



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 7, 2020 – 11:40 AM BST

PDB ID : 1H1M
Title : CRYSTAL STRUCTURE OF QUERCETIN 2,3-DIOXYGENASE ANAEROBICALLY COMPLEXED WITH THE SUBSTRATE KAEMPFEROL
Authors : Steiner, R.A.; Dijkstra, B.W.
Deposited on : 2002-07-19
Resolution : 1.90 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

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

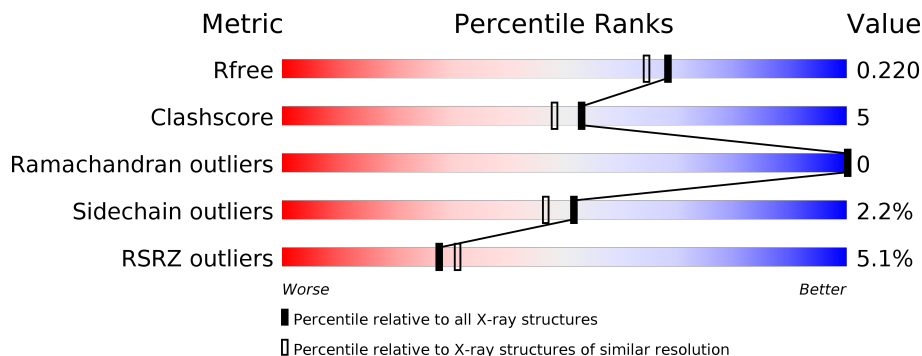
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

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

Mol	Chain	Length	Quality of chain
1	A	350	
1	B	350	
1	C	350	
1	D	350	
2	E	5	
3	F	4	

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Mol	Chain	Length	Quality of chain
3	G	4	
4	H	2	

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
8	MPD	B	1361	-	-	X	-
8	MPD	C	1360	-	-	X	-

2 Entry composition [i](#)

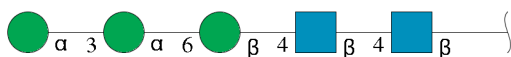
There are 9 unique types of molecules in this entry. The entry contains 12645 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 QUERCETIN 2,3-DIOXYGENASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	344	Total 2657	C 1682	N 431	O 539	S 5	0	5	0
1	B	344	Total 2654	C 1681	N 431	O 537	S 5	0	4	0
1	C	343	Total 2648	C 1678	N 430	O 535	S 5	0	4	0
1	D	343	Total 2648	C 1678	N 430	O 535	S 5	0	5	0

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



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	E	5	Total 61	C 34	N 2	O 25	0	0	0

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



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
3	F	4	Total 50	C 28	N 2	O 20	0	0	0

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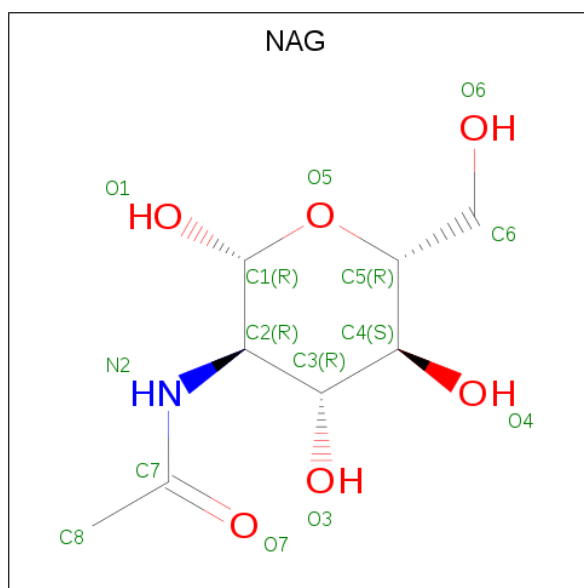
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
3	G	4	50	28	2	20	0	0	0

- 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
			Total	C	N	O			
4	H	2	28	16	2	10	0	0	0

- Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: C₈H₁₅NO₆).



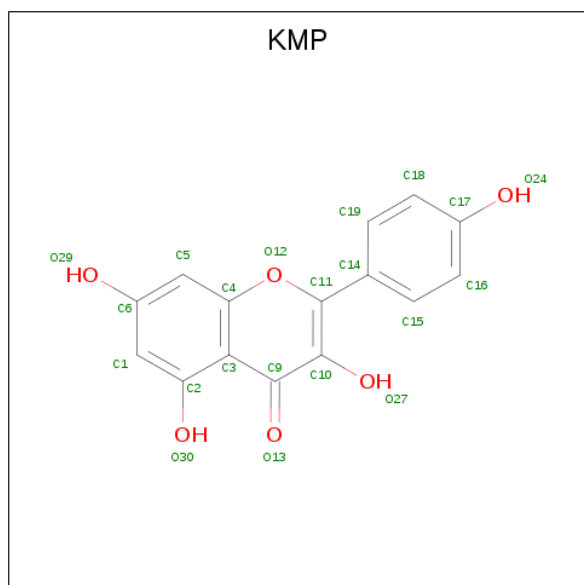
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
5	A	1	14	8	1	5	0	0
5	A	1	14	8	1	5	0	0
5	A	1	14	8	1	5	0	0
5	A	1	14	8	1	5	0	0

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

- Molecule 6 is 3,5,7-TRIHydroxy-2-(4-Hydroxyphenyl)-4H-Chromen-4-one (three-letter code: KMP) (formula: C₁₅H₁₀O₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			21	15	6		
6	B	1	Total	C	O	0	0
			21	15	6		

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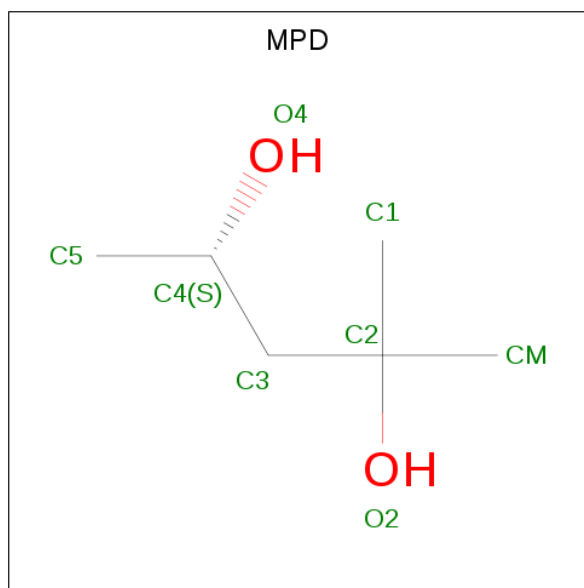
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	C	1	Total	C	O	0	0
			21	15	6		
6	D	1	Total	C	O	0	0
			21	15	6		

- Molecule 7 is COPPER (II) ION (three-letter code: CU) (formula: Cu).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	B	1	Total	Cu	0	0
			1	1		
7	A	1	Total	Cu	0	0
			1	1		
7	D	1	Total	Cu	0	0
			1	1		
7	C	1	Total	Cu	0	0
			1	1		

- Molecule 8 is (4S)-2-METHYL-2,4-PENTANEDIOL (three-letter code: MPD) (formula: C₆H₁₄O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	A	1	Total	C	O	0	0
			8	6	2		
8	B	1	Total	C	O	0	0
			8	6	2		
8	B	1	Total	C	O	0	0
			8	6	2		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	C	1	Total	C	O	0	0
			8	6	2		

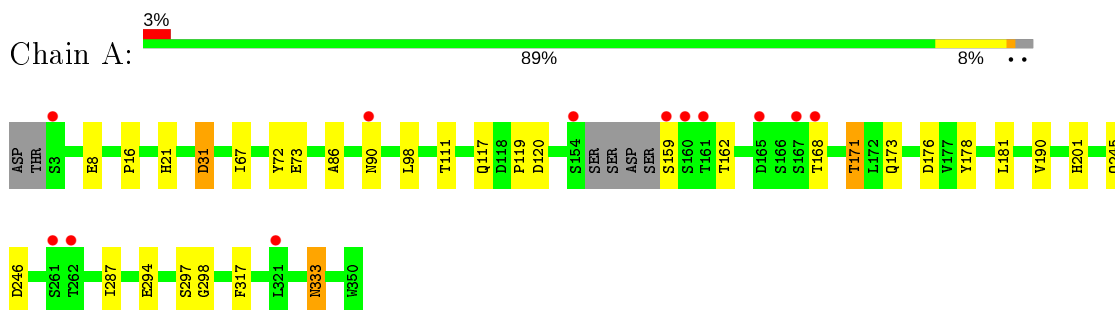
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	416	Total	O	0	0
			416	416		
9	B	387	Total	O	0	0
			387	387		
9	C	381	Total	O	0	0
			381	381		
9	D	363	Total	O	0	0
			363	363		

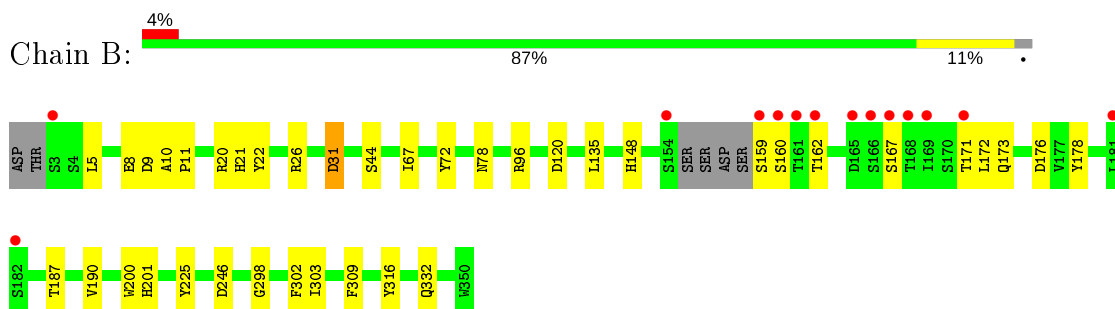
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

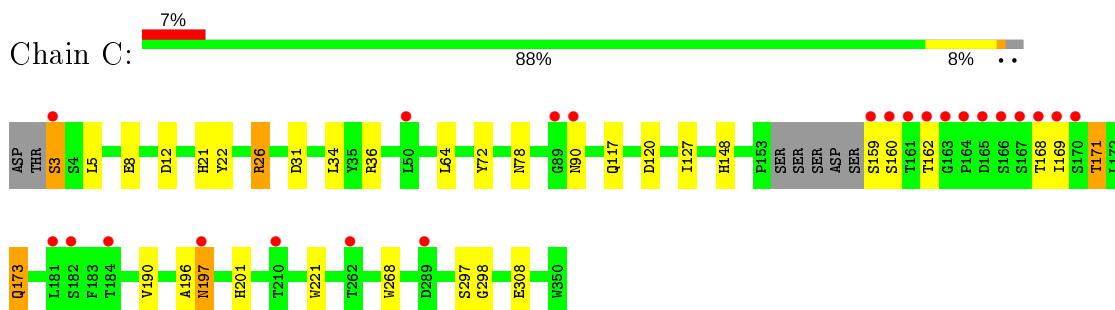
- Molecule 1: QUERCETIN 2,3-DIOXYGENASE



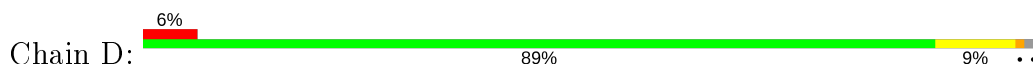
- Molecule 1: QUERCETIN 2,3-DIOXYGENASE

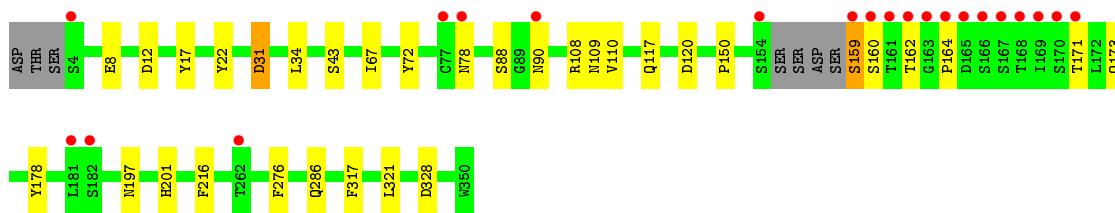


- Molecule 1: QUERCETIN 2,3-DIOXYGENASE



- Molecule 1: QUERCETIN 2,3-DIOXYGENASE





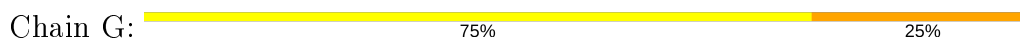
- Molecule 2: alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 3: alpha-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 3: alpha-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



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



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	109.27Å 55.38Å 123.93Å 90.00° 98.33° 90.00°	Depositor
Resolution (Å)	49.39 – 1.90 39.72 – 1.90	Depositor EDS
% Data completeness (in resolution range)	96.5 (49.39-1.90) 96.5 (39.72-1.90)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.59 (at 1.91Å)	Xtrriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.154 , 0.204 0.180 , 0.220	Depositor DCC
R_{free} test set	5608 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	16.5	Xtrriage
Anisotropy	0.571	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 52.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	12645	wwPDB-VP
Average B, all atoms (Å ²)	12.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.38% 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: MPD, BMA, NAG, KMP, CU, MAN

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.95	0/2762	0.94	4/3786 (0.1%)
1	B	0.91	0/2754	0.93	7/3775 (0.2%)
1	C	0.92	0/2748	0.87	2/3767 (0.1%)
1	D	0.90	0/2748	0.88	4/3767 (0.1%)
All	All	0.92	0/11012	0.91	17/15095 (0.1%)

There are no bond length outliers.

