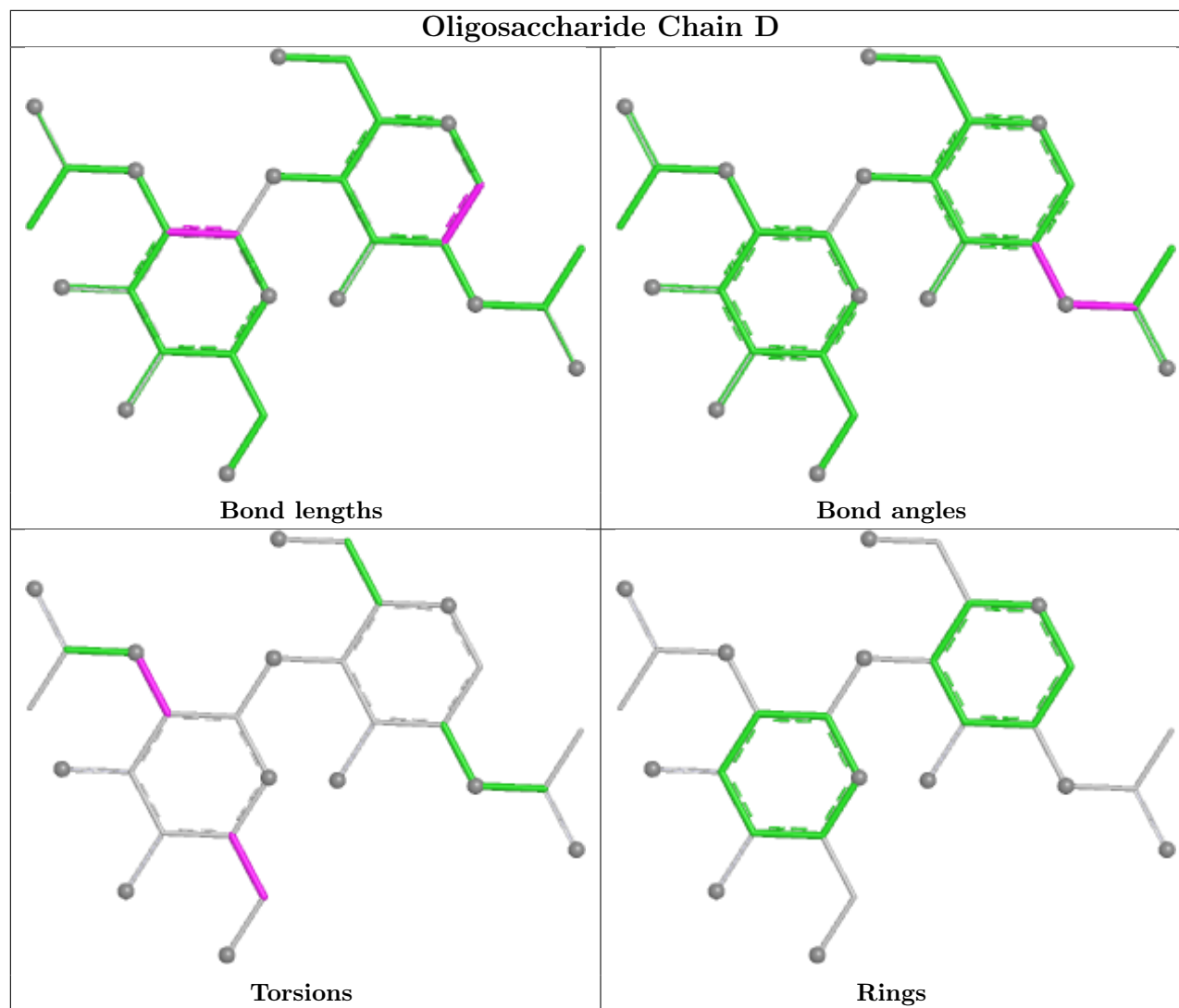
There are no ring outliers.

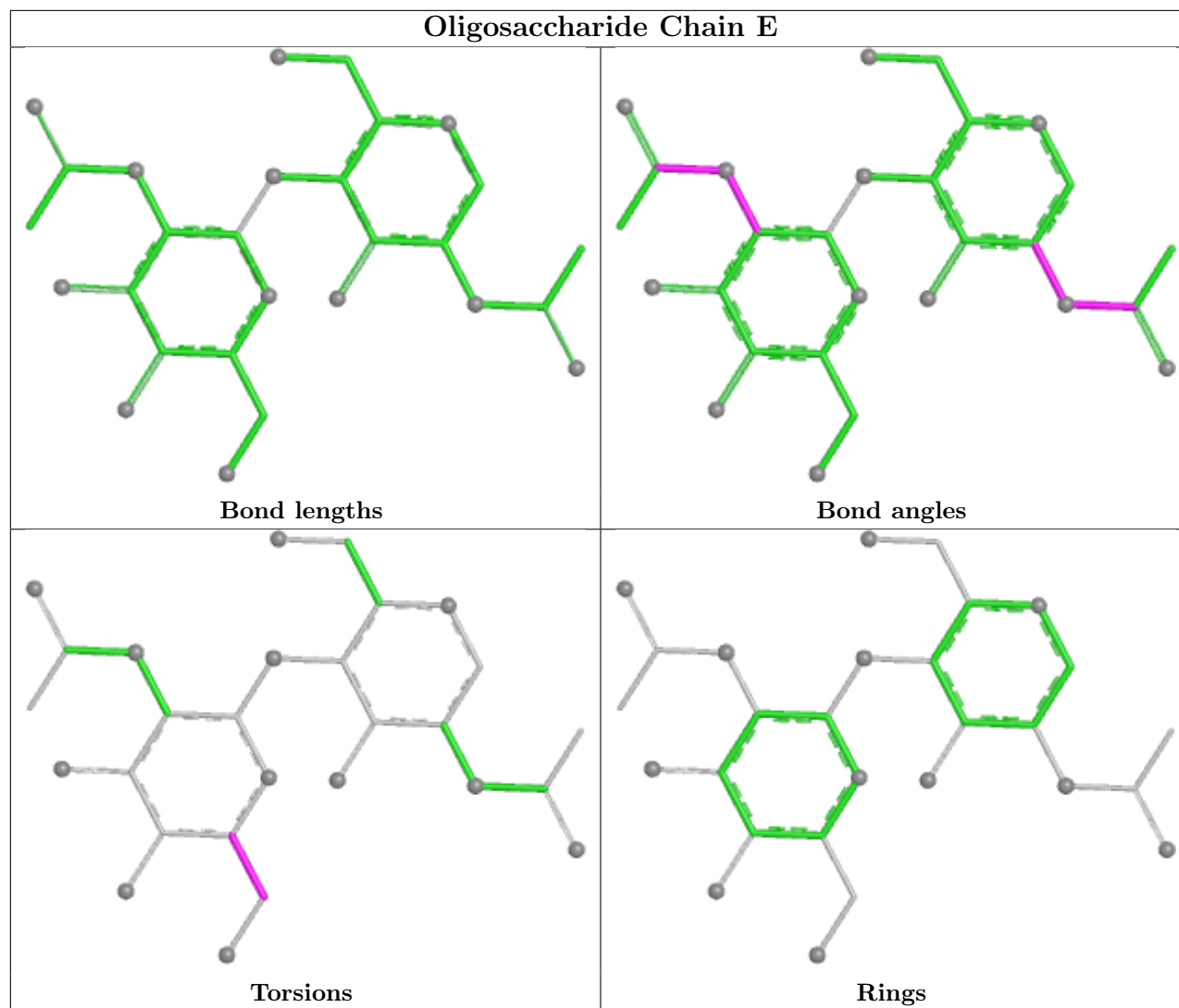
6 monomers are involved in 6 short contacts:

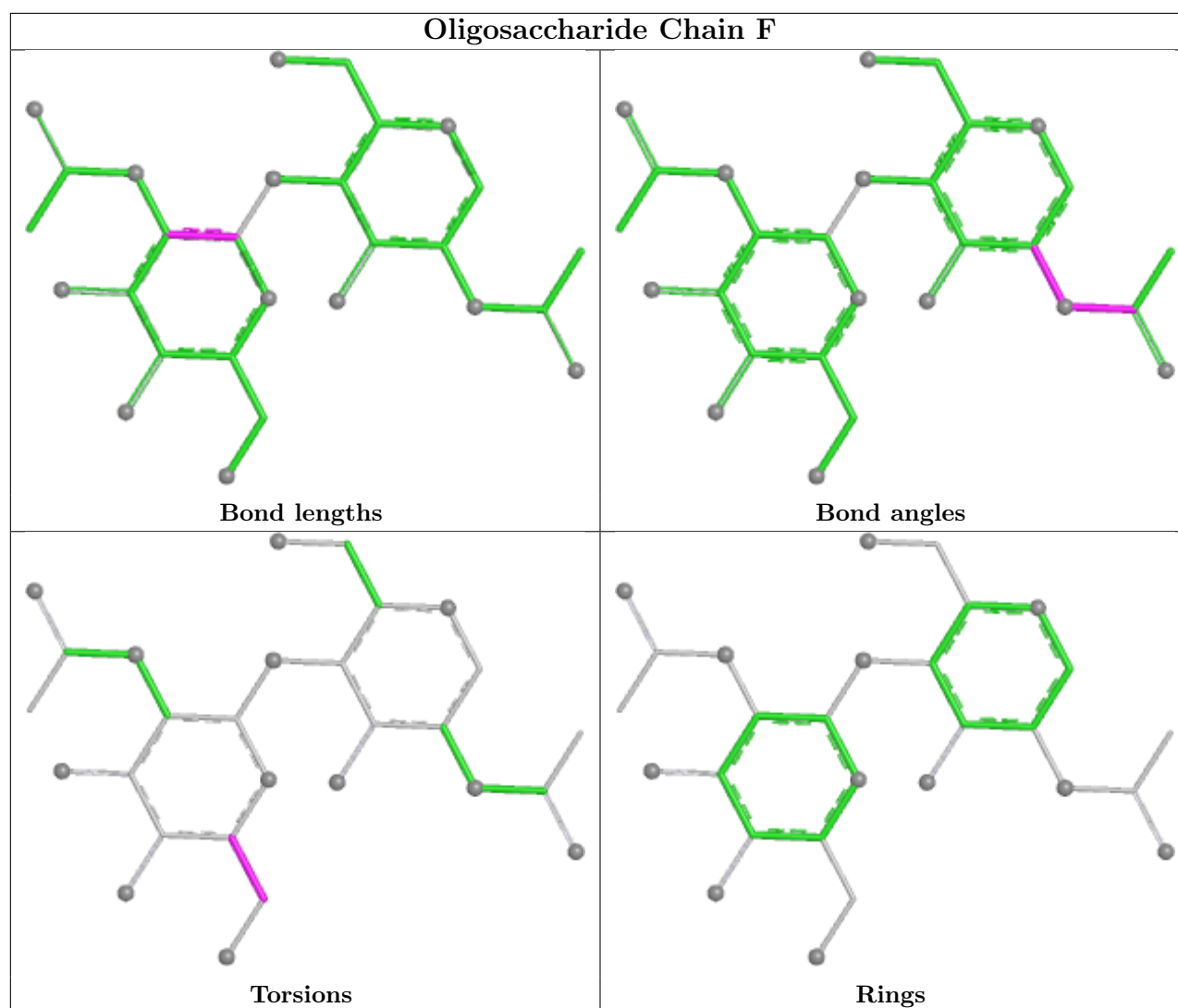
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	2	NAG	2	0
2	F	1	NAG	2	0
2	G	2	NAG	1	0
2	F	2	NAG	2	0
2	D	1	NAG	2	0
2	G	1	NAG	1	0

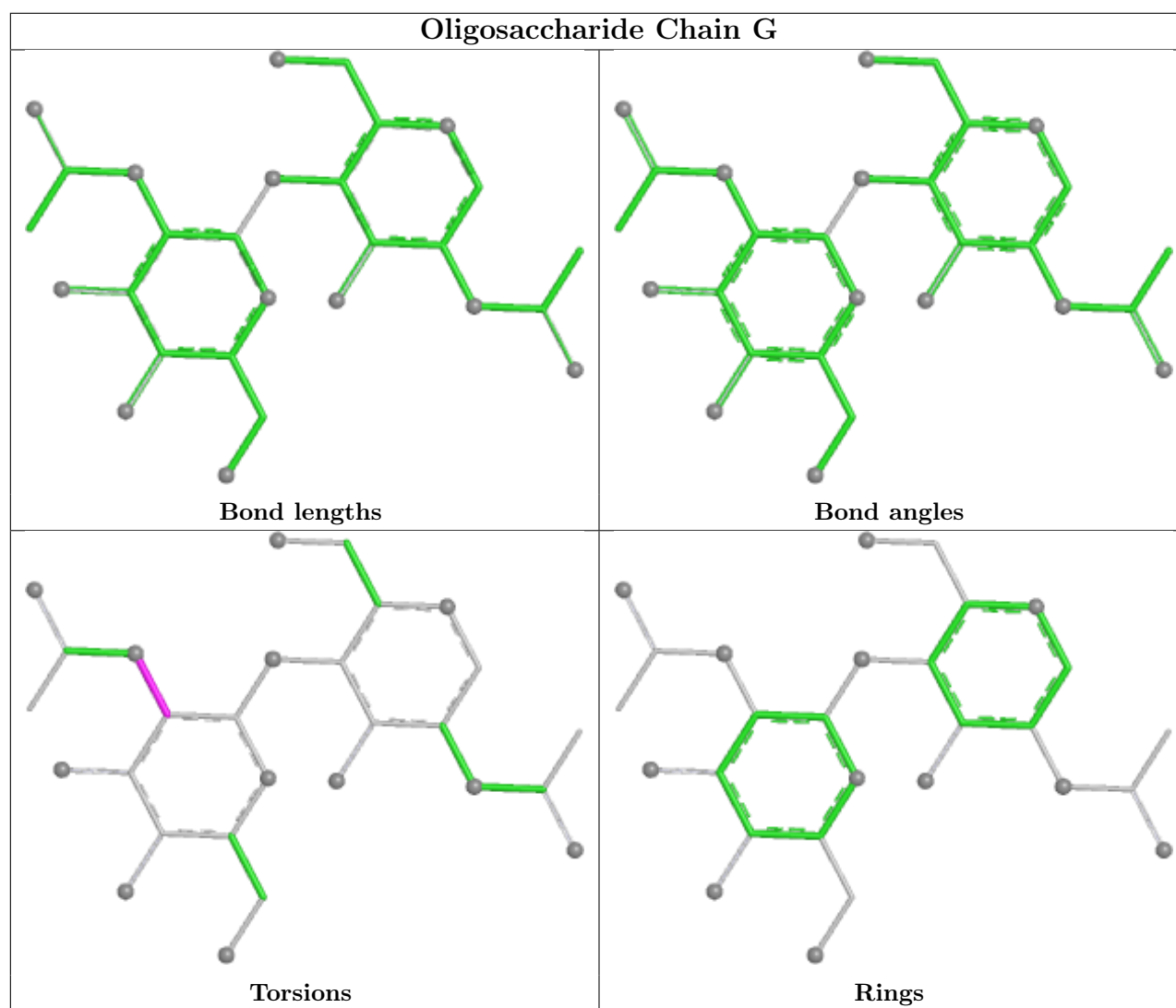
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