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	120	ASP	CB-CG-OD2	8.07	125.56	118.30
1	A	246	ASP	CB-CG-OD2	7.70	125.23	118.30
1	B	9	ASP	CB-CG-OD2	7.60	125.14	118.30
1	A	176	ASP	CB-CG-OD2	7.37	124.93	118.30
1	C	120	ASP	CB-CG-OD2	7.01	124.61	118.30
1	D	31	ASP	CB-CG-OD2	6.62	124.25	118.30
1	A	31	ASP	CB-CG-OD2	6.57	124.21	118.30
1	A	120	ASP	CB-CG-OD2	6.44	124.10	118.30
1	D	120	ASP	CB-CG-OD2	6.36	124.02	118.30
1	B	176	ASP	CB-CG-OD2	5.93	123.64	118.30
1	D	328	ASP	CB-CG-OD2	5.78	123.50	118.30
1	B	20	ARG	NE-CZ-NH2	-5.52	117.54	120.30
1	D	12	ASP	CB-CG-OD2	5.46	123.22	118.30
1	B	31	ASP	CB-CG-OD2	5.46	123.21	118.30
1	C	12	ASP	CB-CG-OD2	5.26	123.04	118.30
1	B	20	ARG	NE-CZ-NH1	5.21	122.91	120.30
1	B	246	ASP	CB-CG-OD2	5.02	122.82	118.30

There are no chirality outliers.

There are no planarity outliers.

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	2657	0	2453	32	0
1	B	2654	0	2454	26	0
1	C	2648	0	2449	26	0
1	D	2648	0	2445	25	0
2	E	61	0	52	1	0
3	F	50	0	43	1	0
3	G	50	0	43	1	0
4	H	28	0	25	1	0
5	A	56	0	52	0	0
5	B	42	0	39	0	0
5	C	42	0	39	4	0
5	D	42	0	39	0	0
6	A	21	0	6	1	0
6	B	21	0	6	1	0
6	C	21	0	6	0	0
6	D	21	0	7	0	0
7	A	1	0	0	0	0
7	B	1	0	0	0	0
7	C	1	0	0	0	0
7	D	1	0	0	0	0
8	A	8	0	14	4	0
8	B	16	0	28	7	0
8	C	8	0	14	6	0
9	A	416	0	0	14	0
9	B	387	0	0	12	0
9	C	381	0	0	6	0
9	D	363	0	0	11	0
All	All	12645	0	10214	115	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 5.

All (115) 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)
8:B:1360:MPD:H32	8:B:1361:MPD:H32	1.26	1.14
1:A:8[A]:GLU:HG2	9:A:2018:HOH:O	1.55	1.05
1:D:117:GLN:HG3	9:D:2141:HOH:O	1.66	0.95
1:D:78[D]:ASN:HB3	9:D:2151:HOH:O	1.66	0.94
8:B:1360:MPD:C3	8:B:1361:MPD:H32	2.02	0.89
1:B:8[A]:GLU:HG2	9:B:2004:HOH:O	1.71	0.89
1:D:78[D]:ASN:CB	9:D:2151:HOH:O	2.20	0.88
1:B:8[A]:GLU:CG	9:B:2004:HOH:O	2.23	0.84
1:D:8[A]:GLU:HG2	9:D:2009:HOH:O	1.81	0.80
8:A:1362:MPD:H32	8:C:1360:MPD:H32	1.65	0.79
9:B:2384:HOH:O	5:C:1355:NAG:H61	1.84	0.77
1:B:8[A]:GLU:OE2	1:B:148:HIS:HE1	1.68	0.77
1:C:78[B]:ASN:ND2	9:C:2117:HOH:O	2.02	0.76
1:B:172:LEU:HA	9:B:2184:HOH:O	1.86	0.74
1:C:8[A]:GLU:HG2	9:C:2009:HOH:O	1.87	0.74
1:C:26:ARG:HH11	1:C:26:ARG:HG3	1.53	0.73
1:D:159:SER:HA	9:D:2182:HOH:O	1.87	0.72
8:A:1362:MPD:C3	8:C:1360:MPD:H32	2.21	0.71
1:D:173:GLN:HG2	1:D:178:TYR:CE2	2.25	0.71
1:B:31:ASP:HB2	1:B:162:THR:HG23	1.73	0.70
8:B:1360:MPD:H12	1:D:201:HIS:CE1	2.26	0.69
1:A:16:PRO:HB3	1:A:287:ILE:HG21	1.74	0.68
1:B:201:HIS:CE1	8:B:1361:MPD:H12	2.29	0.68
1:C:8[A]:GLU:OE2	1:C:148:HIS:HE1	1.76	0.68
1:A:171:THR:CG2	9:A:2215:HOH:O	2.43	0.67
9:B:2384:HOH:O	5:C:1355:NAG:C6	2.43	0.65
1:A:173:GLN:HG3	9:A:2221:HOH:O	1.96	0.63
1:A:201:HIS:HB3	1:C:117:GLN:NE2	2.15	0.62
1:D:31:ASP:HB2	1:D:162:THR:HG23	1.84	0.60
1:A:8[A]:GLU:CG	9:A:2018:HOH:O	2.31	0.59
1:A:201:HIS:CE1	8:A:1362:MPD:H12	2.38	0.59
9:B:2384:HOH:O	5:C:1355:NAG:H4	2.03	0.59
1:A:294:GLU:HG2	9:A:2331:HOH:O	2.03	0.58
1:C:190:VAL:HG11	3:G:1:NAG:H82	1.86	0.58
1:A:117:GLN:HE22	8:C:1360:MPD:HM3	1.69	0.58
1:C:168:THR:O	1:C:171:THR:HB	2.03	0.57
1:C:26:ARG:HH11	1:C:26:ARG:CG	2.18	0.56
1:A:117:GLN:NE2	8:C:1360:MPD:HM3	2.21	0.56
1:D:78[D]:ASN:HB2	9:D:2151:HOH:O	1.97	0.56
1:A:117:GLN:HG3	9:A:2165:HOH:O	2.06	0.55
1:D:108:ARG:O	1:D:109[A]:ASN:HB2	2.05	0.55
8:B:1360:MPD:H4	8:B:1361:MPD:HM1	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:173:GLN:HG2	1:A:178:TYR:CE2	2.42	0.55
1:B:171:THR:O	1:B:171:THR:HG22	2.07	0.54
1:D:171:THR:O	1:D:171:THR:HG22	2.08	0.54
1:C:196:ALA:O	1:C:197:ASN:CB	2.53	0.54
1:B:167:SER:HB3	9:B:2182:HOH:O	2.08	0.54
1:A:90:ASN:HA	9:A:2129:HOH:O	2.09	0.53
1:C:171:THR:HG23	9:C:2202:HOH:O	2.09	0.53
1:A:173:GLN:HG3	9:A:2218:HOH:O	2.09	0.53
6:B:1358:KMP:O27	6:B:1358:KMP:H19	2.09	0.52
1:A:31:ASP:HB2	1:A:162:THR:HG23	1.92	0.52
1:C:5:LEU:HD22	9:C:2309:HOH:O	2.10	0.52
1:A:171:THR:HG22	9:A:2215:HOH:O	2.07	0.52
1:A:168:THR:O	1:A:171:THR:HB	2.10	0.51
1:C:8[A]:GLU:OE2	1:C:148:HIS:CE1	2.61	0.51
1:A:297:SER:HB2	9:A:2111:HOH:O	2.10	0.51
1:C:196:ALA:O	1:C:197:ASN:HB2	2.11	0.51
1:B:332:GLN:NE2	9:B:2346:HOH:O	2.44	0.51
1:B:135:LEU:HD11	1:B:172:LEU:HD13	1.93	0.49
8:B:1360:MPD:H32	8:B:1361:MPD:C3	2.18	0.49
1:D:34:LEU:HD23	1:D:34:LEU:C	2.33	0.49
1:D:173:GLN:NE2	9:D:2195:HOH:O	2.46	0.49
1:A:21:HIS:CG	1:A:298:GLY:HA3	2.49	0.48
1:D:164:PRO:HD2	9:D:2185:HOH:O	2.14	0.48
1:B:8[A]:GLU:OE2	1:B:148:HIS:CE1	2.58	0.48
1:D:8[A]:GLU:CG	9:D:2009:HOH:O	2.48	0.48
1:B:190:VAL:HG11	3:F:1:NAG:H82	1.94	0.48
1:C:31:ASP:HB2	1:C:162:THR:HG23	1.96	0.48
1:D:173:GLN:HG2	1:D:178:TYR:CD2	2.49	0.48
1:D:197:ASN:O	1:D:197:ASN:CG	2.52	0.48
1:A:333[A]:ASN:ND2	9:A:2376:HOH:O	1.96	0.47
1:D:88:SER:HB2	1:D:110:VAL:HG22	1.97	0.47
1:C:173:GLN:H	1:C:173:GLN:HG2	1.45	0.46
1:A:67:ILE:HG12	1:A:178:TYR:HB2	1.96	0.46
1:D:286:GLN:HA	9:D:2293:HOH:O	2.16	0.46
1:B:78[B]:ASN:ND2	9:B:2094:HOH:O	2.48	0.45
1:A:317:PHE:HB3	1:C:22:TYR:CE2	2.52	0.45
8:A:1362:MPD:H31	8:C:1360:MPD:H32	1.99	0.45
1:B:332:GLN:NE2	9:B:2347:HOH:O	2.50	0.45
1:C:90:ASN:ND2	9:C:2132:HOH:O	2.50	0.45
9:B:2384:HOH:O	5:C:1355:NAG:C5	2.65	0.45
1:D:17:TYR:CE2	1:D:43[A]:SER:OG	2.69	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:216:PHE:CD1	1:D:216:PHE:N	2.85	0.45
1:A:245:GLN:HG3	9:A:2293:HOH:O	2.17	0.44
1:A:73:GLU:OE2	6:A:1360:KMP:O27	2.36	0.44
1:C:3:SER:N	9:C:2001:HOH:O	2.50	0.44
1:A:171:THR:HG23	9:A:2216:HOH:O	2.18	0.44
1:A:86:ALA:HA	1:A:111:THR:O	2.17	0.43
1:C:127:ILE:N	1:C:127:ILE:HD12	2.33	0.43
1:D:197:ASN:HB3	9:D:2215:HOH:O	2.18	0.43
1:A:117:GLN:HB2	9:A:2162:HOH:O	2.18	0.43
1:B:187:THR:HB	4:H:2:NAG:O3	2.18	0.43
1:B:22:TYR:CE2	1:D:317:PHE:HB3	2.54	0.43
1:B:5:LEU:HD11	9:B:2295:HOH:O	2.18	0.42
1:B:173:GLN:HG2	1:B:178:TYR:CD2	2.54	0.42
1:A:16:PRO:HB3	1:A:287:ILE:CG2	2.47	0.42
1:A:98:LEU:HD12	1:A:98:LEU:N	2.35	0.42
1:A:201:HIS:HB3	1:C:117:GLN:HE22	1.85	0.42
1:A:119:PRO:HD3	1:C:221:TRP:CE2	2.55	0.41
1:B:316:TYR:HB2	1:D:22:TYR:HB2	2.03	0.41
1:B:10:ALA:HA	1:B:11:PRO:HD3	1.90	0.41
1:B:225:TYR:N	1:B:225:TYR:CD1	2.89	0.41
1:C:201:HIS:CE1	8:C:1360:MPD:H12	2.55	0.41
1:D:276:PHE:HA	1:D:321:LEU:O	2.21	0.41
1:B:201:HIS:ND1	8:B:1361:MPD:H12	2.36	0.41
1:B:21:HIS:CG	1:B:298:GLY:HA3	2.56	0.41
1:C:268:TRP:O	1:C:308:GLU:HA	2.21	0.41
1:C:64:LEU:HB2	1:C:169:ILE:HD13	2.03	0.40
1:B:303:ILE:HG21	1:B:309:PHE:CD1	2.57	0.40
1:C:21:HIS:CG	1:C:298:GLY:HA3	2.56	0.40
1:C:34:LEU:HD22	1:C:36:ARG:NH1	2.37	0.40
1:B:44:SER:HA	1:B:302:PHE:CZ	2.56	0.40
1:B:96:ARG:HD2	1:B:200:TRP:CD2	2.56	0.40
1:A:190:VAL:HG11	2:E:1:NAG:H82	2.03	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	345/350 (99%)	333 (96%)	12 (4%)	0	100	100
1	B	344/350 (98%)	332 (96%)	12 (4%)	0	100	100
1	C	343/350 (98%)	334 (97%)	9 (3%)	0	100	100
1	D	343/350 (98%)	334 (97%)	9 (3%)	0	100	100
All	All	1375/1400 (98%)	1333 (97%)	42 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	293/294 (100%)	287 (98%)	6 (2%)	55	51
1	B	292/294 (99%)	287 (98%)	5 (2%)	60	57
1	C	291/294 (99%)	282 (97%)	9 (3%)	40	32
1	D	291/294 (99%)	285 (98%)	6 (2%)	53	48
All	All	1167/1176 (99%)	1141 (98%)	26 (2%)	52	47