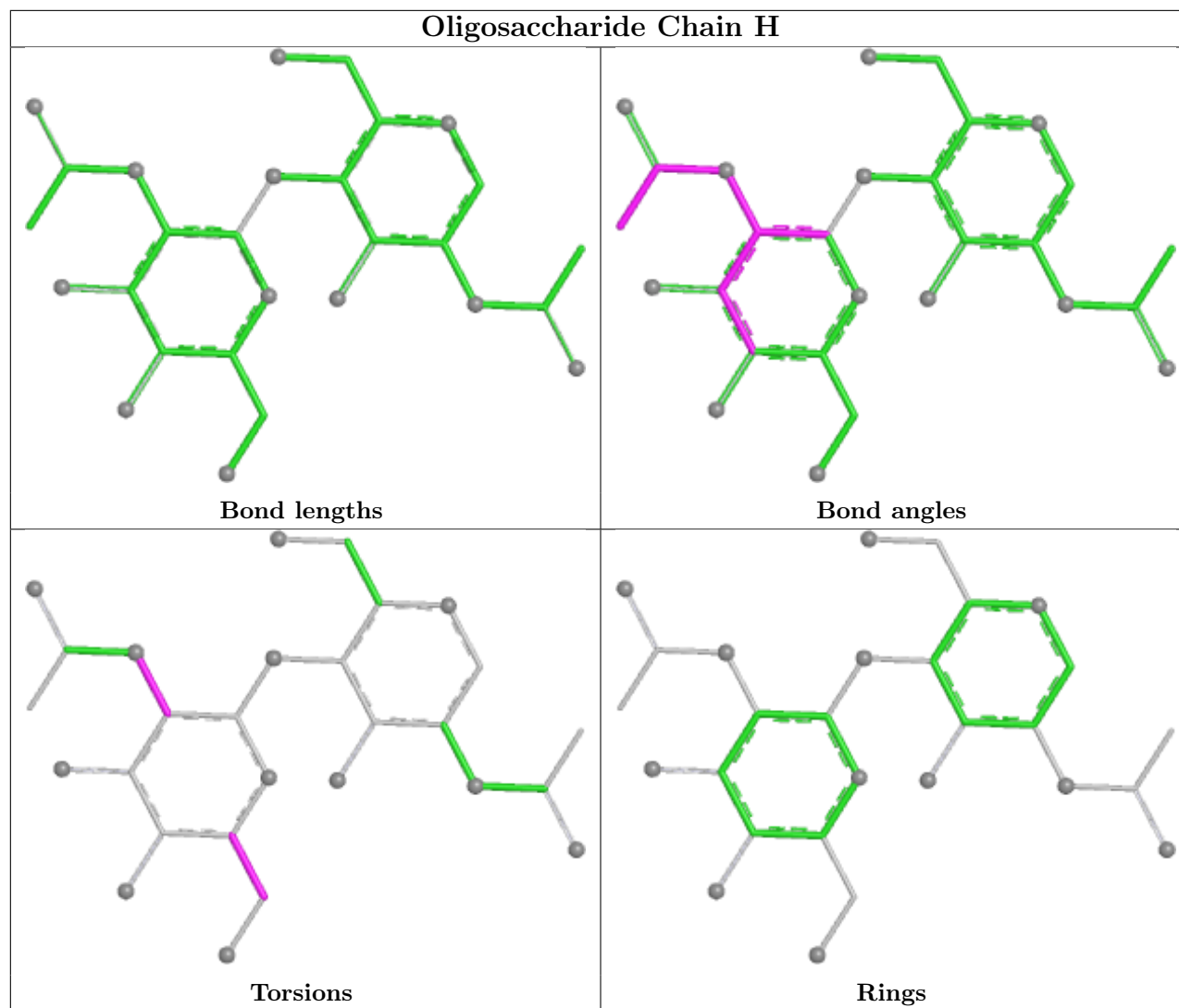


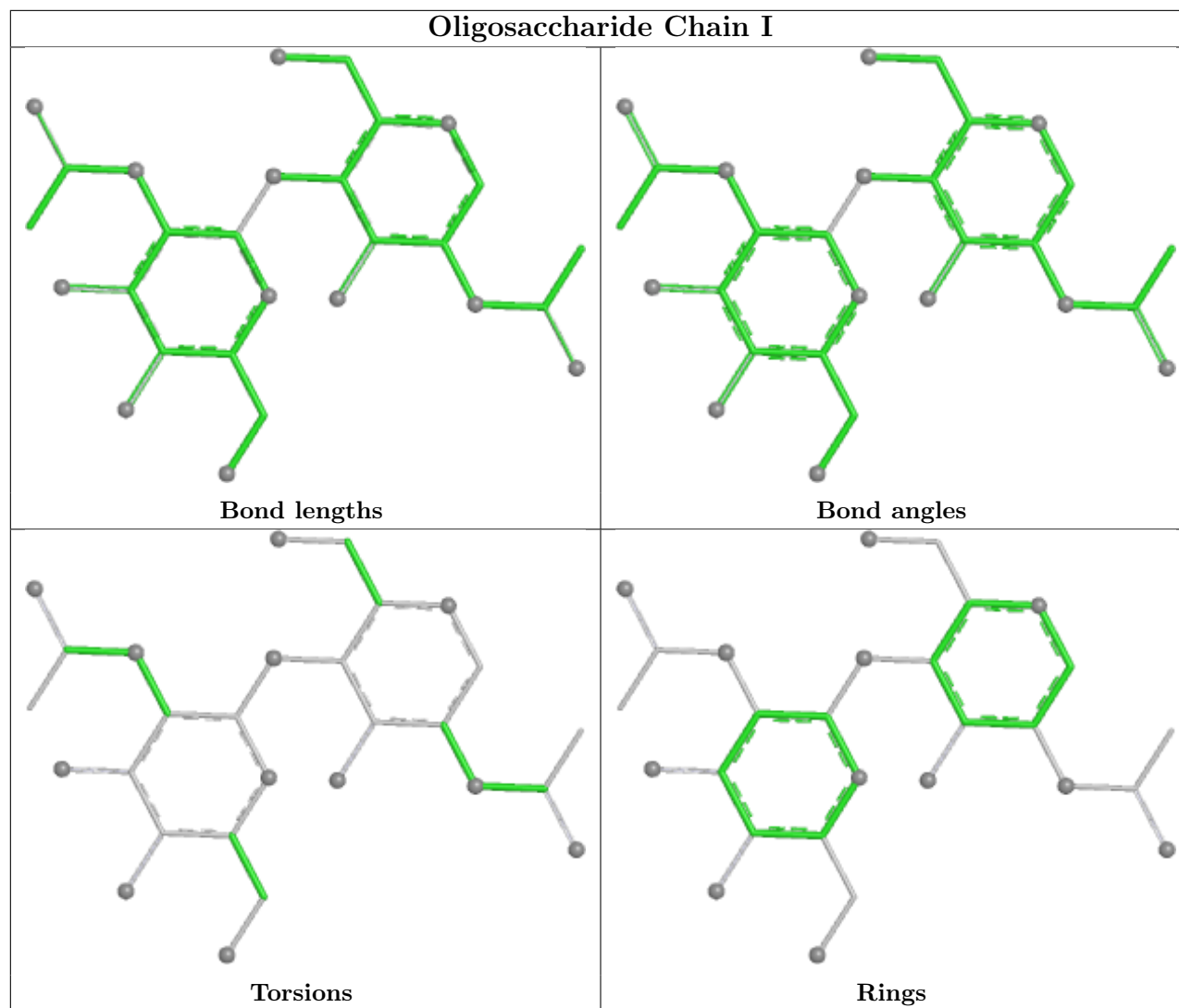


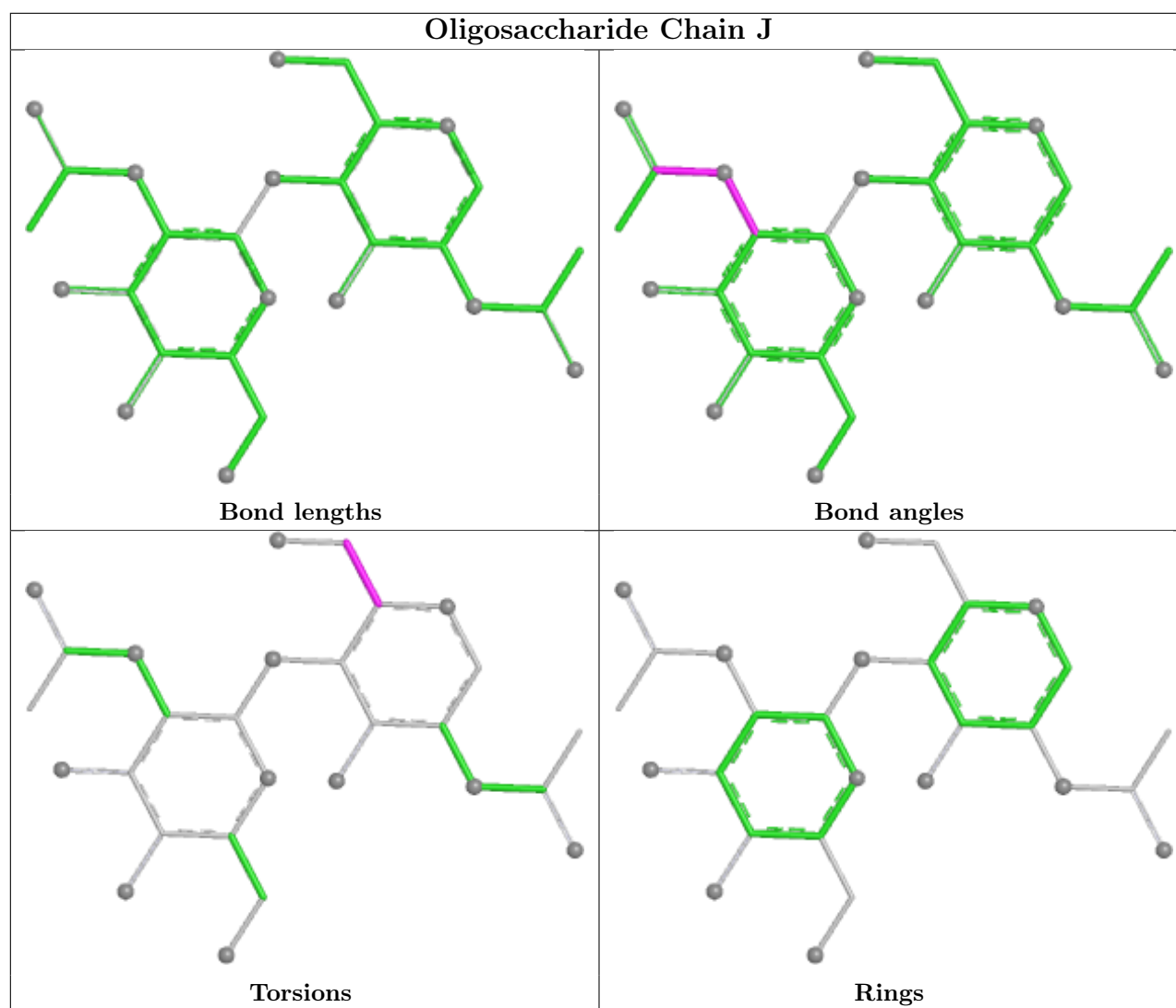


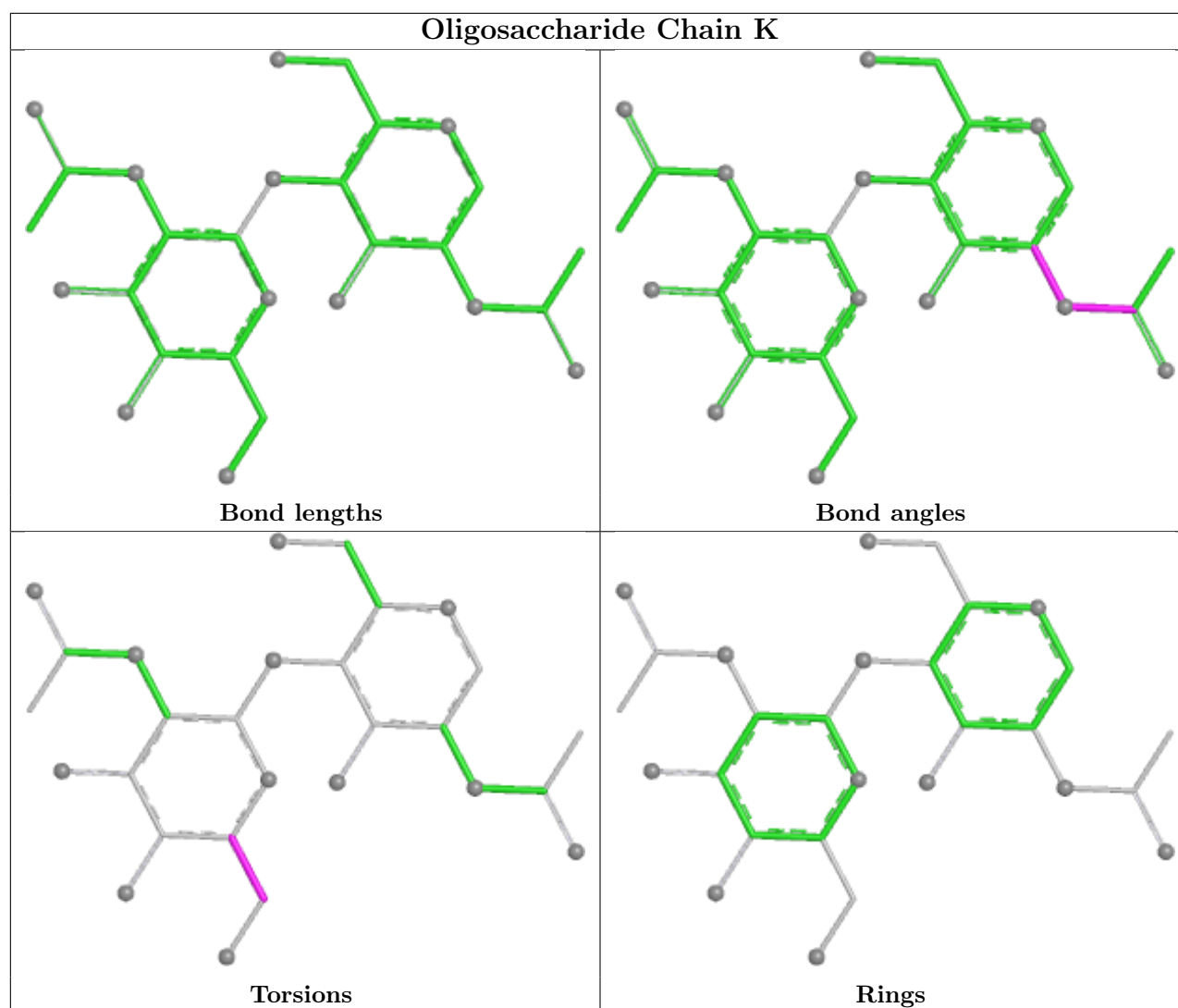












5.6 Ligand geometry [i](#)

Of 9 ligands modelled in this entry, 1 is monoatomic - leaving 8 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	NAG	A	802	1	14,14,15	0.72	0	17,19,21	0.75	1 (5%)
5	317	A	805	-	32,33,33	1.57	5 (15%)	40,49,49	2.26	13 (32%)
5	317	B	804	-	32,33,33	1.74	6 (18%)	40,49,49	2.29	14 (35%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	NAG	B	803	1	14,14,15	0.51	0	17,19,21	1.03	1 (5%)
3	NAG	A	801	1	14,14,15	0.70	0	17,19,21	0.62	0
3	NAG	B	802	1	14,14,15	0.74	0	17,19,21	0.71	1 (5%)
3	NAG	A	803	1	14,14,15	0.77	0	17,19,21	0.69	0
3	NAG	B	801	1	14,14,15	0.81	1 (7%)	17,19,21	0.76	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	A	802	1	-	0/6/23/26	0/1/1/1
5	317	A	805	-	-	6/21/39/39	0/4/4/4
5	317	B	804	-	-	6/21/39/39	0/4/4/4
3	NAG	B	803	1	-	0/6/23/26	0/1/1/1
3	NAG	A	801	1	-	2/6/23/26	0/1/1/1
3	NAG	B	802	1	-	2/6/23/26	0/1/1/1
3	NAG	A	803	1	-	2/6/23/26	0/1/1/1
3	NAG	B	801	1	-	3/6/23/26	0/1/1/1

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	804	317	C4-C3	4.60	1.37	1.32
5	B	804	317	C31-C36	4.44	1.45	1.39
5	A	805	317	C31-C36	4.12	1.45	1.39
5	B	804	317	C28-N29	4.03	1.36	1.30
5	A	805	317	C28-N29	3.61	1.36	1.30
5	A	805	317	C4-C3	3.14	1.36	1.32
5	B	804	317	C33-C34	2.62	1.43	1.38
5	B	804	317	C32-C31	2.41	1.43	1.39
5	A	805	317	C32-C31	2.31	1.43	1.39
3	B	801	NAG	C1-C2	2.25	1.55	1.52
5	B	804	317	C26-N27	2.24	1.33	1.29
5	A	805	317	C2-C3	2.12	1.52	1.49

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	805	317	C26-O30-N29	-6.50	103.66	106.18
5	B	804	317	C26-O30-N29	-6.37	103.71	106.18
5	A	805	317	O30-C26-C1	5.39	122.95	117.66
5	B	804	317	O30-N29-C28	4.88	106.14	103.28
5	B	804	317	O30-C26-C1	4.85	122.43	117.66
5	A	805	317	N27-C28-N29	-4.48	111.20	114.47
5	A	805	317	O30-N29-C28	4.44	105.88	103.28
5	B	804	317	N27-C28-N29	-4.13	111.46	114.47
5	B	804	317	O42-S40-C34	-3.32	105.65	108.23
5	B	804	317	C31-C36-CL47	3.22	125.77	121.01
5	B	804	317	C31-C28-N27	3.18	129.06	122.08
5	A	805	317	C31-C28-N27	3.09	128.86	122.08
5	A	805	317	C32-C31-C28	-3.09	113.12	118.29
5	B	804	317	C31-C28-N29	-3.03	119.42	124.32
5	B	804	317	C32-C31-C28	-3.02	113.24	118.29
5	A	805	317	O41-S40-C43	2.93	112.76	108.47
5	B	804	317	C35-C36-CL47	-2.89	113.72	118.45
5	B	804	317	C7-C18-C20	-2.87	115.72	119.71
5	A	805	317	C31-C28-N29	-2.75	119.87	124.32
5	B	804	317	O41-S40-C43	2.72	112.45	108.47
5	A	805	317	C31-C36-CL47	2.42	124.59	121.01
3	B	803	NAG	C2-N2-C7	-2.28	119.85	122.90
5	A	805	317	C7-C1-C26	-2.27	104.66	109.96
3	B	802	NAG	C2-N2-C7	-2.23	119.91	122.90
5	A	805	317	C35-C36-CL47	-2.23	114.80	118.45
5	A	805	317	C36-C31-C28	2.16	128.59	125.11
5	B	804	317	C7-C1-C26	-2.13	104.98	109.96
5	A	805	317	C1-C2-C3	2.12	114.87	111.33
5	B	804	317	C36-C31-C28	2.11	128.53	125.11
3	A	802	NAG	C2-N2-C7	-2.09	120.10	122.90

There are no chirality outliers.