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	72	TYR
1	A	159	SER
1	A	171	THR
1	A	181	LEU
1	A	333[A]	ASN
1	A	333[B]	ASN
1	B	26	ARG
1	B	67	ILE

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Mol	Chain	Res	Type
1	B	72	TYR
1	B	159	SER
1	B	160	SER
1	C	3	SER
1	C	26	ARG
1	C	72	TYR
1	C	159	SER
1	C	160	SER
1	C	171	THR
1	C	173	GLN
1	C	197	ASN
1	C	297	SER
1	D	67	ILE
1	D	72	TYR
1	D	90	ASN
1	D	150	PRO
1	D	159	SER
1	D	160	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	83	GLN
1	A	117	GLN
1	A	332	GLN
1	B	83	GLN
1	B	117	GLN
1	B	148	HIS
1	B	286	GLN
1	B	332	GLN
1	C	83	GLN
1	C	117	GLN
1	C	148	HIS
1	C	336	ASN
1	D	83	GLN
1	D	90	ASN
1	D	173	GLN
1	D	336	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

15 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	E	1	1,2	14,14,15	0.86	1 (7%)	17,19,21	1.24	2 (11%)
2	NAG	E	2	2	14,14,15	0.72	0	17,19,21	1.19	2 (11%)
2	BMA	E	3	2	11,11,12	0.95	0	15,15,17	0.95	1 (6%)
2	MAN	E	4	2	11,11,12	0.59	0	15,15,17	1.42	2 (13%)
2	MAN	E	5	2	11,11,12	0.78	0	15,15,17	0.85	0
3	NAG	F	1	1,3	14,14,15	0.87	1 (7%)	17,19,21	1.50	1 (5%)
3	NAG	F	2	3	14,14,15	0.62	0	17,19,21	1.93	6 (35%)
3	BMA	F	3	3	11,11,12	0.71	0	15,15,17	1.72	3 (20%)
3	MAN	F	4	3	11,11,12	0.85	0	15,15,17	1.12	0
3	NAG	G	1	1,3	14,14,15	0.46	0	17,19,21	1.66	5 (29%)
3	NAG	G	2	3	14,14,15	0.50	0	17,19,21	1.51	4 (23%)
3	BMA	G	3	3	11,11,12	1.10	1 (9%)	15,15,17	0.97	0
3	MAN	G	4	3	11,11,12	0.68	0	15,15,17	1.23	1 (6%)
4	NAG	H	1	1,4	14,14,15	0.74	1 (7%)	17,19,21	0.98	1 (5%)
4	NAG	H	2	4	14,14,15	0.86	0	17,19,21	1.79	4 (23%)

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	E	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	E	2	2	-	0/6/23/26	0/1/1/1
2	BMA	E	3	2	-	0/2/19/22	0/1/1/1
2	MAN	E	4	2	-	2/2/19/22	0/1/1/1
2	MAN	E	5	2	-	2/2/19/22	0/1/1/1
3	NAG	F	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	F	2	3	-	0/6/23/26	0/1/1/1
3	BMA	F	3	3	-	0/2/19/22	0/1/1/1
3	MAN	F	4	3	-	0/2/19/22	0/1/1/1
3	NAG	G	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	G	2	3	-	0/6/23/26	0/1/1/1
3	BMA	G	3	3	-	0/2/19/22	0/1/1/1
3	MAN	G	4	3	-	0/2/19/22	0/1/1/1
4	NAG	H	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	H	2	4	-	0/6/23/26	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	G	3	BMA	O5-C1	-2.98	1.39	1.43
2	E	1	NAG	C1-C2	2.22	1.55	1.52
4	H	1	NAG	C1-C2	2.22	1.55	1.52
3	F	1	NAG	C1-C2	2.00	1.55	1.52

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	1	NAG	C1-O5-C5	5.32	119.40	112.19
3	F	2	NAG	C1-O5-C5	4.49	118.28	112.19
3	F	3	BMA	C1-O5-C5	4.19	117.86	112.19
2	E	4	MAN	C1-O5-C5	4.13	117.78	112.19
4	H	2	NAG	O7-C7-C8	-4.12	114.41	122.06
3	G	4	MAN	O5-C5-C6	3.69	112.98	107.20
3	G	2	NAG	O5-C1-C2	-3.63	105.56	111.29
3	G	1	NAG	O5-C5-C6	3.23	112.26	107.20
4	H	2	NAG	C1-C2-N2	3.02	115.65	110.49
4	H	2	NAG	C8-C7-N2	2.98	121.15	116.10
4	H	2	NAG	C1-O5-C5	2.91	116.14	112.19
3	G	1	NAG	O5-C1-C2	-2.89	106.72	111.29
2	E	2	NAG	C1-O5-C5	2.88	116.10	112.19
3	F	2	NAG	O6-C6-C5	-2.82	101.61	111.29

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	1	NAG	O5-C1-C2	-2.75	106.94	111.29
3	F	2	NAG	C3-C4-C5	2.63	114.93	110.24
3	F	2	NAG	C8-C7-N2	-2.61	111.69	116.10
2	E	4	MAN	O5-C5-C6	2.52	111.15	107.20
3	F	3	BMA	O5-C5-C6	2.36	110.91	107.20
3	G	1	NAG	O7-C7-C8	-2.36	117.68	122.06
3	G	2	NAG	O4-C4-C3	-2.26	105.12	110.35
3	F	3	BMA	C3-C4-C5	2.25	114.25	110.24
3	G	2	NAG	C1-C2-N2	2.23	114.30	110.49
3	F	2	NAG	O7-C7-N2	2.20	126.00	121.95
3	F	2	NAG	C6-C5-C4	-2.20	107.85	113.00
4	H	1	NAG	C8-C7-N2	2.20	119.82	116.10
3	G	2	NAG	O7-C7-N2	2.19	125.98	121.95
3	G	1	NAG	O6-C6-C5	-2.18	103.81	111.29
3	G	1	NAG	C4-C3-C2	2.18	114.21	111.02
2	E	3	BMA	O3-C3-C2	2.05	113.92	109.99
2	E	1	NAG	C4-C3-C2	-2.01	108.07	111.02
2	E	2	NAG	O5-C1-C2	-2.00	108.13	111.29

There are no chirality outliers.

All (4) torsion outliers are listed below:

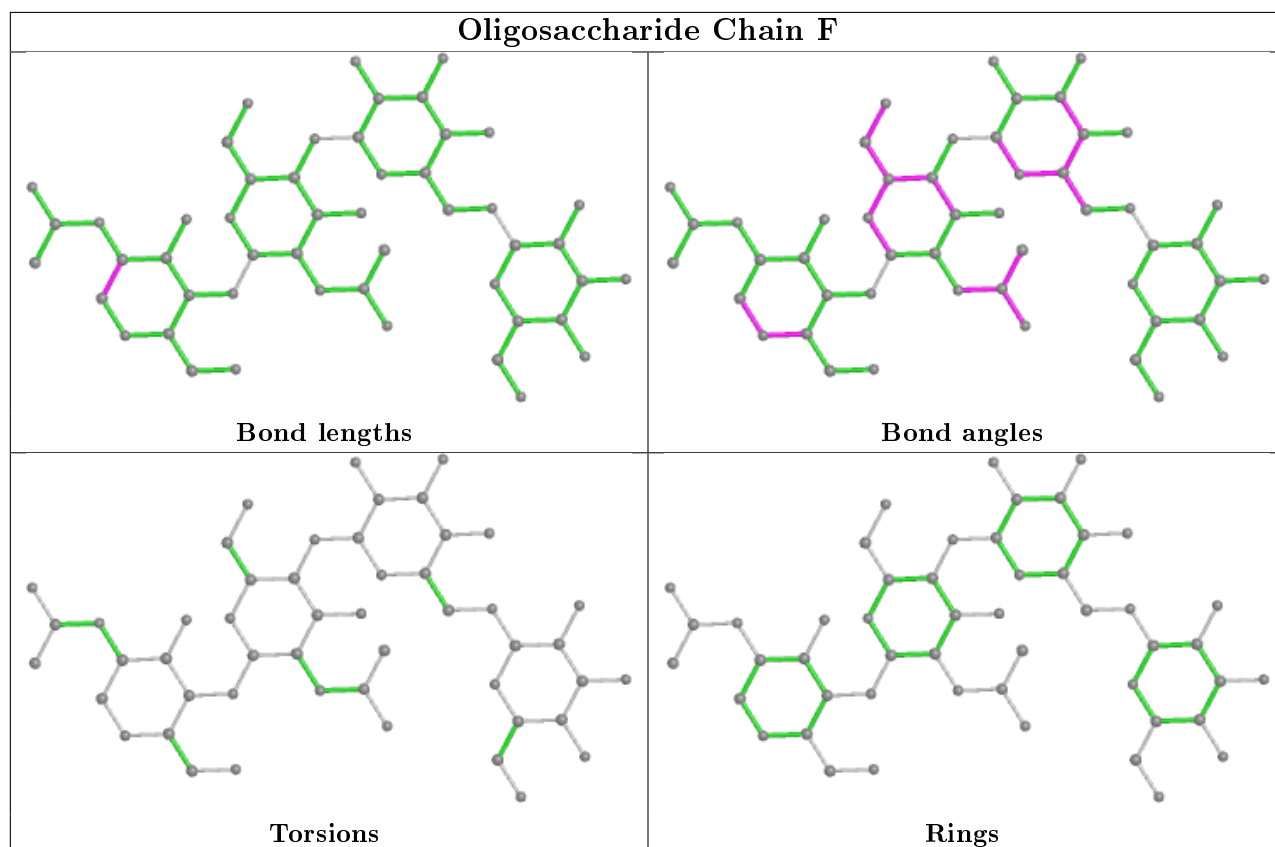
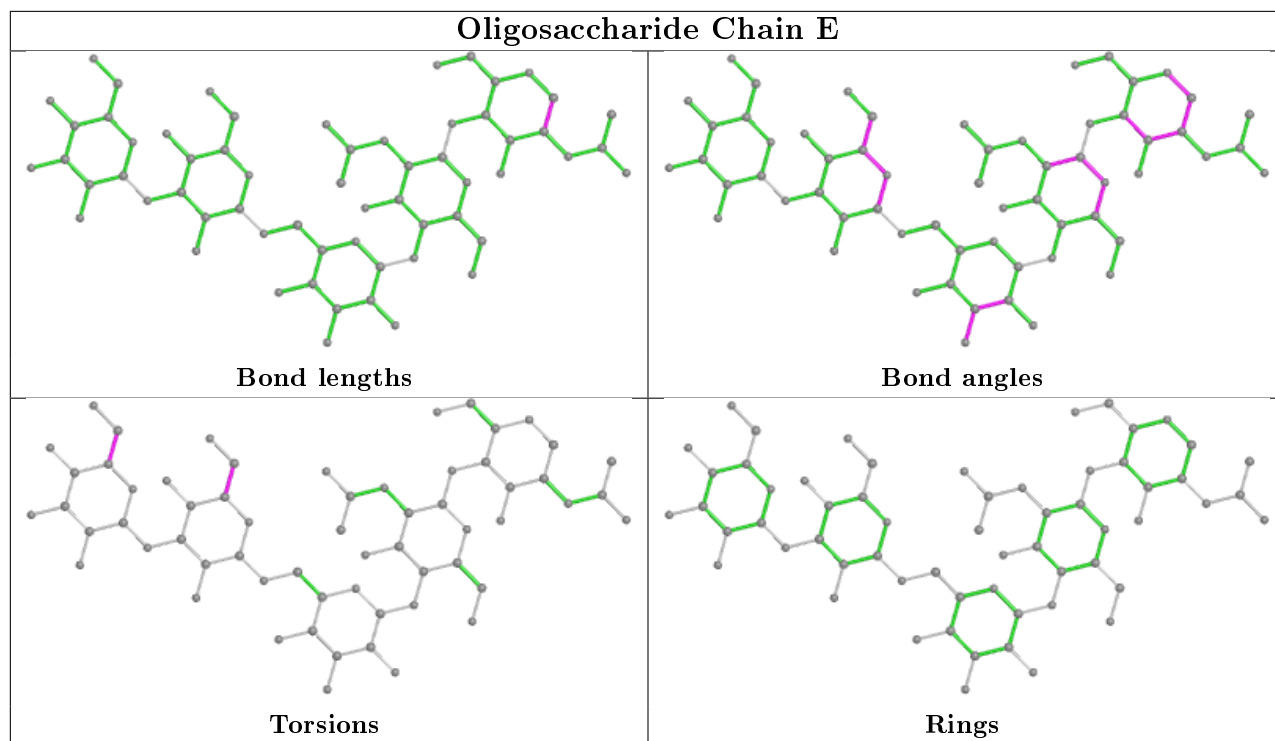
Mol	Chain	Res	Type	Atoms
2	E	5	MAN	O5-C5-C6-O6
2	E	5	MAN	C4-C5-C6-O6
2	E	4	MAN	C4-C5-C6-O6
2	E	4	MAN	O5-C5-C6-O6

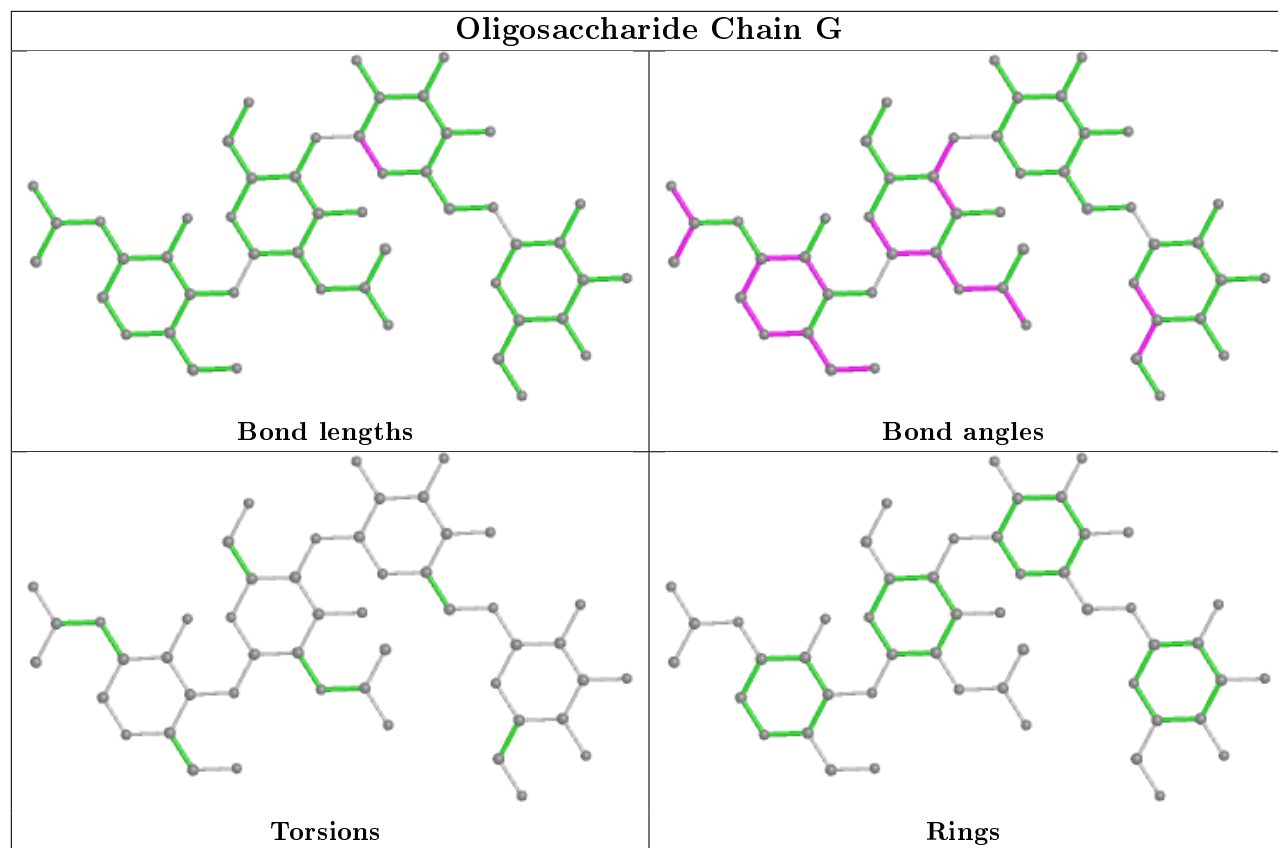
There are no ring outliers.

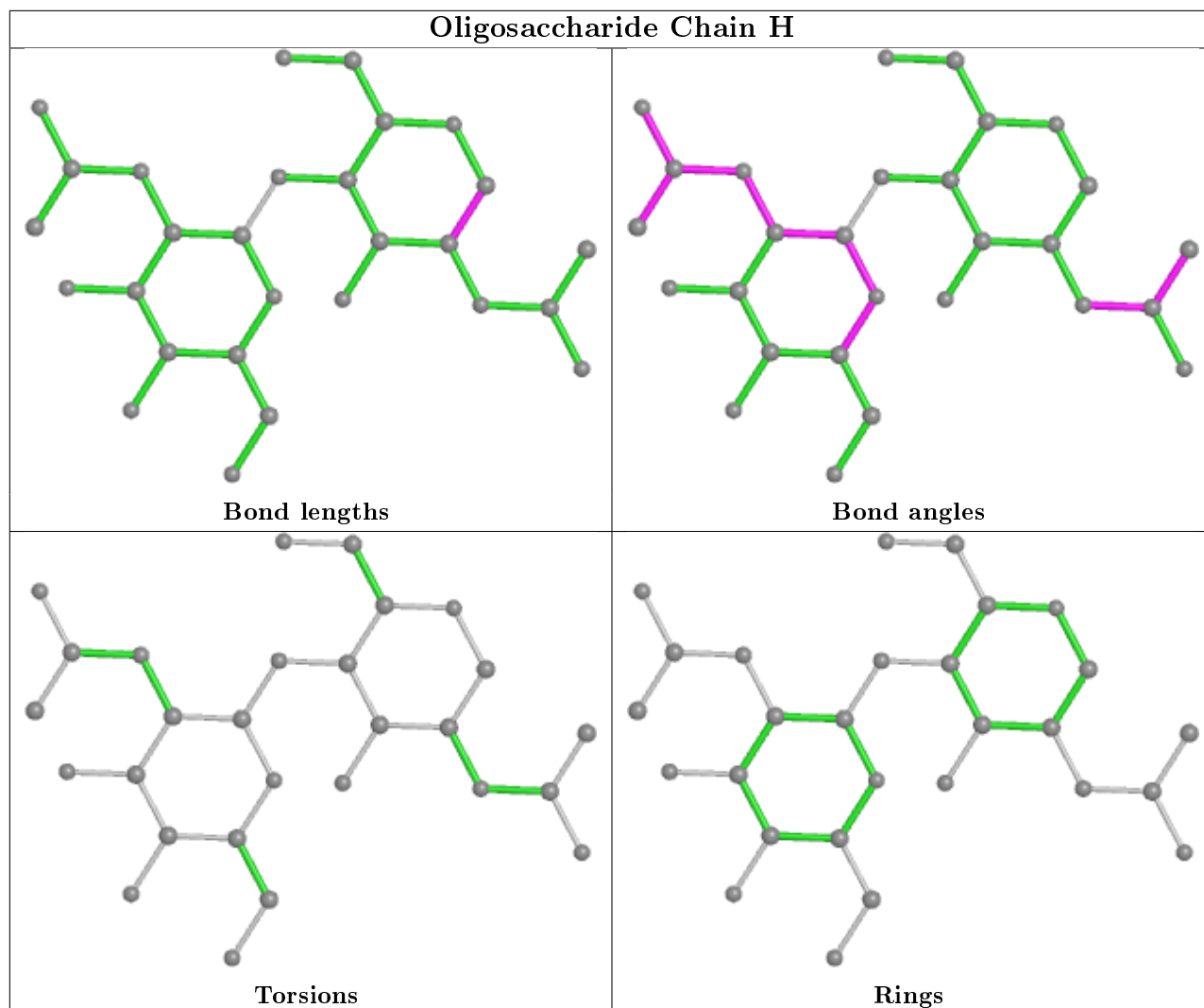
4 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	H	2	NAG	1	0
3	F	1	NAG	1	0
2	E	1	NAG	1	0
3	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 25 ligands modelled in this entry, 4 are monoatomic - leaving 21 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
8	MPD	C	1360	-	7,7,7	0.66	0	9,10,10	1.52	1 (11%)
5	NAG	B	1355	1	14,14,15	0.77	0	17,19,21	1.17	2 (11%)
5	NAG	B	1351	1	14,14,15	0.49	0	17,19,21	1.01	1 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	KMP	D	1356	7	20,23,23	1.94	3 (15%)	26,34,34	1.85	8 (30%)
6	KMP	A	1360	7	20,23,23	1.83	5 (25%)	26,34,34	1.92	5 (19%)
5	NAG	D	1355	1	14,14,15	0.82	1 (7%)	17,19,21	1.63	3 (17%)
5	NAG	A	1356	1	14,14,15	0.94	0	17,19,21	1.50	1 (5%)
8	MPD	B	1360	-	7,7,7	0.65	0	9,10,10	2.05	3 (33%)
6	KMP	B	1358	7	20,23,23	1.64	5 (25%)	26,34,34	1.60	8 (30%)
8	MPD	A	1362	-	7,7,7	0.74	0	9,10,10	1.92	1 (11%)
5	NAG	C	1355	1	14,14,15	0.76	0	17,19,21	2.06	6 (35%)
5	NAG	A	1355	1	14,14,15	0.98	1 (7%)	17,19,21	1.50	3 (17%)
5	NAG	C	1351	1	14,14,15	0.53	0	17,19,21	1.08	0
5	NAG	A	1354	1	14,14,15	0.72	0	17,19,21	1.50	3 (17%)
5	NAG	D	1354	1	14,14,15	1.07	1 (7%)	17,19,21	1.91	2 (11%)
8	MPD	B	1361	-	7,7,7	0.71	0	9,10,10	1.93	2 (22%)
5	NAG	B	1354	1	14,14,15	0.73	0	17,19,21	1.31	3 (17%)
5	NAG	D	1351	1	14,14,15	0.79	0	17,19,21	1.46	3 (17%)
5	NAG	C	1354	1	14,14,15	0.90	1 (7%)	17,19,21	1.44	3 (17%)
5	NAG	A	1351	1	14,14,15	0.65	0	17,19,21	1.04	1 (5%)
6	KMP	C	1358	7	20,23,23	1.89	3 (15%)	26,34,34	1.79	7 (26%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
8	MPD	C	1360	-	-	2/5/5/5	-
5	NAG	B	1355	1	-	0/6/23/26	0/1/1/1
5	NAG	B	1351	1	-	0/6/23/26	0/1/1/1
6	KMP	D	1356	7	-	2/4/4/4	0/3/3/3
6	KMP	A	1360	7	-	2/4/4/4	0/3/3/3
5	NAG	D	1355	1	-	2/6/23/26	0/1/1/1
5	NAG	A	1356	1	-	5/6/23/26	0/1/1/1
8	MPD	B	1360	-	-	1/5/5/5	-
6	KMP	B	1358	7	-	2/4/4/4	0/3/3/3
8	MPD	A	1362	-	-	1/5/5/5	-
5	NAG	C	1355	1	-	1/6/23/26	0/1/1/1
5	NAG	A	1355	1	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	NAG	C	1351	1	-	0/6/23/26	0/1/1/1
5	NAG	A	1354	1	-	0/6/23/26	0/1/1/1
5	NAG	D	1354	1	-	3/6/23/26	0/1/1/1
8	MPD	B	1361	-	-	2/5/5/5	-
5	NAG	B	1354	1	-	2/6/23/26	0/1/1/1
5	NAG	D	1351	1	-	4/6/23/26	0/1/1/1
5	NAG	C	1354	1	-	2/6/23/26	0/1/1/1
5	NAG	A	1351	1	-	0/6/23/26	0/1/1/1
6	KMP	C	1358	7	-	2/4/4/4	0/3/3/3