All (21) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	803	NAG	C1-C2-N2-C7
5	B	804	317	C20-C18-C7-C1
3	B	801	NAG	O5-C5-C6-O6
3	A	801	NAG	C4-C5-C6-O6
3	A	801	NAG	O5-C5-C6-O6
3	B	801	NAG	C4-C5-C6-O6
5	B	804	317	C35-C34-S40-O42
5	A	805	317	C35-C34-S40-O42

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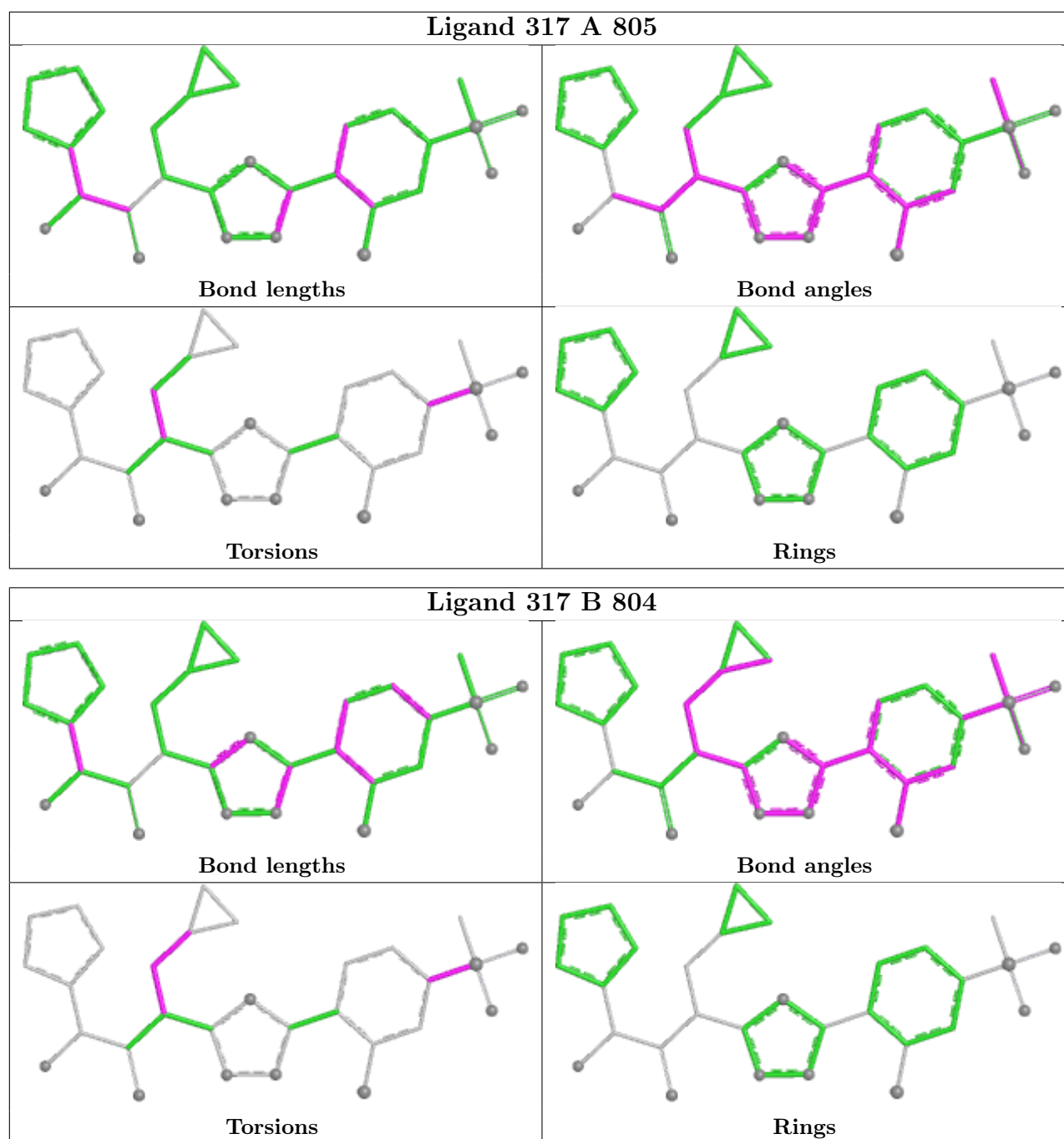
Mol	Chain	Res	Type	Atoms
5	A	805	317	C33-C34-S40-O42
5	B	804	317	C33-C34-S40-O42
5	B	804	317	C35-C34-S40-C43
5	A	805	317	C35-C34-S40-C43
5	A	805	317	C33-C34-S40-C43
5	B	804	317	C33-C34-S40-C43
3	B	802	NAG	C4-C5-C6-O6
3	B	802	NAG	O5-C5-C6-O6
5	A	805	317	C26-C1-C7-C18
5	B	804	317	C2-C1-C7-C18
3	A	803	NAG	C3-C2-N2-C7
3	B	801	NAG	C3-C2-N2-C7
5	A	805	317	C2-C1-C7-C18

There are no ring outliers.

5 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	805	317	4	0
5	B	804	317	2	0
3	A	801	NAG	2	0
3	A	803	NAG	2	0
3	B	801	NAG	1	0

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



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	728/728 (100%)	-0.12	13 (1%) 67 69	15, 25, 44, 57	0
1	B	728/728 (100%)	-0.12	12 (1%) 70 72	16, 26, 44, 56	0
All	All	1456/1456 (100%)	-0.12	25 (1%) 69 70	15, 26, 44, 57	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	766	PRO	3.8
1	A	83	TYR	3.2
1	A	39	THR	3.1
1	B	73	GLU	2.7
1	B	766	PRO	2.7
1	B	39	THR	2.6
1	B	74	ASN	2.6
1	A	142	LEU	2.6
1	B	91	GLU	2.5
1	A	92	ASN	2.5
1	A	521	GLU	2.4
1	A	75	ASN	2.4
1	B	66	HIS	2.4
1	A	141	GLN	2.3
1	B	75	ASN	2.3
1	A	332	GLU	2.3
1	A	278	SER	2.2
1	B	379	GLU	2.2
1	B	97	GLU	2.1
1	B	143	ILE	2.1
1	A	40	ARG	2.1
1	A	74	ASN	2.1
1	B	99	GLY	2.1
1	B	522	THR	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	73	GLU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