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	D	1356	KMP	C3-C4	5.63	1.49	1.41
6	C	1358	KMP	C3-C4	5.49	1.48	1.41
6	B	1358	KMP	C3-C4	4.69	1.47	1.41
6	A	1360	KMP	C3-C4	4.43	1.47	1.41
6	A	1360	KMP	C14-C11	3.80	1.52	1.46
6	C	1358	KMP	C14-C11	3.76	1.52	1.46
6	D	1356	KMP	C14-C11	3.39	1.52	1.46
6	A	1360	KMP	C2-C3	2.94	1.48	1.43
5	D	1354	NAG	C1-C2	2.78	1.56	1.52
6	B	1358	KMP	C1-C6	2.74	1.43	1.39
5	C	1354	NAG	C1-C2	2.73	1.56	1.52
6	D	1356	KMP	C2-C3	2.71	1.48	1.43
6	B	1358	KMP	C2-C3	2.68	1.48	1.43
5	D	1355	NAG	C1-C2	2.55	1.56	1.52
6	C	1358	KMP	C2-C3	2.52	1.47	1.43
5	A	1355	NAG	C1-C2	2.40	1.55	1.52
6	A	1360	KMP	O24-C17	-2.20	1.31	1.37
6	B	1358	KMP	O29-C6	-2.16	1.32	1.37
6	A	1360	KMP	C1-C2	2.16	1.42	1.37
6	B	1358	KMP	C14-C11	2.14	1.50	1.46

All (66) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	1360	KMP	O12-C4-C5	4.97	121.93	116.11
8	A	1362	MPD	CM-C2-C1	-4.93	100.29	110.57
5	D	1354	NAG	O5-C1-C2	-4.93	103.51	111.29
5	A	1356	NAG	C1-O5-C5	4.86	118.78	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	B	1361	MPD	CM-C2-C1	-4.71	100.77	110.57
5	C	1355	NAG	C1-O5-C5	-4.70	105.82	112.19
6	D	1356	KMP	O12-C4-C5	4.65	121.56	116.11
6	A	1360	KMP	C5-C4-C3	-4.51	118.06	123.05
8	B	1360	MPD	CM-C2-C1	-4.44	101.33	110.57
5	D	1354	NAG	C1-O5-C5	4.43	118.20	112.19
6	C	1358	KMP	C19-C14-C11	3.78	125.56	120.29
8	C	1360	MPD	CM-C2-C1	-3.68	102.90	110.57
6	C	1358	KMP	O12-C4-C5	3.67	120.40	116.11
5	D	1351	NAG	C1-O5-C5	3.67	117.16	112.19
5	D	1355	NAG	O5-C1-C2	-3.57	105.65	111.29
5	A	1354	NAG	C2-N2-C7	-3.54	117.86	122.90
5	D	1355	NAG	C2-N2-C7	-3.53	117.88	122.90
6	C	1358	KMP	C1-C2-C3	-3.47	116.58	120.61
6	A	1360	KMP	C15-C16-C17	-3.47	116.07	119.88
5	A	1355	NAG	C1-O5-C5	3.44	116.85	112.19
5	C	1355	NAG	C6-C5-C4	3.40	120.97	113.00
6	B	1358	KMP	C1-C2-C3	-3.39	116.68	120.61
6	D	1356	KMP	C10-C9-C3	-3.38	116.65	121.38
5	B	1355	NAG	C2-N2-C7	3.24	127.51	122.90
6	C	1358	KMP	C2-C1-C6	3.23	122.62	119.70
5	D	1355	NAG	C1-O5-C5	3.23	116.57	112.19
5	A	1354	NAG	O5-C1-C2	-3.20	106.24	111.29
5	C	1355	NAG	O4-C4-C3	-3.11	103.16	110.35
6	C	1358	KMP	C15-C14-C11	-3.09	115.98	120.29
6	B	1358	KMP	O12-C4-C5	3.01	119.63	116.11
6	D	1356	KMP	C19-C18-C17	2.91	123.07	119.88
6	D	1356	KMP	C5-C6-C1	2.89	123.27	120.94
5	C	1354	NAG	C2-N2-C7	-2.89	118.79	122.90
5	C	1354	NAG	O5-C5-C6	2.87	111.71	107.20
5	B	1354	NAG	O5-C5-C6	2.78	111.56	107.20
6	D	1356	KMP	C5-C4-C3	-2.75	120.01	123.05
5	A	1355	NAG	O7-C7-C8	-2.73	116.98	122.06
6	B	1358	KMP	O12-C4-C3	-2.71	118.36	121.03
8	B	1360	MPD	O2-C2-CM	-2.65	99.57	108.08
6	D	1356	KMP	O12-C4-C3	-2.64	118.43	121.03
6	C	1358	KMP	C10-C9-C3	-2.54	117.82	121.38
8	B	1361	MPD	O2-C2-C1	2.54	116.23	108.08
6	D	1356	KMP	C19-C14-C11	2.52	123.81	120.29
5	C	1355	NAG	O5-C5-C4	-2.51	104.72	110.83
8	B	1360	MPD	O2-C2-C1	2.48	116.04	108.08
5	C	1355	NAG	C2-N2-C7	-2.42	119.46	122.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	1351	NAG	C3-C4-C5	-2.41	105.94	110.24
6	B	1358	KMP	C15-C14-C11	-2.40	116.94	120.29
6	C	1358	KMP	C5-C4-C3	-2.39	120.41	123.05
6	B	1358	KMP	C5-C6-C1	2.35	122.84	120.94
6	B	1358	KMP	C19-C14-C11	2.33	123.55	120.29
5	C	1354	NAG	O5-C1-C2	-2.29	107.67	111.29
5	B	1354	NAG	O5-C1-C2	-2.26	107.73	111.29
5	B	1354	NAG	C2-N2-C7	-2.21	119.75	122.90
6	D	1356	KMP	C15-C14-C11	-2.15	117.29	120.29
5	C	1355	NAG	O5-C5-C6	2.10	110.50	107.20
6	B	1358	KMP	C2-C1-C6	2.10	121.60	119.70
5	D	1351	NAG	C2-N2-C7	-2.09	119.93	122.90
5	A	1355	NAG	O5-C1-C2	-2.08	108.00	111.29
6	B	1358	KMP	C10-C9-C3	-2.07	118.48	121.38
5	A	1351	NAG	O6-C6-C5	-2.07	104.20	111.29
5	A	1354	NAG	O5-C5-C4	-2.06	105.81	110.83
5	D	1351	NAG	O7-C7-C8	-2.06	118.23	122.06
5	B	1355	NAG	O5-C1-C2	-2.06	108.04	111.29
6	A	1360	KMP	O29-C6-C5	-2.05	115.11	120.98
6	A	1360	KMP	C18-C19-C14	-2.03	118.21	121.13

There are no chirality outliers.

All (35) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	1360	KMP	C10-C11-C14-C15
6	A	1360	KMP	C10-C11-C14-C19
8	B	1361	MPD	C2-C3-C4-C5
6	D	1356	KMP	C10-C11-C14-C15
6	D	1356	KMP	C10-C11-C14-C19
6	B	1358	KMP	C10-C11-C14-C15
6	B	1358	KMP	C10-C11-C14-C19
6	C	1358	KMP	C10-C11-C14-C15
6	C	1358	KMP	C10-C11-C14-C19
5	D	1351	NAG	O5-C5-C6-O6
5	A	1355	NAG	C4-C5-C6-O6
5	D	1351	NAG	C4-C5-C6-O6
5	D	1351	NAG	C8-C7-N2-C2
5	D	1351	NAG	O7-C7-N2-C2
5	A	1355	NAG	O5-C5-C6-O6
5	A	1356	NAG	O5-C5-C6-O6
5	C	1354	NAG	C4-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
5	C	1354	NAG	O5-C5-C6-O6
5	A	1356	NAG	C8-C7-N2-C2
5	A	1356	NAG	C3-C2-N2-C7
5	C	1355	NAG	C4-C5-C6-O6
5	A	1356	NAG	O7-C7-N2-C2
5	A	1356	NAG	C4-C5-C6-O6
8	C	1360	MPD	C2-C3-C4-C5
5	B	1354	NAG	C4-C5-C6-O6
5	D	1354	NAG	O7-C7-N2-C2
8	B	1361	MPD	C1-C2-C3-C4
8	C	1360	MPD	C1-C2-C3-C4
5	B	1354	NAG	O5-C5-C6-O6
5	D	1354	NAG	C3-C2-N2-C7
5	D	1355	NAG	C8-C7-N2-C2
5	D	1354	NAG	C8-C7-N2-C2
8	B	1360	MPD	C2-C3-C4-C5
8	A	1362	MPD	C2-C3-C4-C5
5	D	1355	NAG	O7-C7-N2-C2

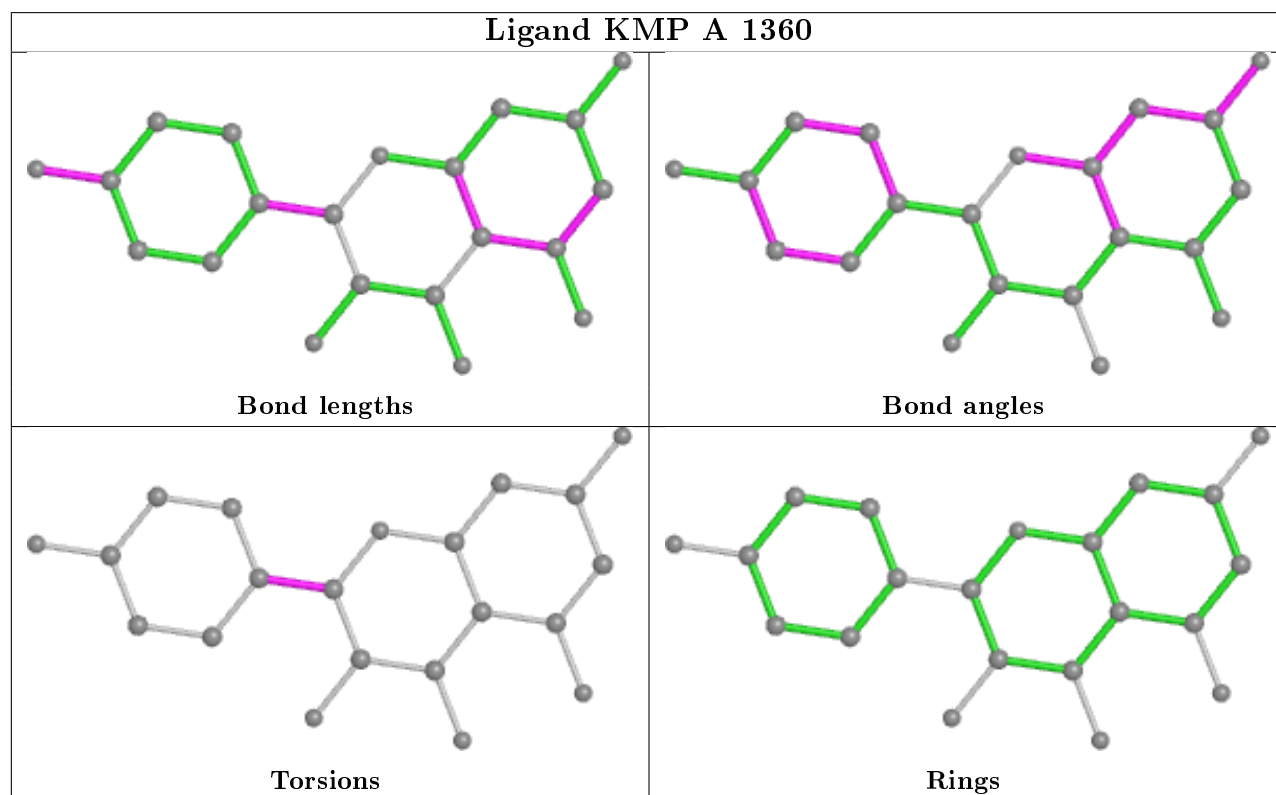
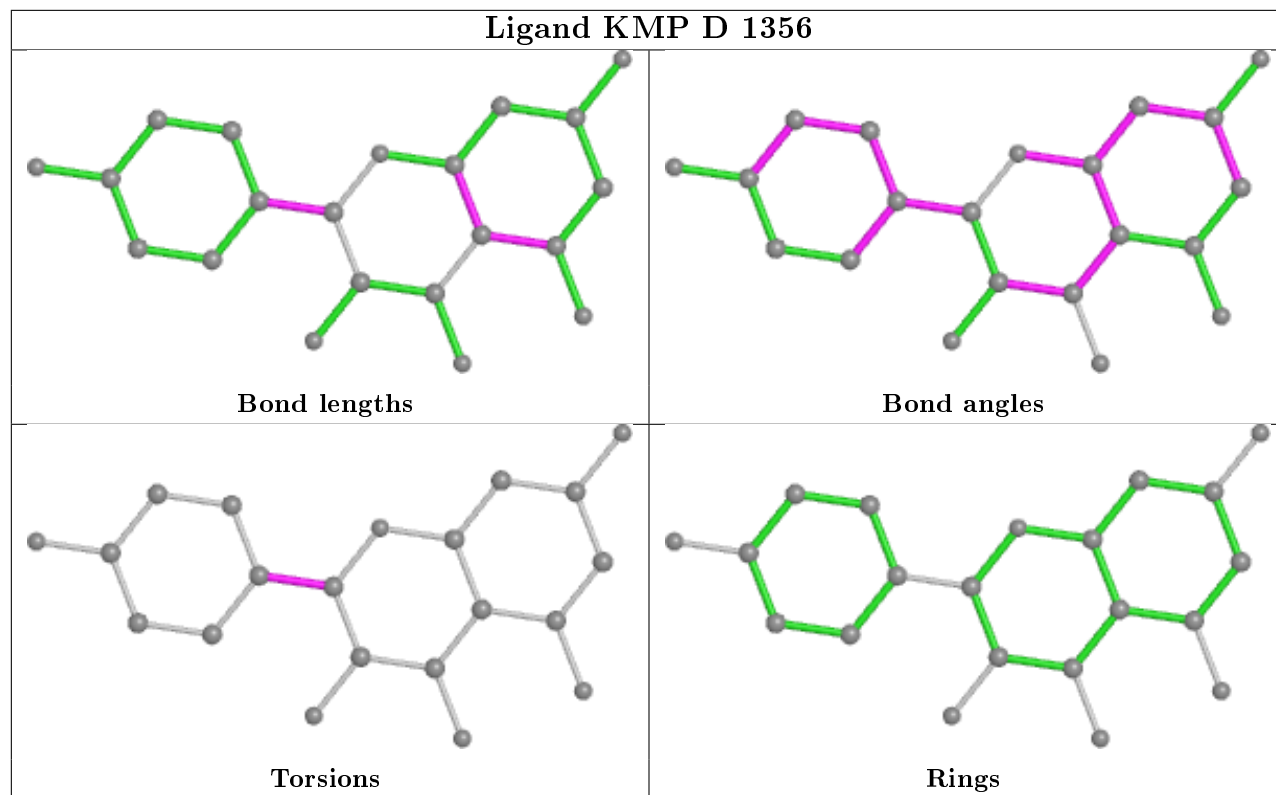
There are no ring outliers.

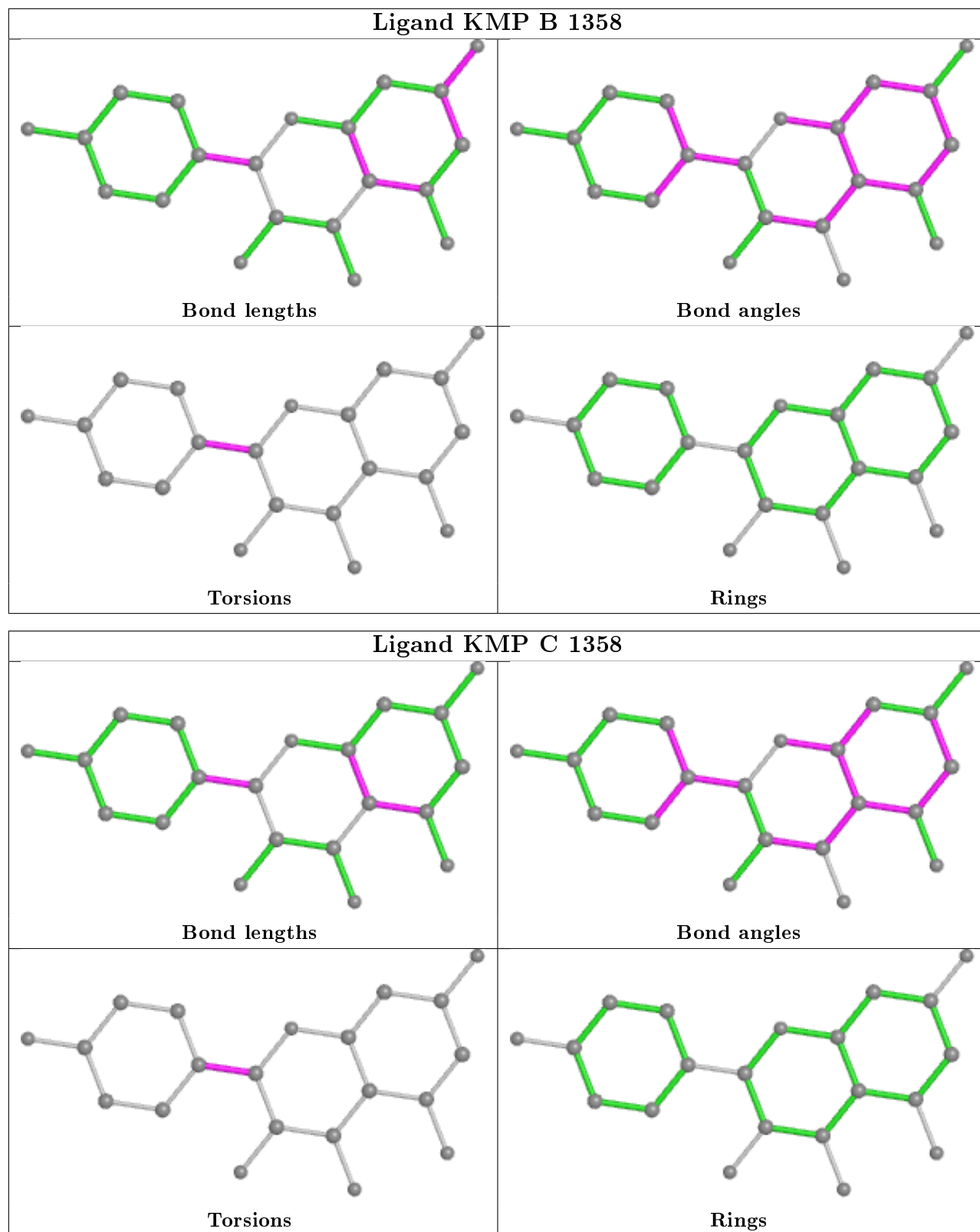
7 monomers are involved in 20 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	C	1360	MPD	6	0
6	A	1360	KMP	1	0
8	B	1360	MPD	5	0
6	B	1358	KMP	1	0
8	A	1362	MPD	4	0
5	C	1355	NAG	4	0
8	B	1361	MPD	6	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

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	344/350 (98%)	-0.13	12 (3%) 44 47	4, 9, 16, 30	0
1	B	344/350 (98%)	-0.15	14 (4%) 37 40	3, 9, 17, 31	0
1	C	343/350 (98%)	0.13	23 (6%) 17 20	3, 9, 17, 31	0
1	D	343/350 (98%)	0.08	21 (6%) 21 24	4, 9, 16, 31	1 (0%)
All	All	1374/1400 (98%)	-0.02	70 (5%) 28 31	3, 9, 17, 31	1 (0%)

All (70) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	169	ILE	7.2
1	C	159	SER	5.9
1	B	3	SER	5.8
1	D	159	SER	5.4
1	C	160	SER	5.2
1	D	160	SER	5.2
1	C	169	ILE	5.1
1	C	166	SER	4.6
1	B	159	SER	4.5
1	C	197	ASN	4.4
1	D	181	LEU	4.4
1	D	166	SER	4.4
1	B	181	LEU	4.3
1	A	90	ASN	4.1
1	A	160	SER	4.1
1	B	160	SER	4.0
1	A	159	SER	3.9
1	B	167	SER	3.9
1	C	163	GLY	3.7
1	B	168	THR	3.7
1	C	167	SER	3.6