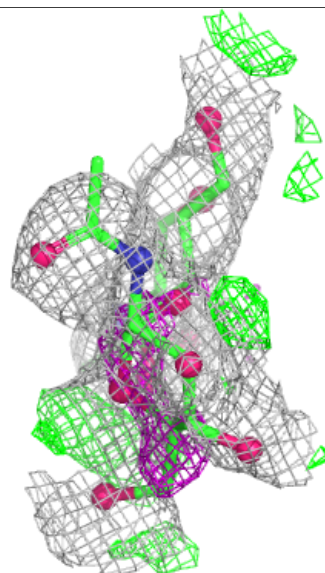
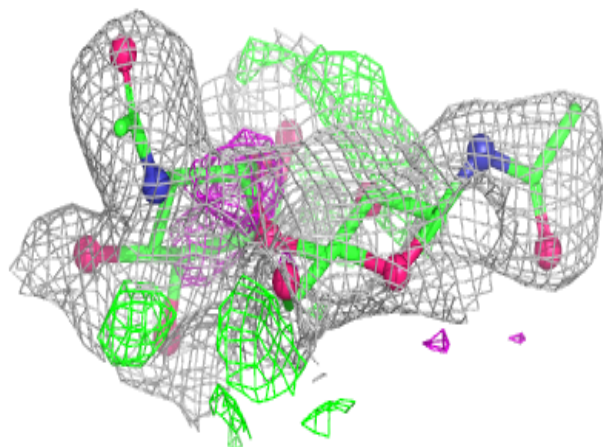
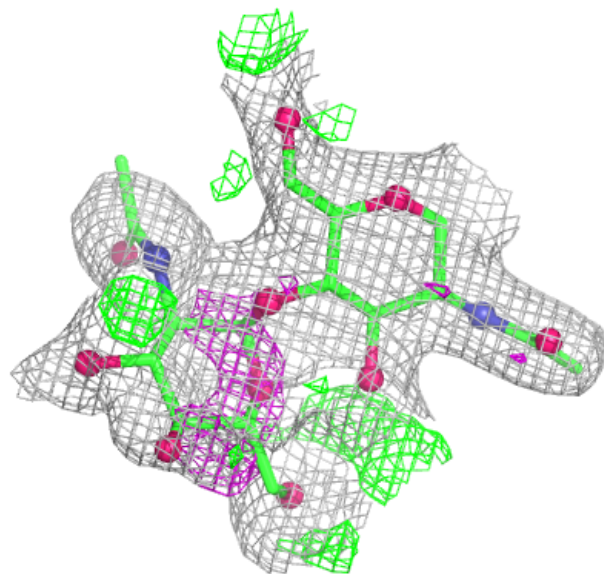
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	NAG	G	2	14/15	0.45	0.20	55,57,59,60	0
2	NAG	H	2	14/15	0.52	0.19	55,57,58,58	0
2	NAG	C	2	14/15	0.54	0.17	56,57,58,58	0
2	NAG	D	2	14/15	0.59	0.17	53,56,58,58	0
2	NAG	D	1	14/15	0.60	0.17	50,53,55,55	0
2	NAG	F	2	14/15	0.60	0.19	50,53,56,57	0
2	NAG	K	2	14/15	0.67	0.19	48,51,53,54	0
2	NAG	E	2	14/15	0.76	0.14	48,50,52,52	0
2	NAG	G	1	14/15	0.78	0.13	41,45,47,52	0
2	NAG	C	1	14/15	0.80	0.13	45,47,52,54	0
2	NAG	J	2	14/15	0.81	0.11	42,44,46,47	0
2	NAG	H	1	14/15	0.84	0.11	42,47,49,52	0
2	NAG	I	2	14/15	0.84	0.12	45,47,49,49	0
2	NAG	E	1	14/15	0.86	0.11	40,43,48,48	0
2	NAG	F	1	14/15	0.91	0.09	34,37,42,45	0
2	NAG	K	1	14/15	0.92	0.09	34,37,41,45	0
2	NAG	J	1	14/15	0.93	0.07	29,30,34,38	0
2	NAG	I	1	14/15	0.93	0.08	35,39,43,43	0

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

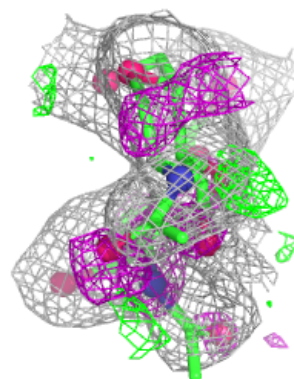
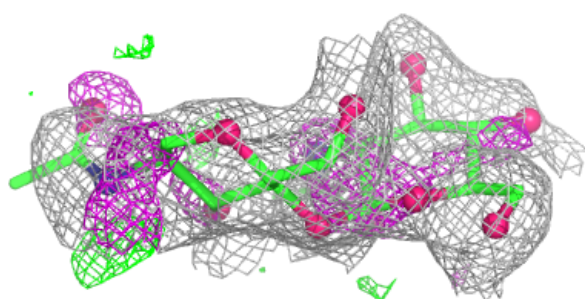
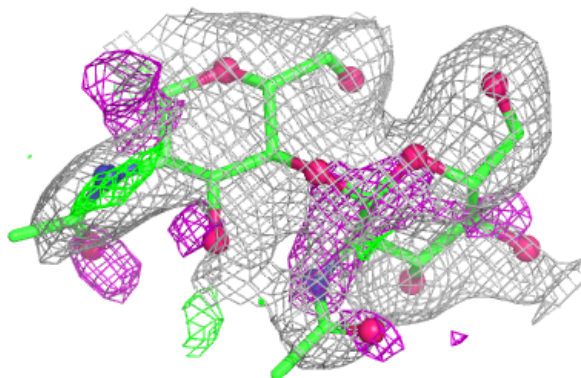
Electron density around Chain C:

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

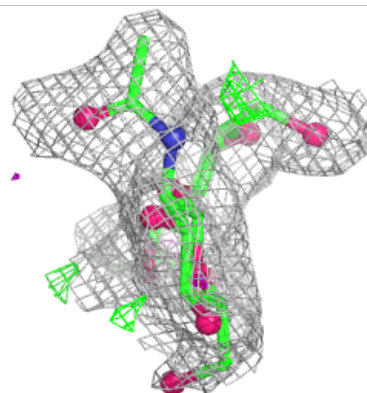
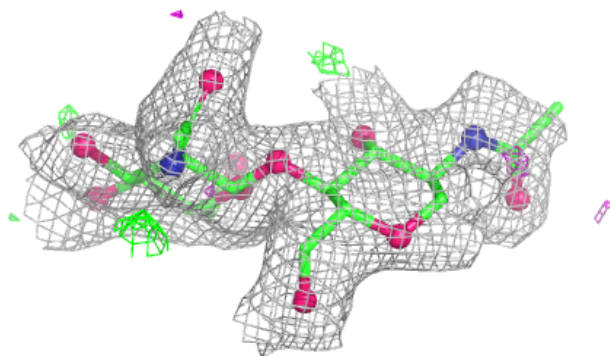
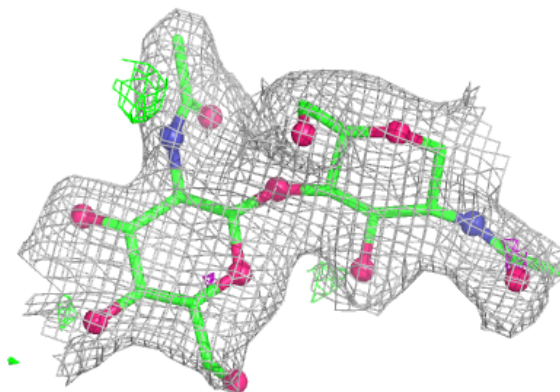


Electron density around Chain D:

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

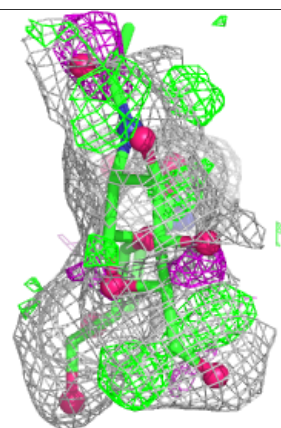
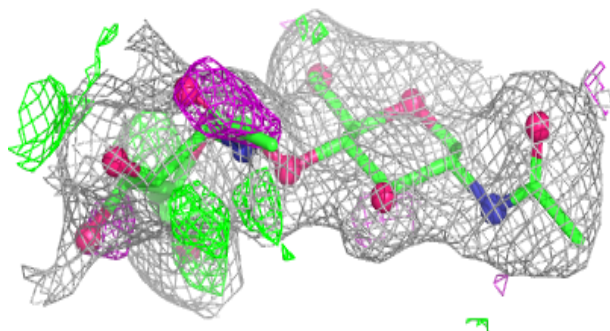
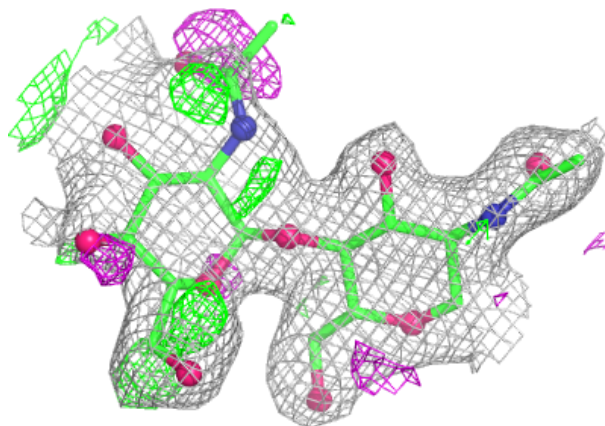
**Electron density around Chain E:**