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Mol	Chain	Res	Type	RSRZ
1	C	90	ASN	3.4
1	C	181	LEU	3.1
1	D	182	SER	3.1
1	B	161	THR	3.1
1	C	165	ASP	3.1
1	C	164	PRO	3.0
1	D	162	THR	3.0
1	D	78[D]	ASN	3.0
1	A	3	SER	3.0
1	D	167	SER	3.0
1	D	4	SER	2.9
1	C	289	ASP	2.9
1	D	262	THR	2.9
1	A	261	SER	2.8
1	C	170	SER	2.8
1	C	168	THR	2.8
1	C	161	THR	2.7
1	A	168	THR	2.7
1	A	167	SER	2.7
1	D	168	THR	2.7
1	C	182	SER	2.7
1	D	169	ILE	2.6
1	C	210	THR	2.6
1	C	3	SER	2.6
1	A	262	THR	2.5
1	B	182	SER	2.5
1	C	162	THR	2.4
1	D	171	THR	2.4
1	D	170	SER	2.4
1	B	171	THR	2.4
1	B	154	SER	2.4
1	C	262	THR	2.4
1	D	90	ASN	2.4
1	B	165	ASP	2.4
1	D	165	ASP	2.4
1	A	154	SER	2.3
1	B	162	THR	2.3
1	A	165	ASP	2.2
1	B	166	SER	2.2
1	C	89	GLY	2.2
1	C	184	THR	2.2
1	D	161	THR	2.2

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Mol	Chain	Res	Type	RSRZ
1	D	77	CYS	2.2
1	D	154	SER	2.2
1	D	163	GLY	2.1
1	A	161	THR	2.1
1	A	321	LEU	2.1
1	D	164	PRO	2.1
1	C	50	LEU	2.0

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

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

6.3 Carbohydrates [i](#)

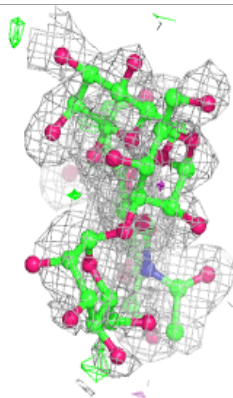
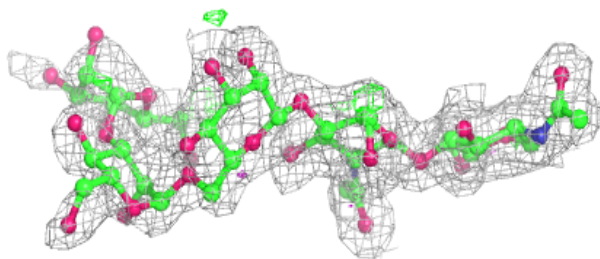
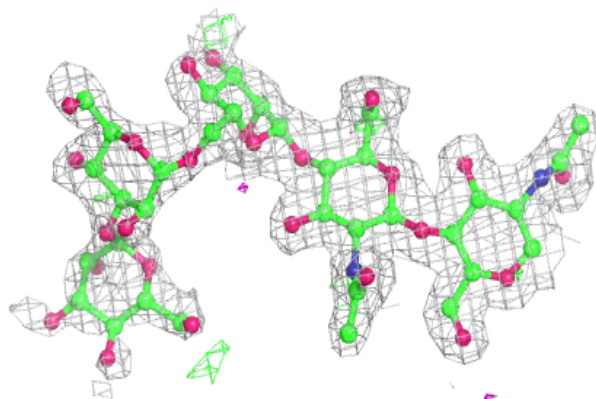
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	MAN	F	4	11/12	0.57	0.27	61,62,64,65	0
3	MAN	G	4	11/12	0.61	0.20	43,46,48,49	0
3	BMA	F	3	11/12	0.65	0.22	46,48,54,59	0
2	MAN	E	5	11/12	0.73	0.31	67,68,68,69	0
2	MAN	E	4	11/12	0.79	0.26	60,61,62,65	0
3	NAG	F	2	14/15	0.82	0.15	24,32,40,44	0
3	BMA	G	3	11/12	0.86	0.15	35,41,43,44	0
2	BMA	E	3	11/12	0.86	0.14	45,48,54,58	0
4	NAG	H	2	14/15	0.93	0.10	18,29,34,35	0
3	NAG	F	1	14/15	0.94	0.10	21,24,29,31	0
3	NAG	G	2	14/15	0.94	0.08	21,25,30,33	0
2	NAG	E	2	14/15	0.94	0.09	24,29,35,39	0
3	NAG	G	1	14/15	0.95	0.08	19,23,34,36	0
2	NAG	E	1	14/15	0.95	0.09	19,25,28,30	0
4	NAG	H	1	14/15	0.96	0.09	20,25,31,33	0

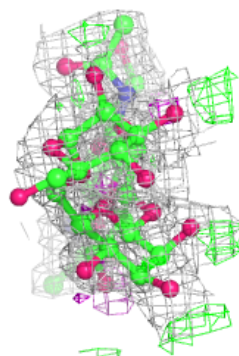
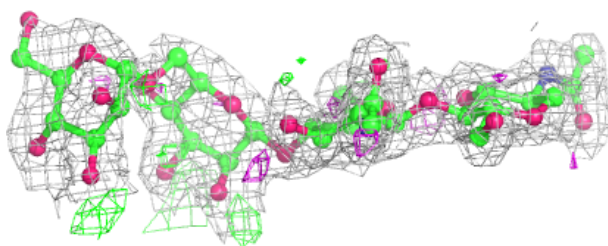
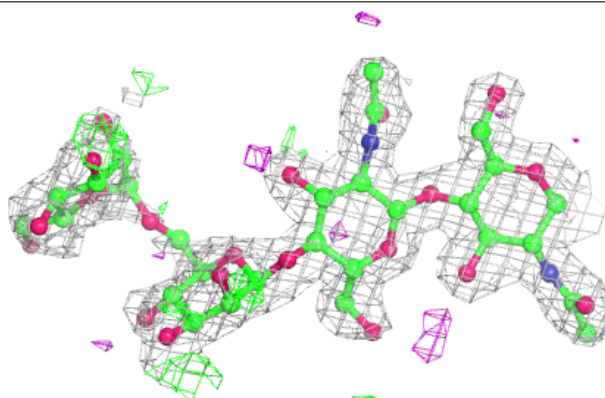
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Chain E:

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

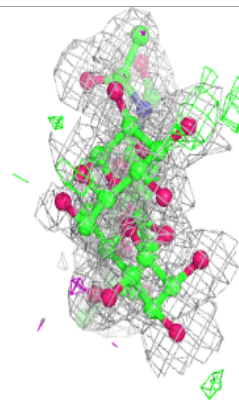
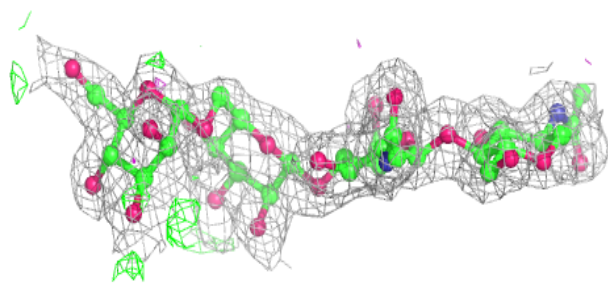
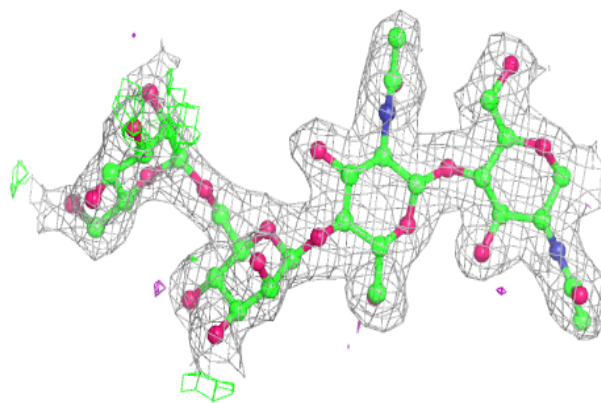
**Electron density around Chain F:**

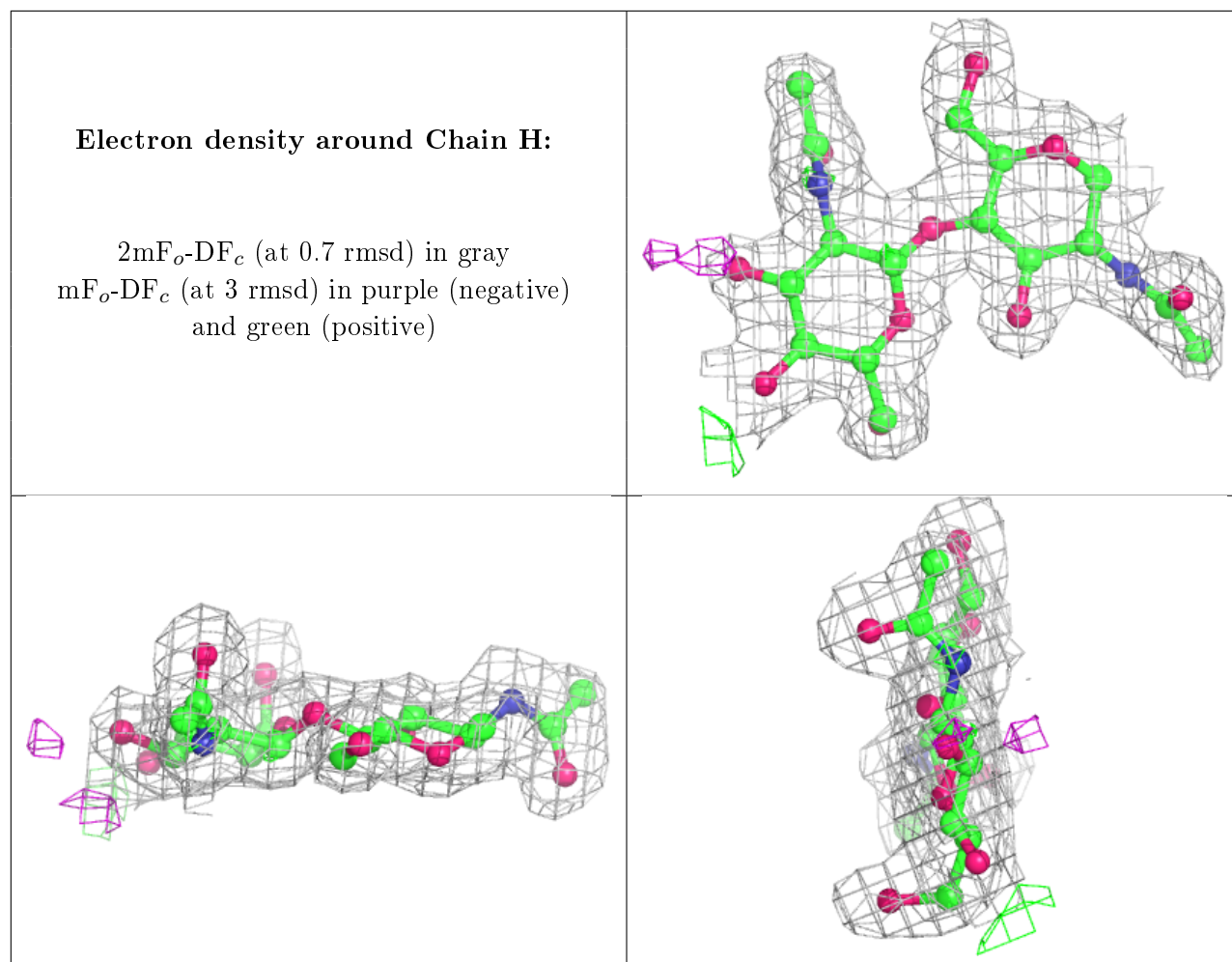
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



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)





6.4 Ligands [i](#)

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

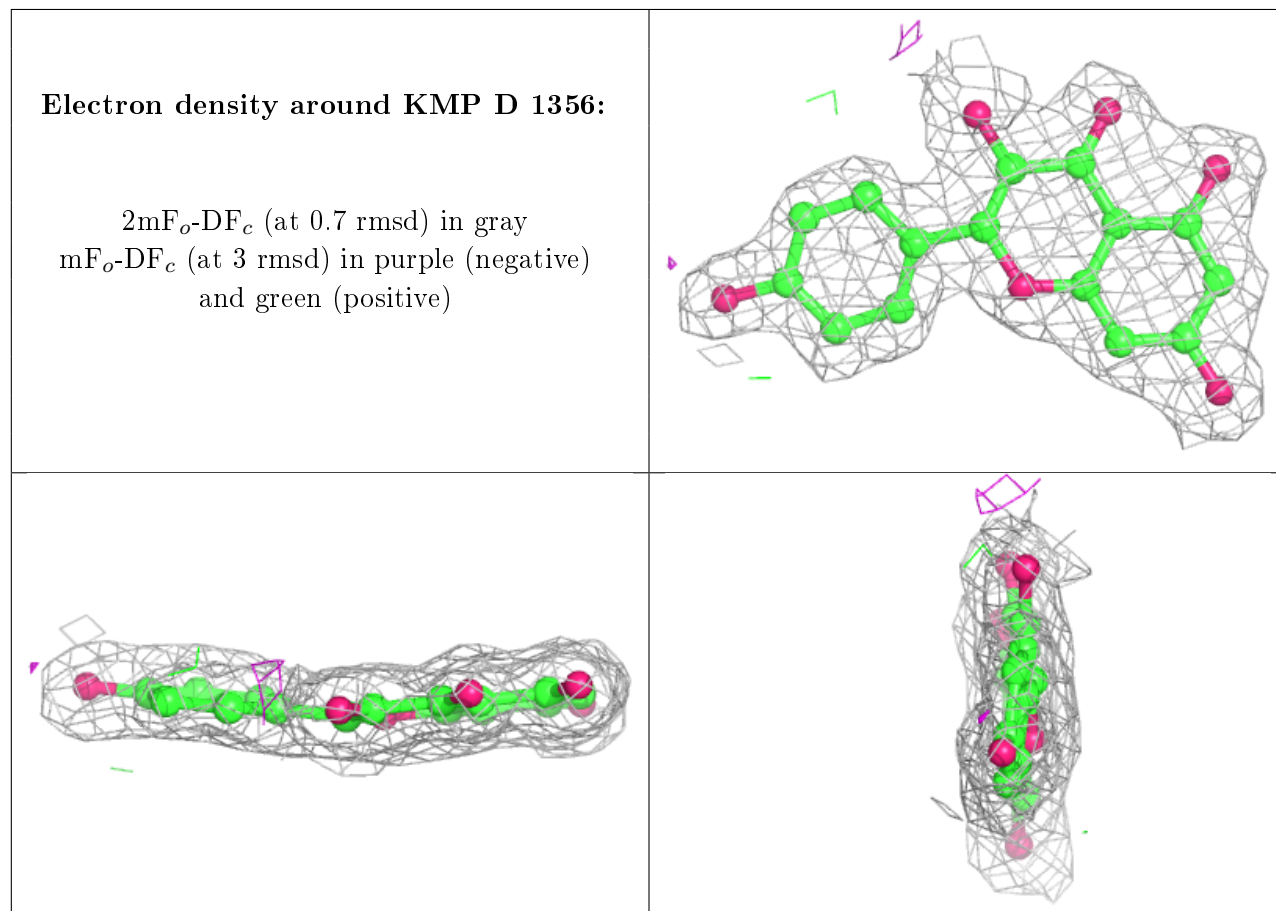
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
8	MPD	B	1361	8/8	0.70	0.33	26,36,39,43	0
8	MPD	B	1360	8/8	0.75	0.31	24,34,37,39	0
5	NAG	A	1356	14/15	0.80	0.15	26,33,36,39	0
8	MPD	C	1360	8/8	0.80	0.23	24,34,38,42	0
5	NAG	C	1355	14/15	0.83	0.15	34,44,49,49	0
8	MPD	A	1362	8/8	0.84	0.23	20,29,31,34	0
5	NAG	D	1351	14/15	0.85	0.17	30,40,42,44	0
5	NAG	B	1355	14/15	0.86	0.16	31,41,44,47	0
5	NAG	A	1351	14/15	0.87	0.14	26,32,39,39	0
5	NAG	C	1354	14/15	0.88	0.14	28,36,46,47	0

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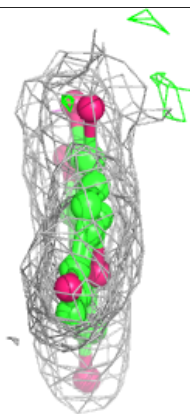
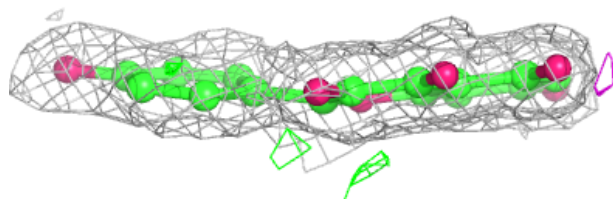
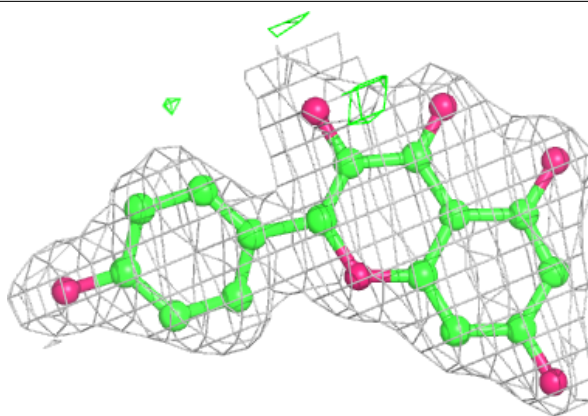
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
5	NAG	D	1355	14/15	0.88	0.13	27,35,44,46	0
5	NAG	B	1354	14/15	0.89	0.18	22,31,42,43	0
5	NAG	D	1354	14/15	0.89	0.13	22,28,32,36	0
5	NAG	A	1354	14/15	0.91	0.10	20,32,41,41	0
5	NAG	A	1355	14/15	0.91	0.24	28,37,44,46	0
5	NAG	B	1351	14/15	0.91	0.15	23,32,34,35	0
5	NAG	C	1351	14/15	0.92	0.13	27,34,40,43	0
6	KMP	D	1356	21/21	0.93	0.09	12,16,18,19	0
6	KMP	C	1358	21/21	0.93	0.10	17,18,20,23	0
6	KMP	B	1358	21/21	0.94	0.10	15,18,21,21	0
6	KMP	A	1360	21/21	0.96	0.08	11,14,19,20	0
7	CU	D	1357	1/1	0.99	0.02	11,11,11,11	0
7	CU	C	1359	1/1	0.99	0.02	11,11,11,11	0
7	CU	A	1361	1/1	1.00	0.06	9,9,9,9	0
7	CU	B	1359	1/1	1.00	0.02	10,10,10,10	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.

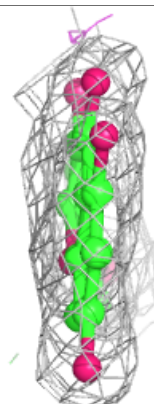
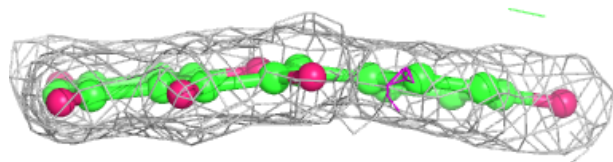
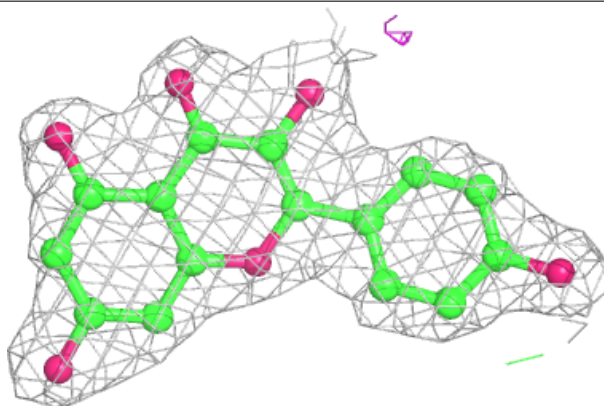


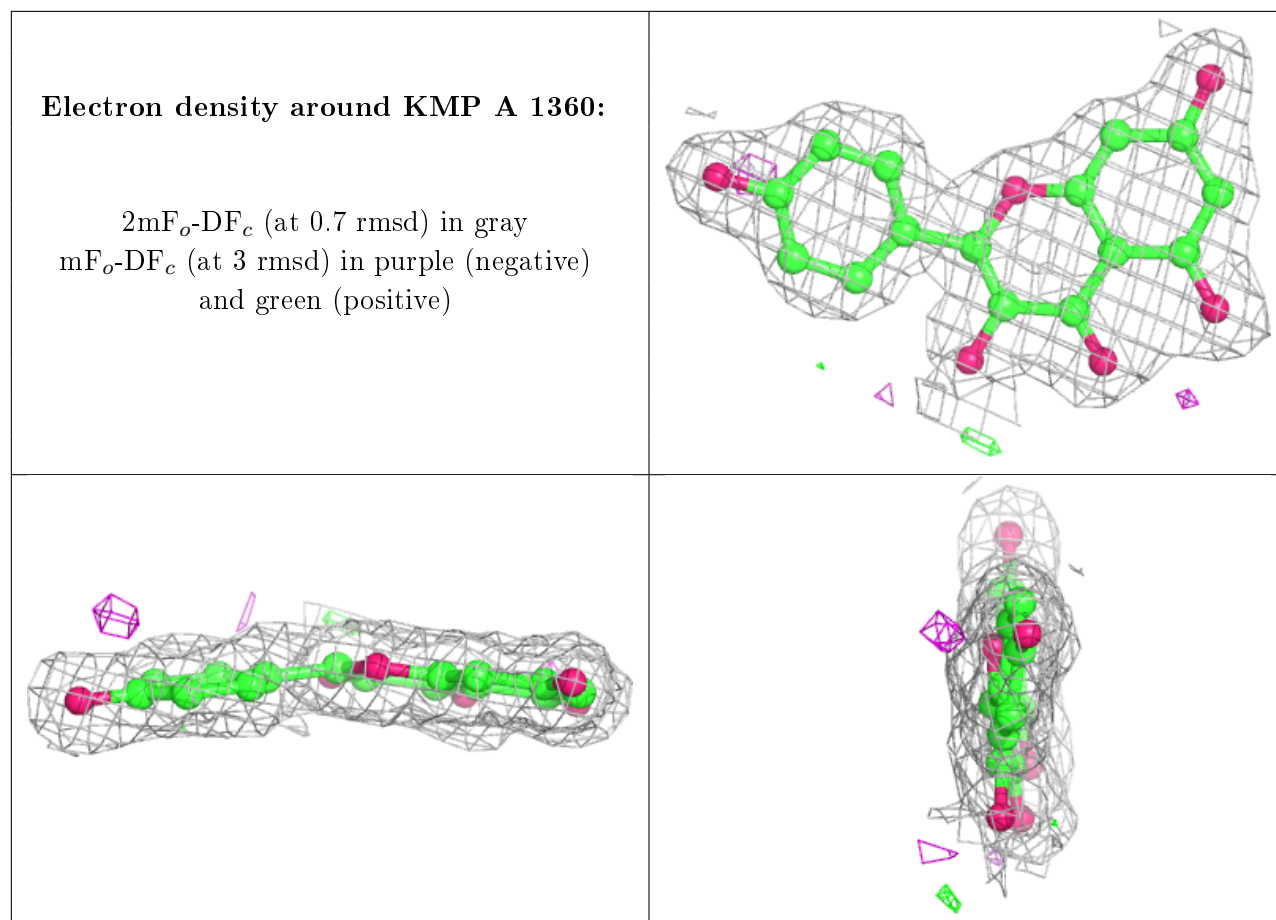
Electron density around KMP C 1358:

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

**Electron density around KMP B 1358:**

$2mF_o-DF_c$ (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.