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



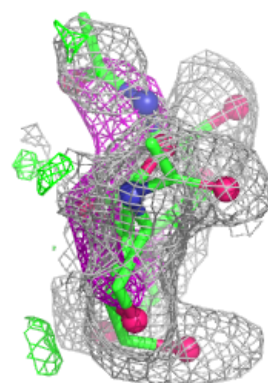
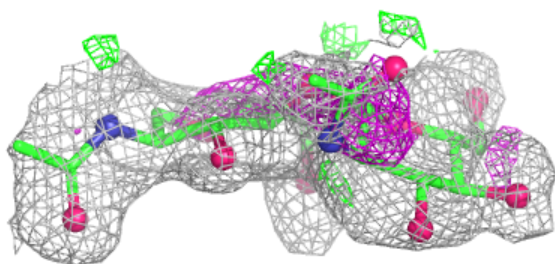
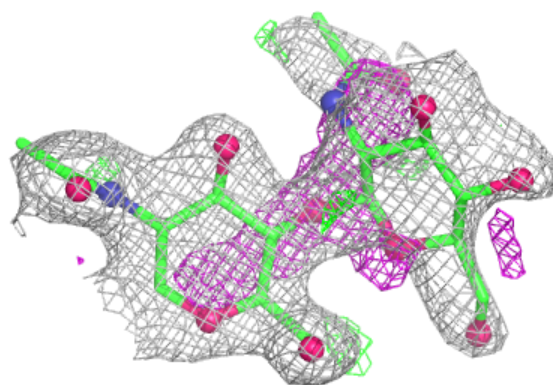
Electron density around Chain F:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

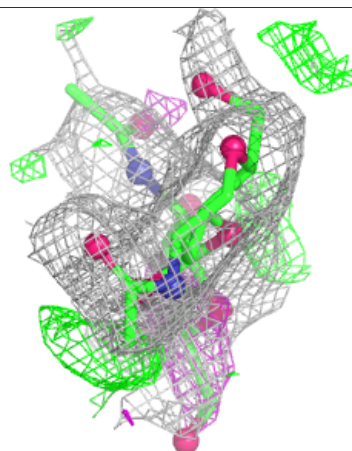
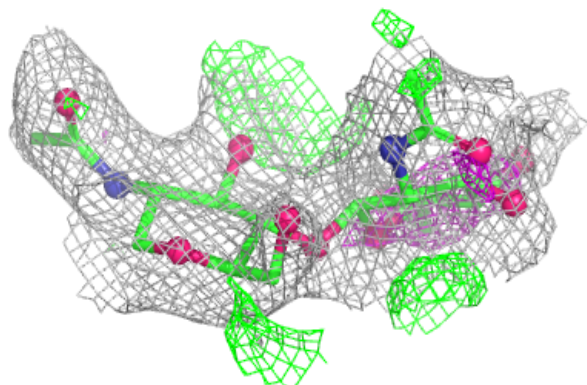
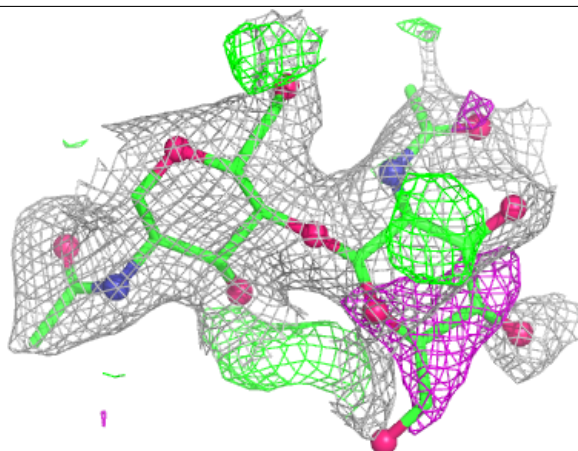


Electron density around Chain G:

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

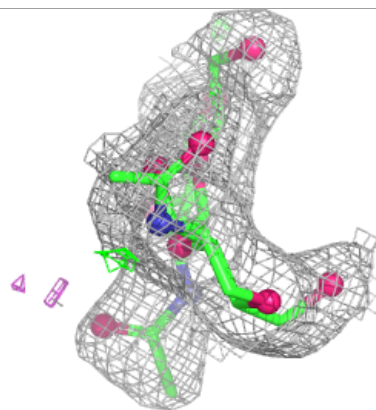
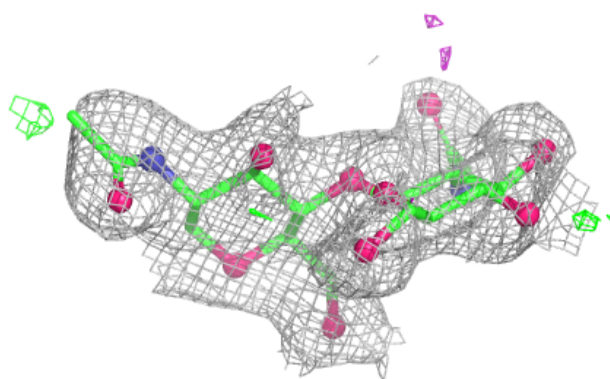
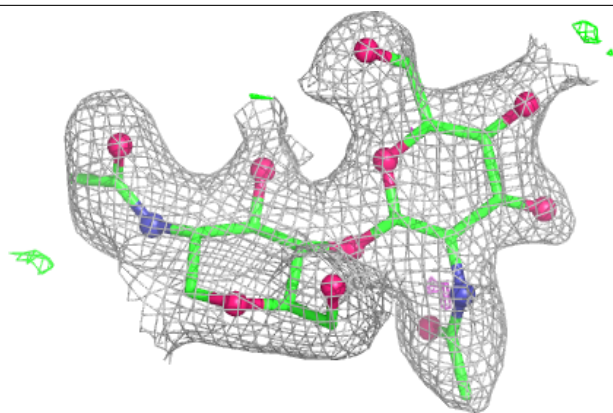
**Electron density around Chain H:**

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



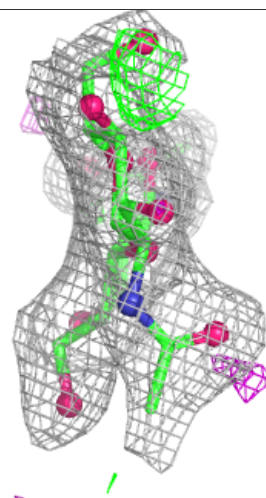
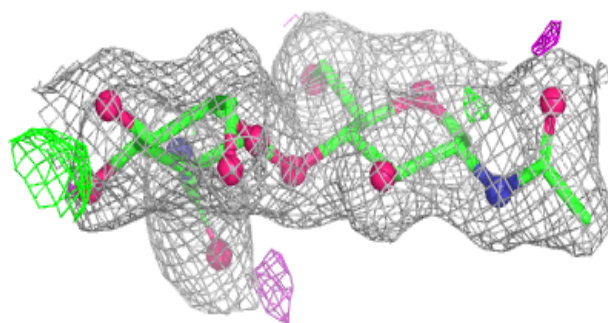
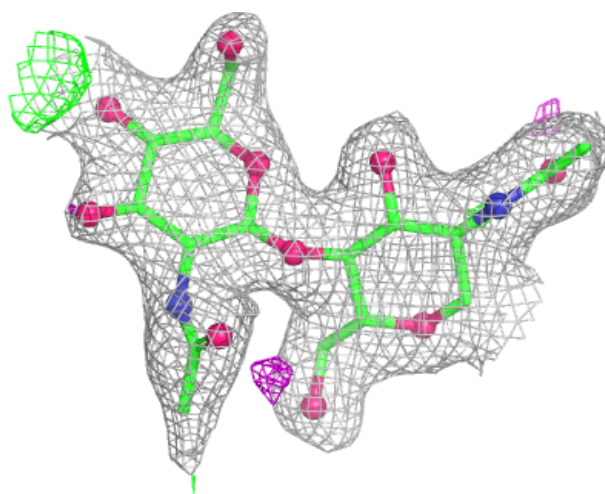
Electron density around Chain I:

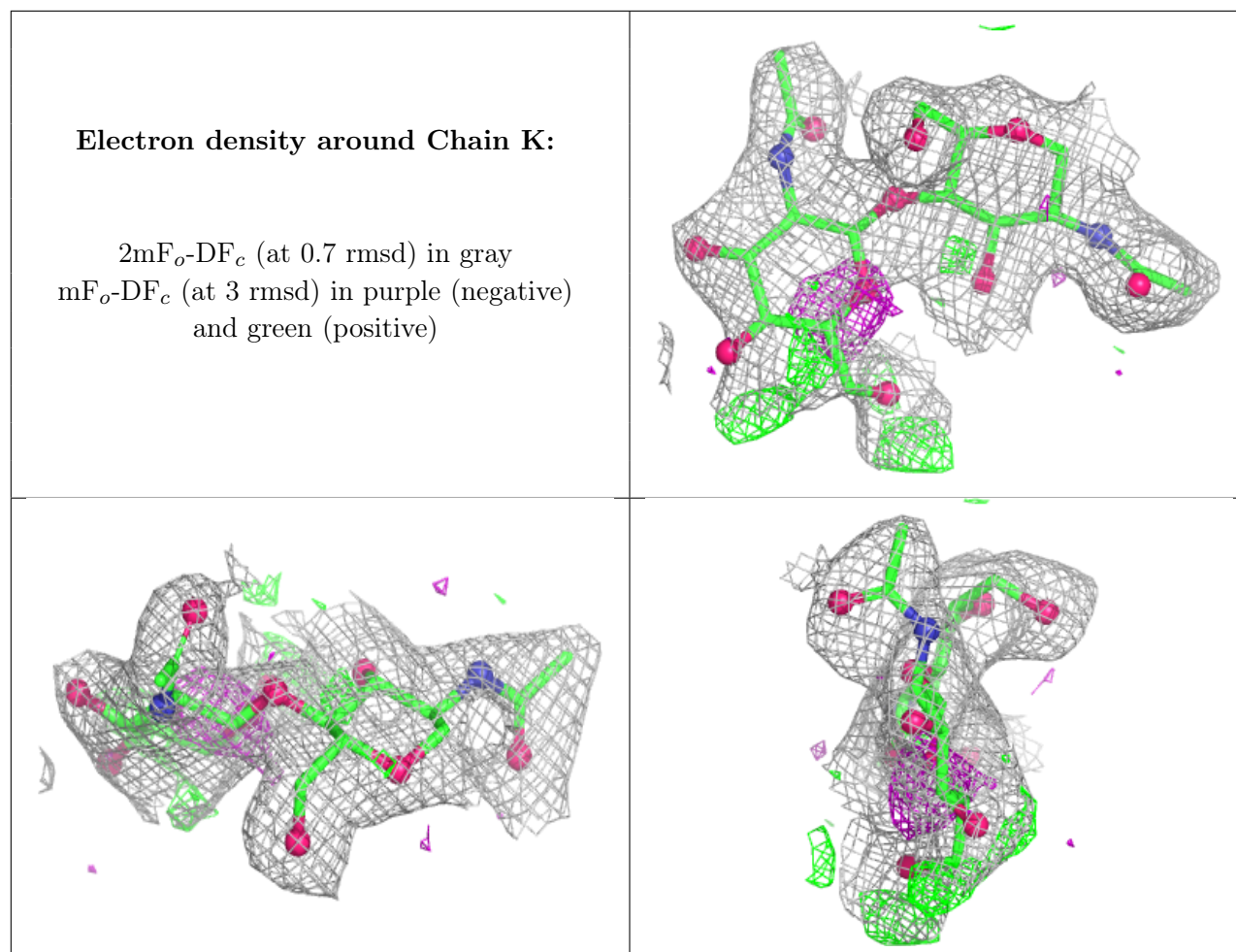
$2mF_o - DF_c$ (at 0.7 rmsd) in gray
 $mF_o - DF_c$ (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain J:

$2mF_o-DF_c$ (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.

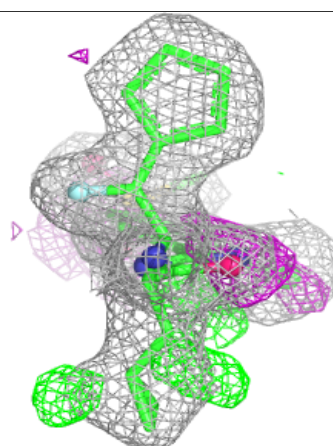
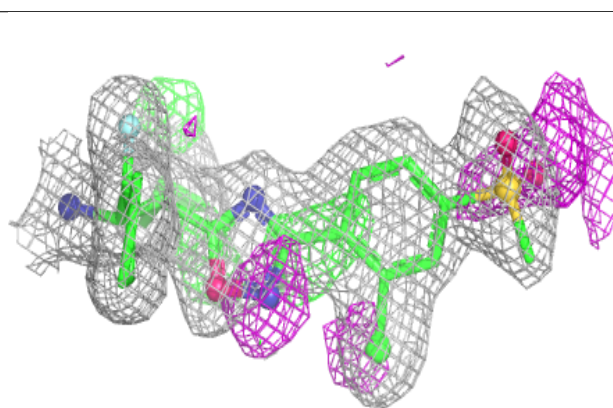
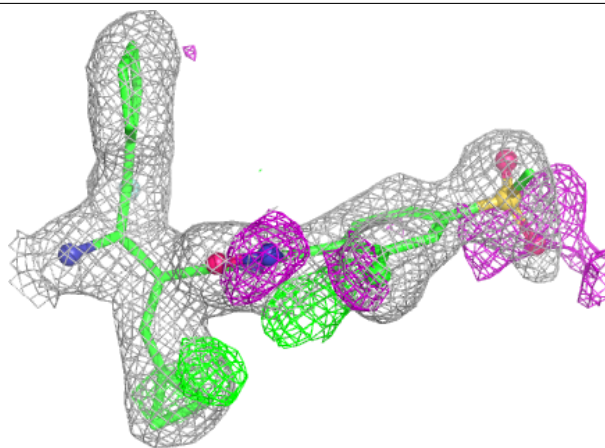
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	NAG	A	801	14/15	0.49	0.17	58,59,59,59	0
3	NAG	B	801	14/15	0.49	0.16	57,57,58,58	0
3	NAG	A	803	14/15	0.56	0.17	54,55,56,57	0
3	NAG	B	802	14/15	0.78	0.12	49,51,53,53	0
3	NAG	B	803	14/15	0.81	0.12	38,41,44,44	0
3	NAG	A	802	14/15	0.82	0.12	45,47,48,48	0
5	317	B	804	30/30	0.88	0.12	21,32,38,41	0
5	317	A	805	30/30	0.91	0.11	20,32,35,35	0
4	NA	A	804	1/1	0.93	0.06	31,31,31,31	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.

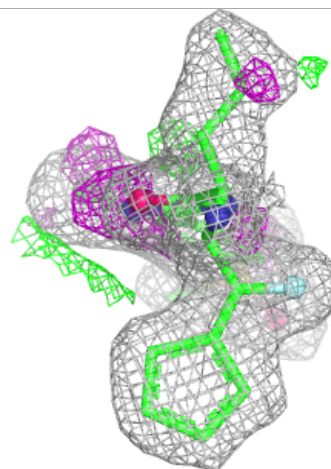
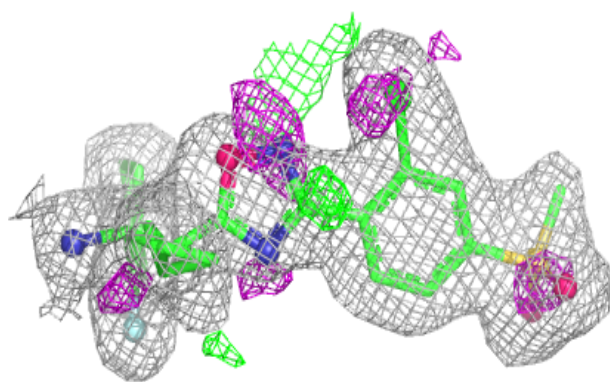
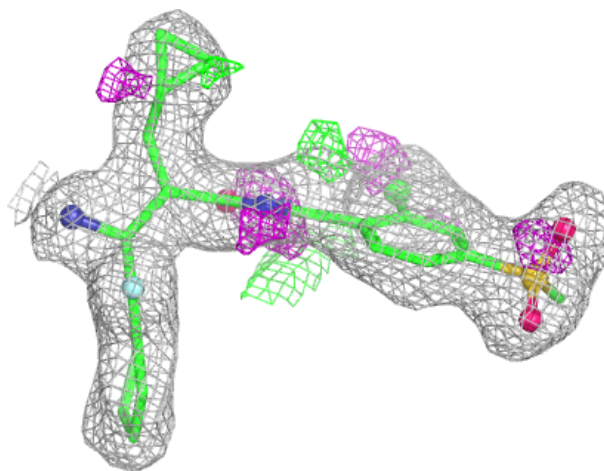
Electron density around 317 B 804:

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



Electron density around 317 A 805:

$2mF_o-DF_c$ (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.